



Full wwPDB EM Validation Report ⓘ

Aug 26, 2026 – 11:00 am BST

PDB ID : 8C2S / pdb_00008c2s
EMDB ID : EMD-16398
Title : Cryo-EM structure NDUFS4 knockout complex I from Mus musculus heart (Class 1).
Authors : Yin, Z.; Bridges, H.R.; Agip, A.N.A.; Hirst, J.
Deposited on : 2022-12-22
Resolution : 3.90 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

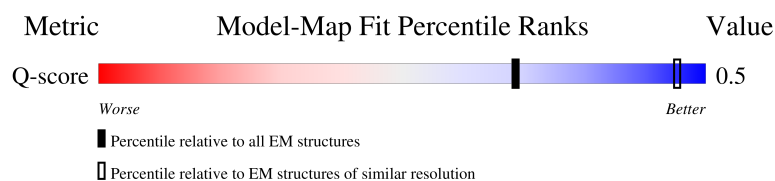
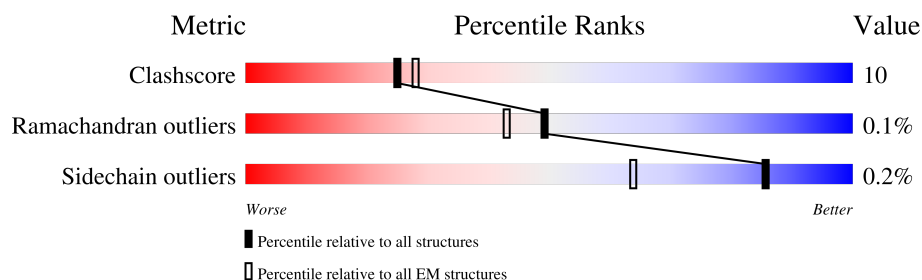
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	8855 (3.40 - 4.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	115	 71% 29%
2	B	224	 52% 18% 30%
3	C	263	 7% 68% 11% 21%
4	D	463	 75% 18% 7%

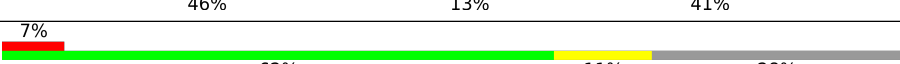
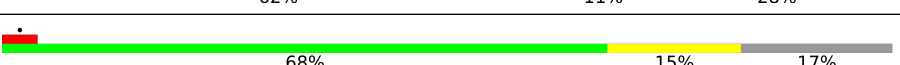
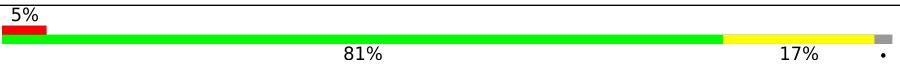
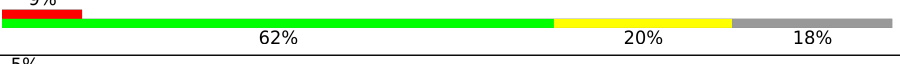
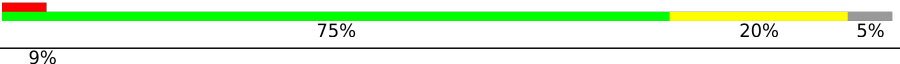
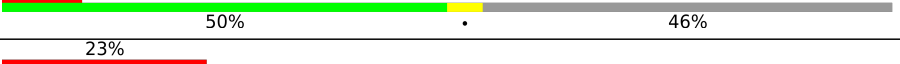
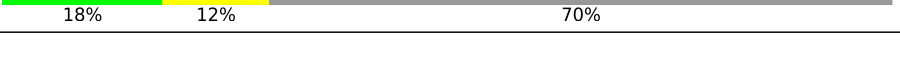
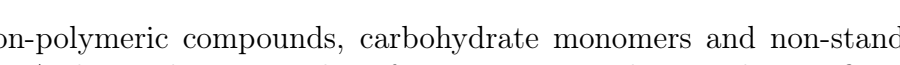
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Mol	Chain	Length	Quality of chain
5	E	245	
6	F	464	
7	G	727	
8	H	318	
9	I	212	
10	J	171	
11	K	98	
12	L	606	
13	M	459	
14	N	345	
15	O	355	
16	P	377	
17	R	116	
18	S	99	
19	T	156	
19	U	156	
20	V	116	
21	W	131	
22	X	172	
23	Y	143	
24	Z	144	
25	a	70	
26	b	84	
27	c	76	
28	d	120	

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Mol	Chain	Length	Quality of chain
29	e	106	
30	f	57	
31	g	151	
32	h	189	
33	i	127	
34	j	105	
35	k	104	
36	l	186	
37	m	129	
38	n	179	
39	o	137	
40	p	176	
41	r	113	
42	s	104	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
43	SF4	G	802	-	-	X	-
43	SF4	I	202	-	-	X	-
46	FMN	F	502	-	X	-	-

2 Entry composition

There are 53 unique types of molecules in this entry. The entry contains 63667 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called NADH-ubiquinone oxidoreductase chain 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	115	Total	C	N	O	S	0	0
			933	633	133	160	7		

- Molecule 2 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	156	Total	C	N	O	S	0	0
			1247	796	223	214	14		

- Molecule 3 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	207	Total	C	N	O	S	0	0
			1721	1111	296	311	3		

- Molecule 4 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	430	Total	C	N	O	S	0	0
			3463	2215	595	629	24		

- Molecule 5 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	210	Total	C	N	O	S	0	0
			1639	1043	275	310	11		

- Molecule 6 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	428	Total	C	N	O	S	0	0
			3300	2080	589	609	22		

- Molecule 7 is a protein called NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	688	Total	C	N	O	S	0	0
			5248	3294	910	1003	41		

- Molecule 8 is a protein called NADH-ubiquinone oxidoreductase chain 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	318	Total	C	N	O	S	0	0
			2540	1706	384	428	22		

- Molecule 9 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	178	Total	C	N	O	S	0	0
			1431	898	245	276	12		

- Molecule 10 is a protein called NADH-ubiquinone oxidoreductase chain 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	171	Total	C	N	O	S	0	0
			1300	874	185	226	15		

- Molecule 11 is a protein called NADH-ubiquinone oxidoreductase chain 4L.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	98	Total	C	N	O	S	0	0
			737	477	112	137	11		

- Molecule 12 is a protein called NADH-ubiquinone oxidoreductase chain 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	606	Total	C	N	O	S	0	0
			4800	3182	746	827	45		

- Molecule 13 is a protein called NADH-ubiquinone oxidoreductase chain 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	459	Total	C	N	O	S	0	0
			3632	2408	567	617	40		

- Molecule 14 is a protein called NADH-ubiquinone oxidoreductase chain 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	344	Total	C	N	O	S	0	0
			2696	1791	416	452	37		

- Molecule 15 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	320	Total	C	N	O	S	0	0
			2607	1674	431	492	10		

- Molecule 16 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 9, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	325	Total	C	N	O	S	0	0
			2626	1702	456	461	7		

- Molecule 17 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	R	55	Total	C	N	O	S	0	0
			413	258	74	78	3		

- Molecule 18 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	S	82	Total	C	N	O	S	0	0
			594	374	108	110	2		

- Molecule 19 is a protein called Acyl carrier protein, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	T	76	Total	C	N	O	S	0	0
			611	392	90	124	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
19	U	86	Total	C	N	O	S	0	0
			692	446	102	139	5		

- Molecule 20 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	V	112	Total	C	N	O	S	0	0
			915	596	152	164	3		

- Molecule 21 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	W	109	Total	C	N	O	S	0	0
			935	598	175	156	6		

- Molecule 22 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	X	171	Total	C	N	O	S	0	0
			1396	889	250	247	10		

- Molecule 23 is a protein called MCG5603.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	Y	140	Total	C	N	O	S	0	0
			1037	662	175	192	8		

- Molecule 24 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 13.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	Z	141	Total	C	N	O	S	0	0
			1167	750	207	202	8		

- Molecule 25 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	a	68	Total	C	N	O	S	0	0
			556	360	99	93	4		

- Molecule 26 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	b	80	Total	C	N	O	S	0	0
			628	414	99	111	4		

- Molecule 27 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	c	48	Total	C	N	O	S	0	0
			398	261	69	67	1		

- Molecule 28 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C2.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	d	120	Total	C	N	O	S	0	0
			996	651	171	165	9		

- Molecule 29 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	e	105	Total	C	N	O	S	0	0
			877	555	162	152	8		

- Molecule 30 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	f	53	Total	C	N	O	S	0	0
			456	295	82	77	2		

- Molecule 31 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 11, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	g	101	Total	C	N	O	S	0	0
			850	549	136	161	4		

- Molecule 32 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 5, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	h	138	Total	C	N	O	S	0	0
			1162	762	194	203	3		

- Molecule 33 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	i	94	Total	C	N	O	S	0	0
			787	515	134	135	3		

- Molecule 34 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	j	62	Total	C	N	O	S	0	0
			537	355	88	93	1		

- Molecule 35 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	k	75	Total	C	N	O	S	0	0
			609	404	103	100	2		

- Molecule 36 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	l	154	Total	C	N	O	S	0	0
			1294	834	215	234	11		

- Molecule 37 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	m	126	Total	C	N	O		0	0
			1050	676	189	185			

- Molecule 38 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	n	177	Total	C	N	O	S	0	0
			1534	981	275	267	11		

- Molecule 39 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	o	113	Total	C	N	O	S	0	0
			979	617	184	170	8		

- Molecule 40 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	p	168	Total	C	N	O	S	0	0
			1424	896	256	264	8		

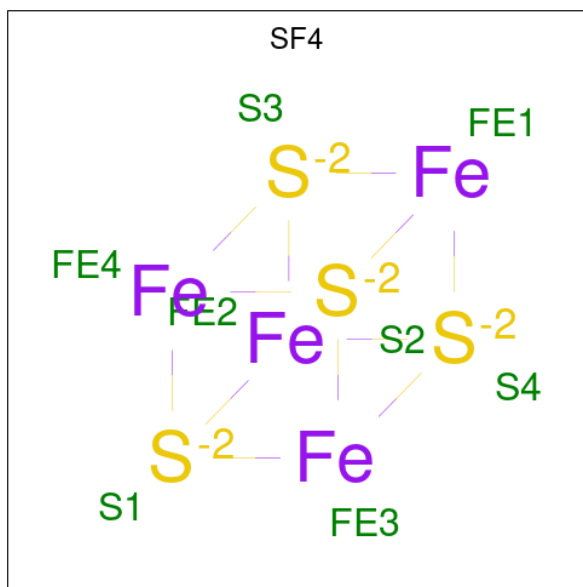
- Molecule 41 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	r	61	Total	C	N	O	S	0	0
			487	311	87	86	3		

- Molecule 42 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 3, mitochondrial.

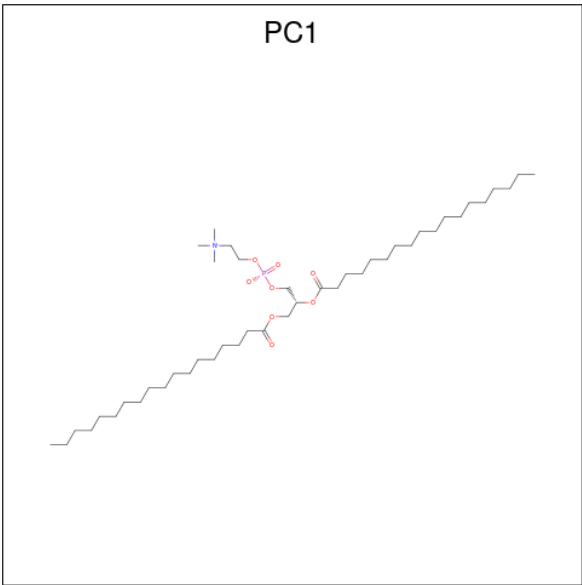
Mol	Chain	Residues	Atoms				AltConf	Trace
42	s	31	Total	C	N	O	0	0
			269	174	45	50		

- Molecule 43 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



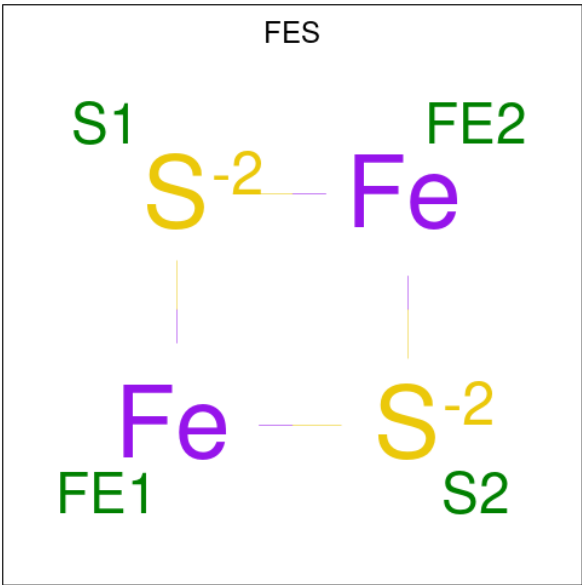
Mol	Chain	Residues	Atoms			AltConf
43	B	1	Total	Fe	S	0
			8	4	4	
43	F	1	Total	Fe	S	0
			8	4	4	
43	G	1	Total	Fe	S	0
			8	4	4	
43	G	1	Total	Fe	S	0
			8	4	4	
43	I	1	Total	Fe	S	0
			8	4	4	
43	I	1	Total	Fe	S	0
			8	4	4	

- Molecule 44 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PC1) (formula: C₄₄H₈₈NO₈P).



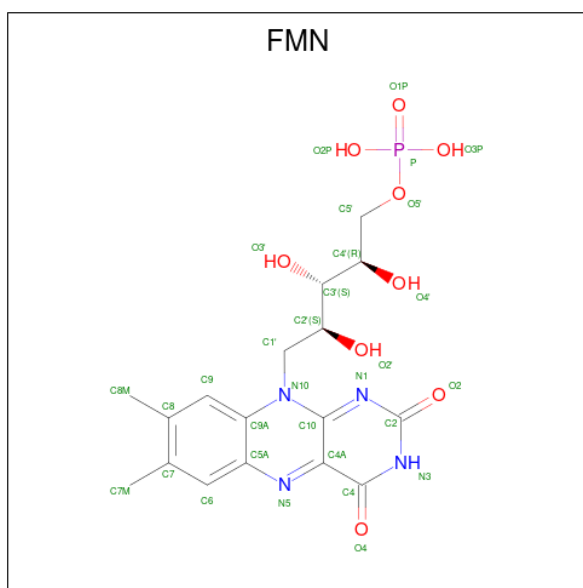
Mol	Chain	Residues	Atoms					AltConf
44	B	1	Total	C	N	O	P	0
			41	31	1	8	1	
44	H	1	Total	C	N	O	P	0
			42	32	1	8	1	

- Molecule 45 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



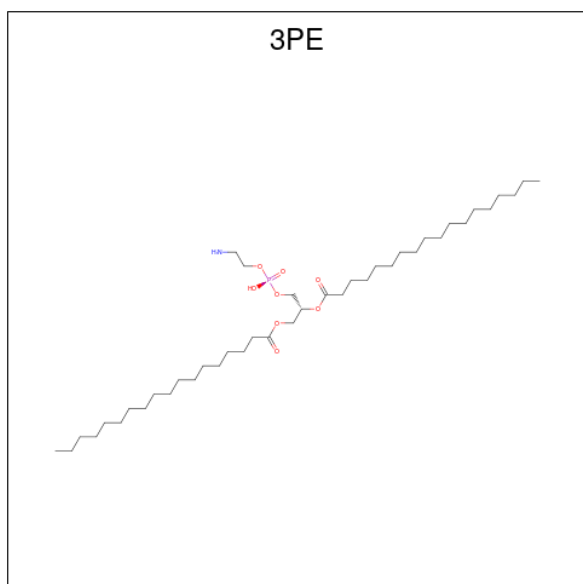
Mol	Chain	Residues	Atoms			AltConf
45	E	1	Total	Fe	S	0
			4	2	2	
45	G	1	Total	Fe	S	0
			4	2	2	

- Molecule 46 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					AltConf
46	F	1	Total	C	N	O	P	0
			31	17	4	9	1	

- Molecule 47 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (CCD ID: 3PE) (formula: $C_{41}H_{82}NO_8P$).



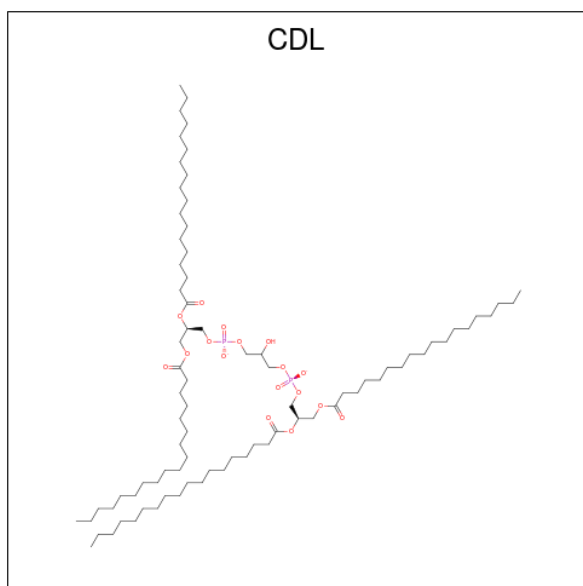
Mol	Chain	Residues	Atoms					AltConf
47	H	1	Total	C	N	O	P	0
			44	34	1	8	1	

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Mol	Chain	Residues	Atoms					AltConf
47	I	1	Total	C	N	O	P	0
			51	41	1	8	1	
47	J	1	Total	C	N	O	P	0
			33	23	1	8	1	
47	L	1	Total	C	N	O	P	0
			49	39	1	8	1	
47	M	1	Total	C	N	O	P	0
			42	32	1	8	1	
47	Y	1	Total	C	N	O	P	0
			41	31	1	8	1	
47	i	1	Total	C	N	O	P	0
			42	32	1	8	1	

- Molecule 48 is CARDIOLIPIN (CCD ID: CDL) (formula: $C_{81}H_{156}O_{17}P_2$).



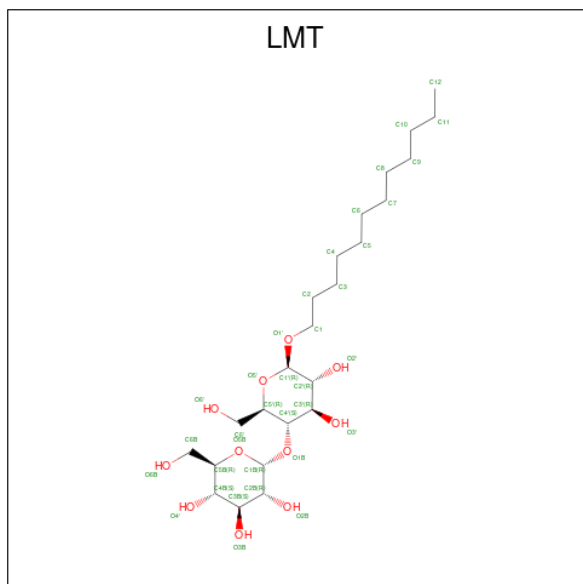
Mol	Chain	Residues	Atoms				AltConf
48	L	1	Total	C	O	P	0
			74	55	17	2	
48	M	1	Total	C	O	P	0
			59	41	16	2	
48	N	1	Total	C	O	P	0
			65	46	17	2	
48	d	1	Total	C	O	P	0
			67	48	17	2	
48	d	1	Total	C	O	P	0
			63	44	17	2	

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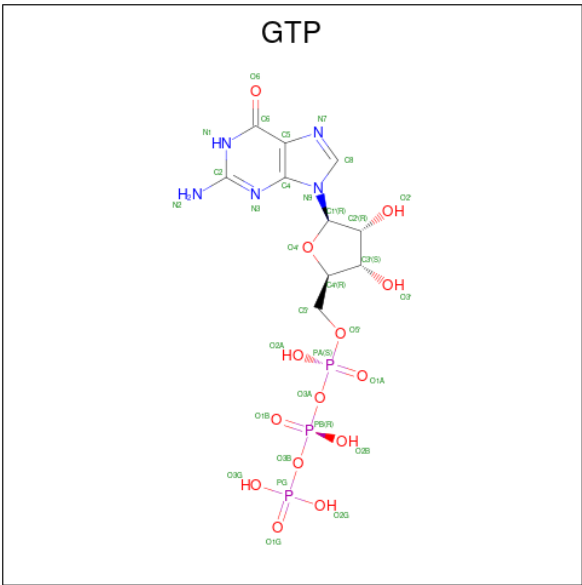
Mol	Chain	Residues	Atoms				AltConf
48	g	1	Total	C	O	P	0
			70	51	17	2	

- Molecule 49 is DODECYL-BETA-D-MALTOSE (CCD ID: LMT) (formula: $C_{24}H_{46}O_{11}$).



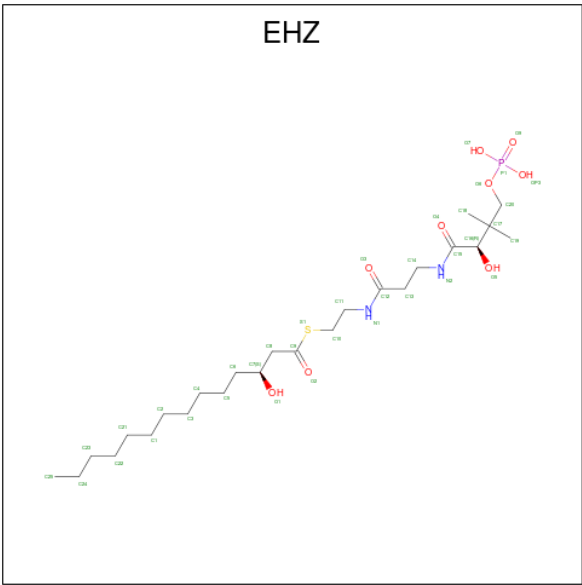
Mol	Chain	Residues	Atoms				AltConf
49	L	1	Total	C	O		0
			35	24	11		
49	L	1	Total	C	O		0
			35	24	11		

- Molecule 50 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



Mol	Chain	Residues	Atoms		AltConf
52	R	1	Total	Zn	0
			1	1	

- Molecule 53 is {S}-[2-[3-[[2 {R}]-3,3-dimethyl-2-oxidanyl-4-phosphonooxy-butanoyl]amino]propanoylamino]ethyl] (3 {S})-3-oxidanyltetradecanethioate (CCD ID: EHZ) (formula: C₂₅H₄₉N₂O₉PS).



Mol	Chain	Residues	Atoms						AltConf
53	T	1	Total	C	N	O	P	S	0
			37	25	2	8	1	1	
53	U	1	Total	C	N	O	P	S	0
			37	25	2	8	1	1	

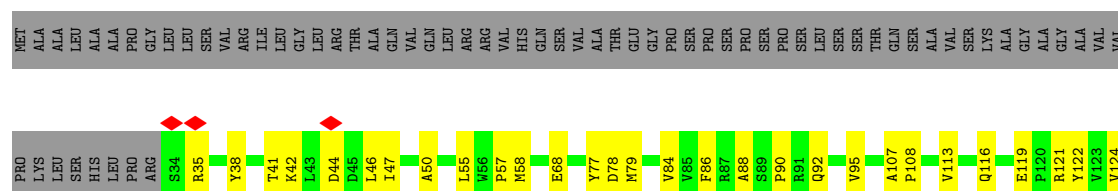
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

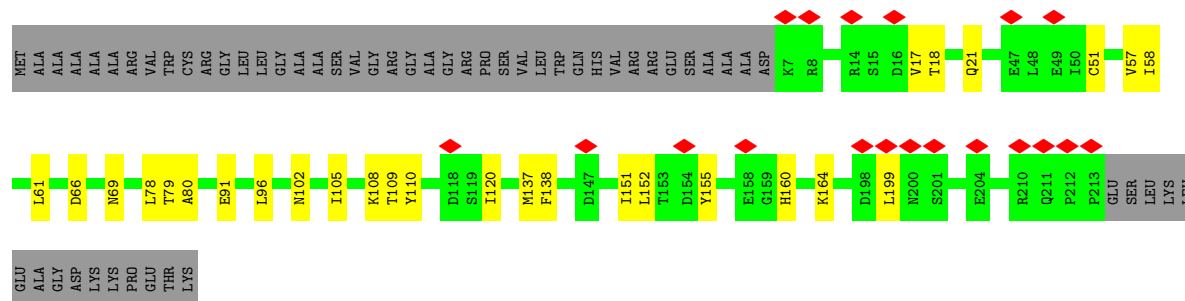
- Molecule 1: NADH-ubiquinone oxidoreductase chain 3



- Molecule 2: NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial

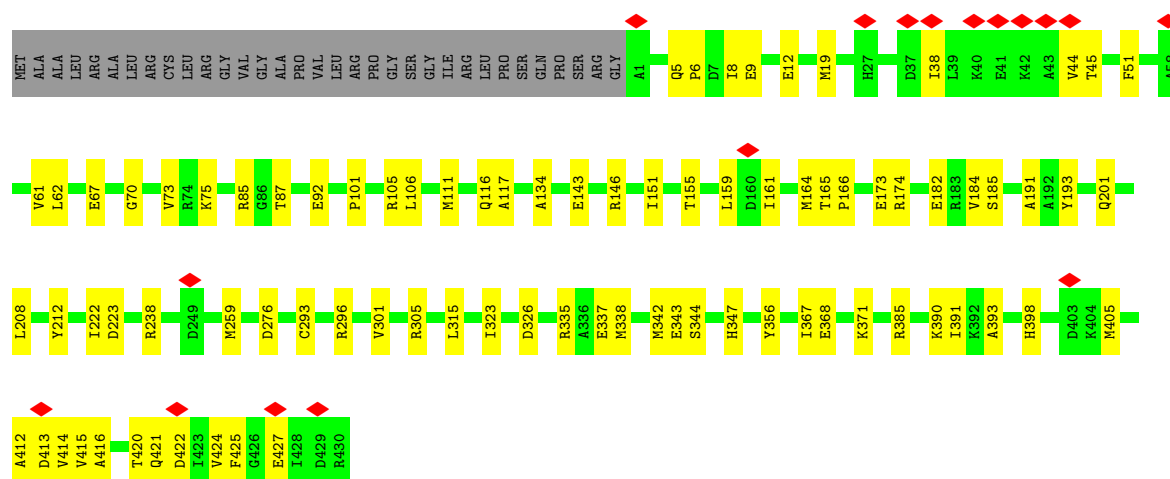


- Molecule 3: NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial



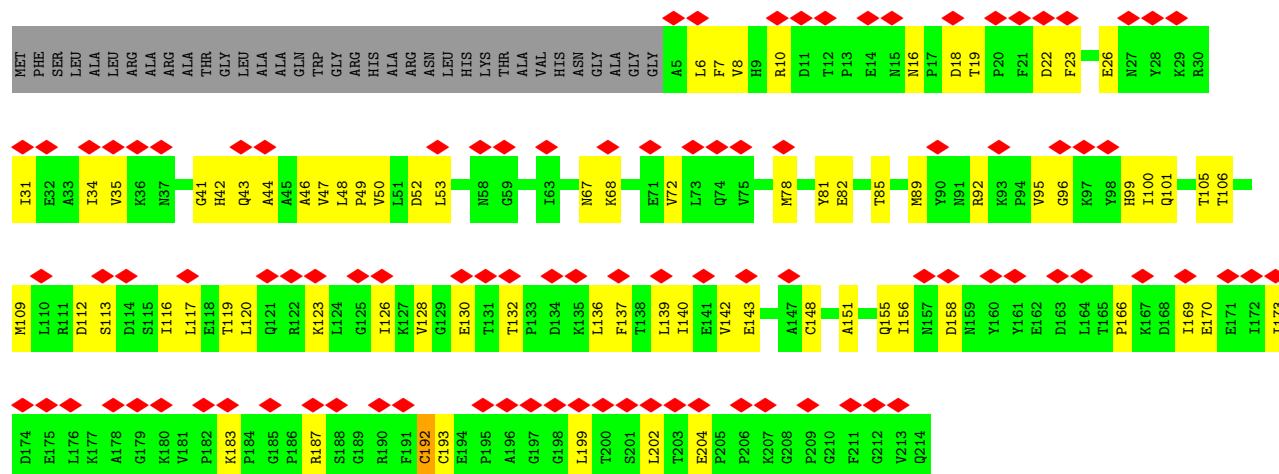
- Molecule 4: NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial

Chain D: 



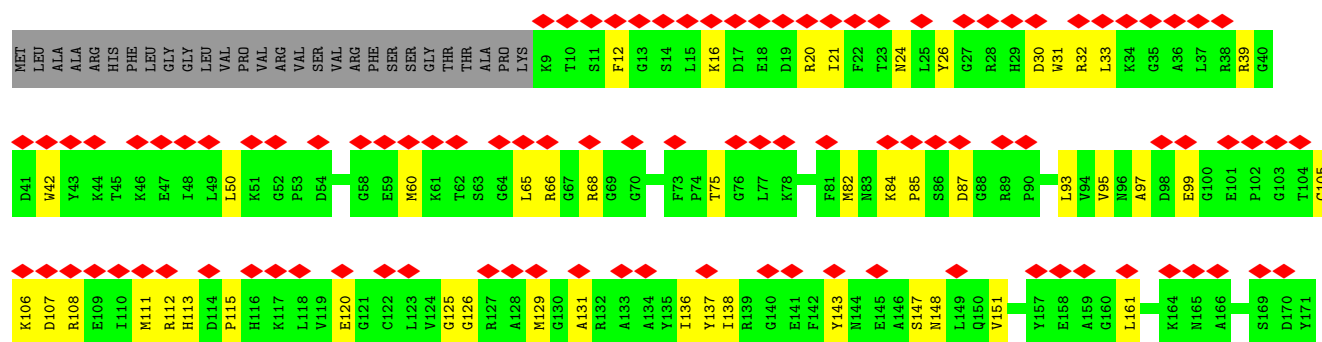
• Molecule 5: NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial

Chain E: 

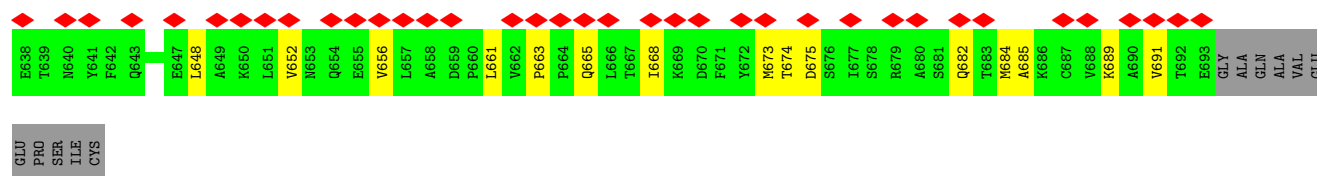


• Molecule 6: NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial

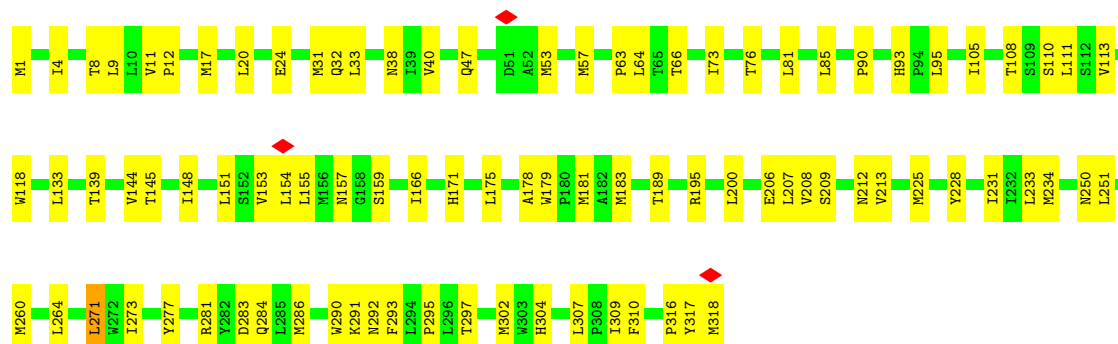
Chain F: 



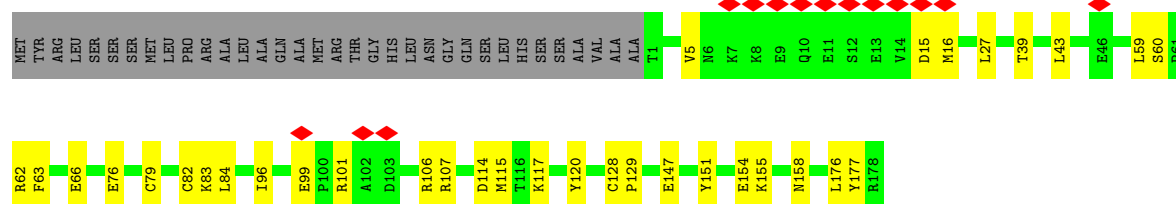




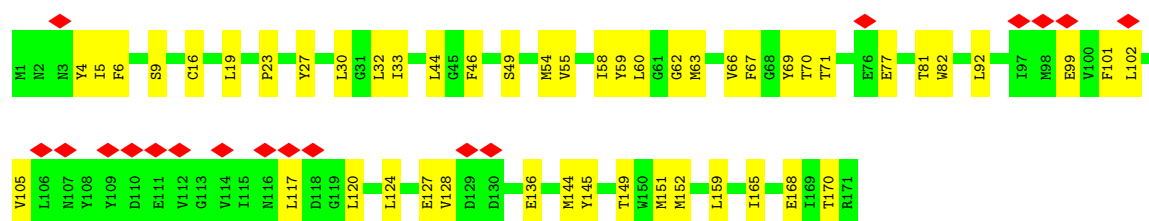
• Molecule 8: NADH-ubiquinone oxidoreductase chain 1



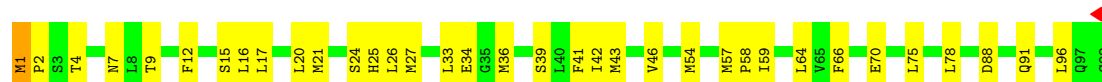
• Molecule 9: NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial



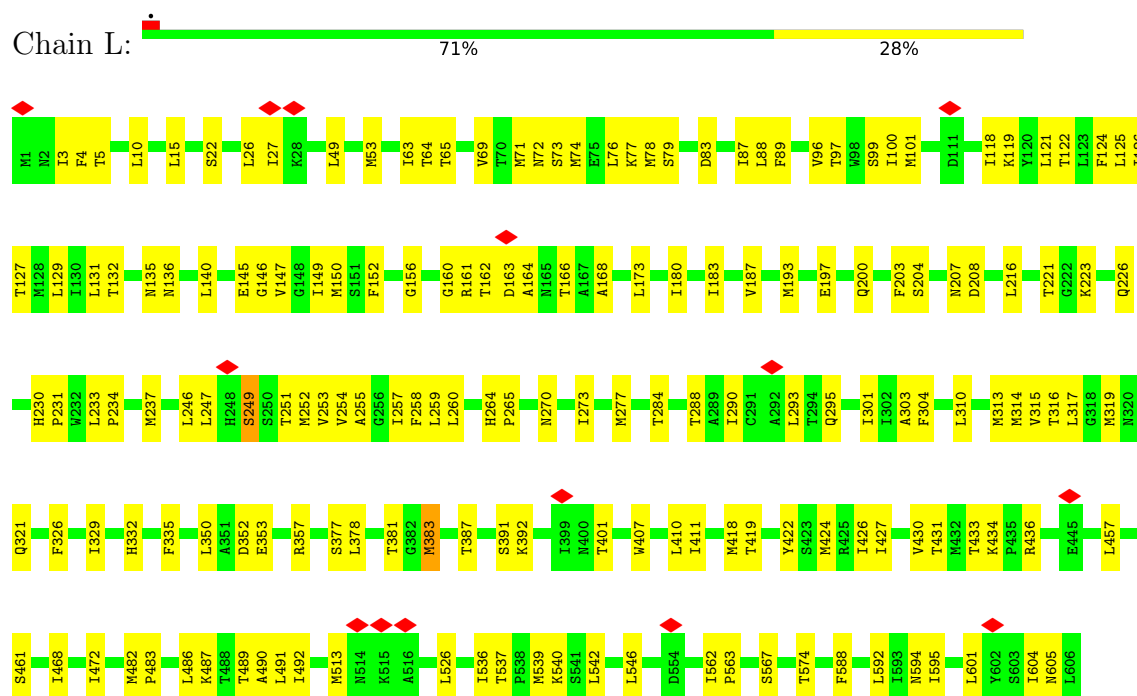
• Molecule 10: NADH-ubiquinone oxidoreductase chain 6



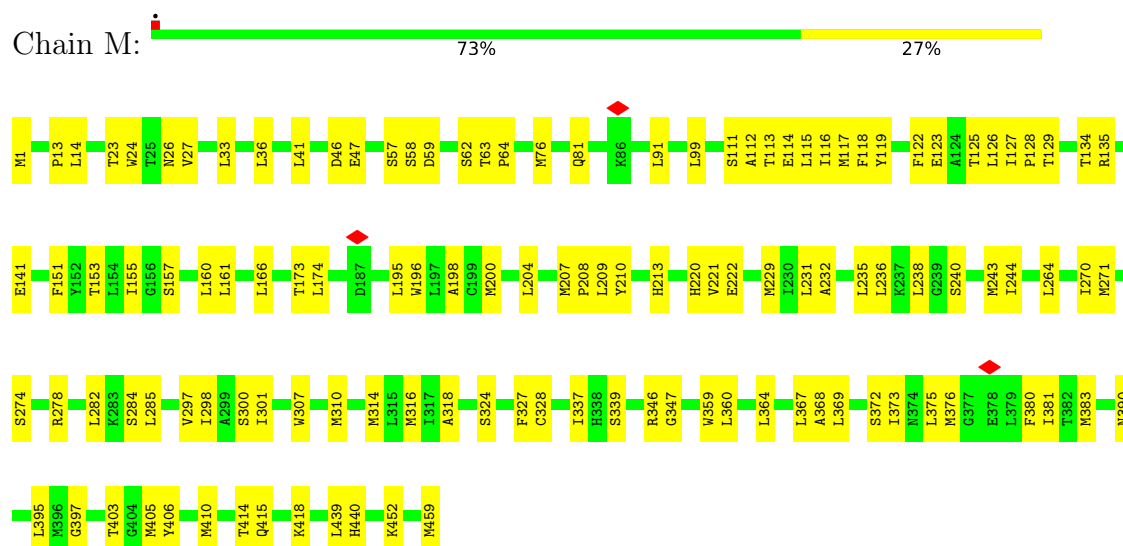
• Molecule 11: NADH-ubiquinone oxidoreductase chain 4L



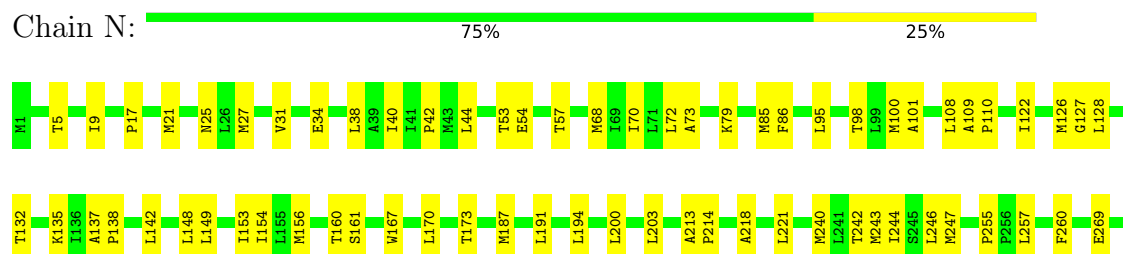
- Molecule 12: NADH-ubiquinone oxidoreductase chain 5



- Molecule 13: NADH-ubiquinone oxidoreductase chain 4

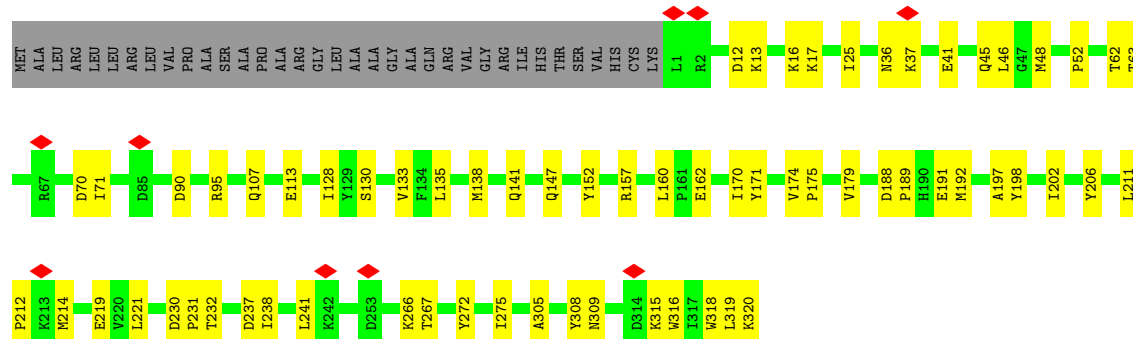


- Molecule 14: NADH-ubiquinone oxidoreductase chain 2

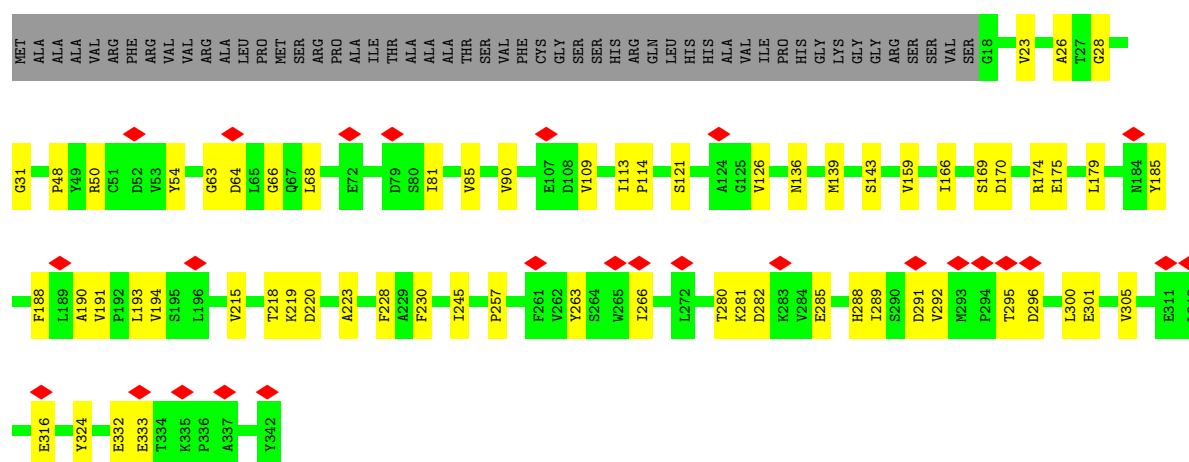




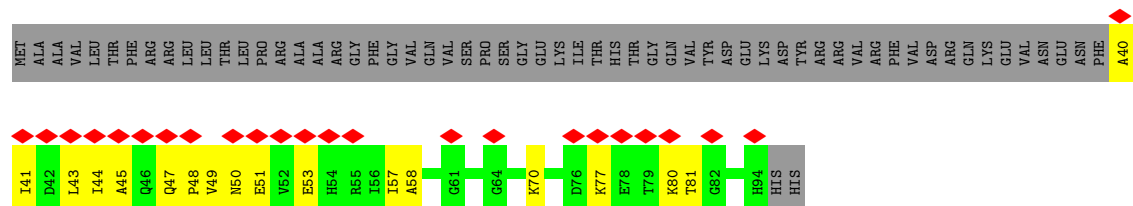
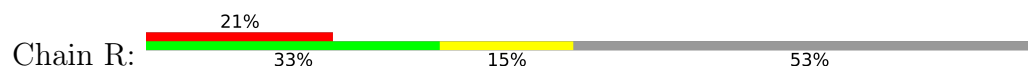
- Molecule 15: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial



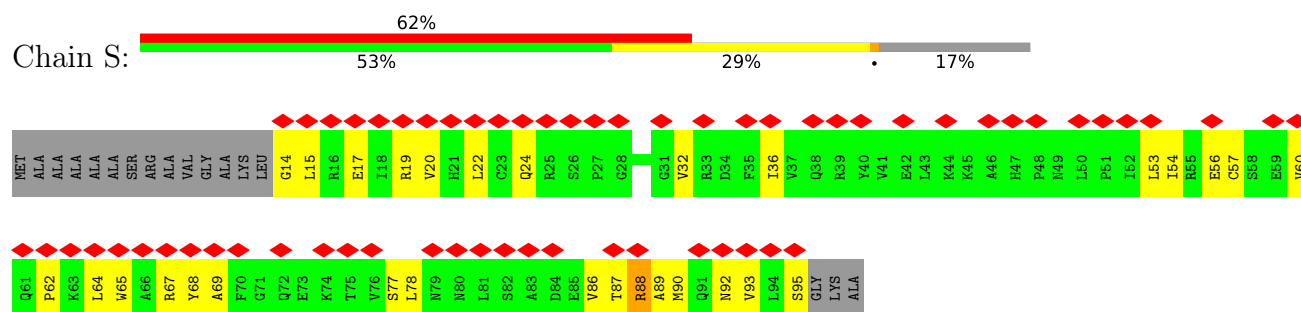
- Molecule 16: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 9, mitochondrial



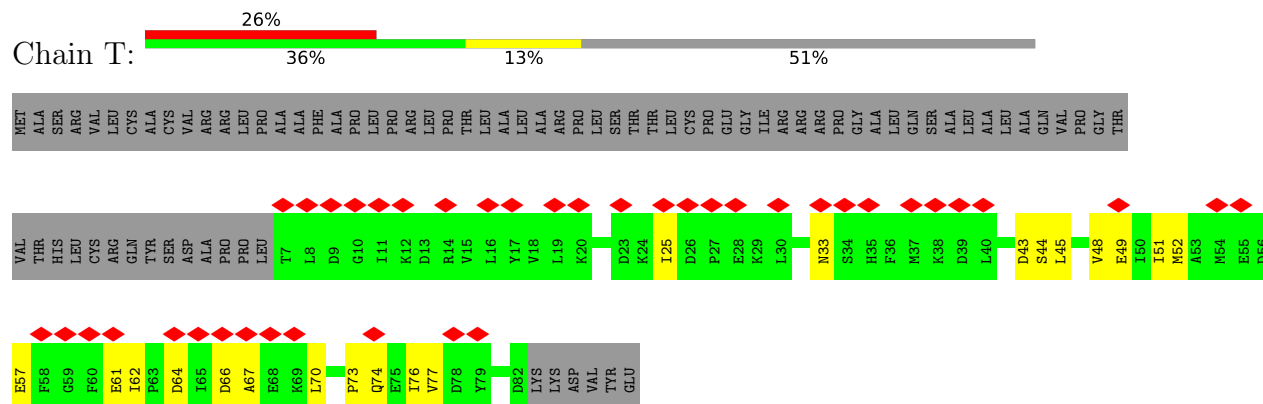
- Molecule 17: NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial



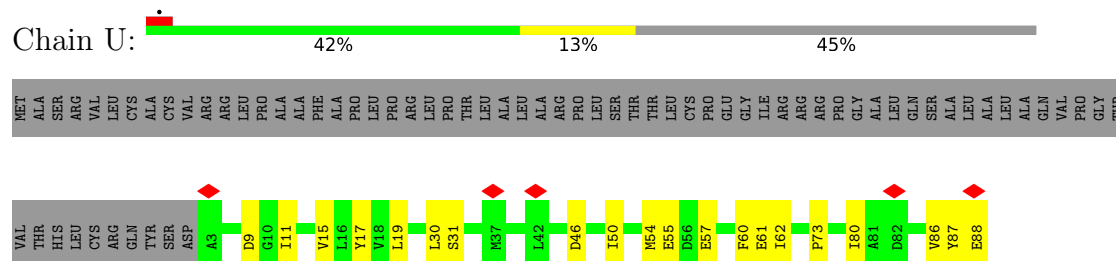
- Molecule 18: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2



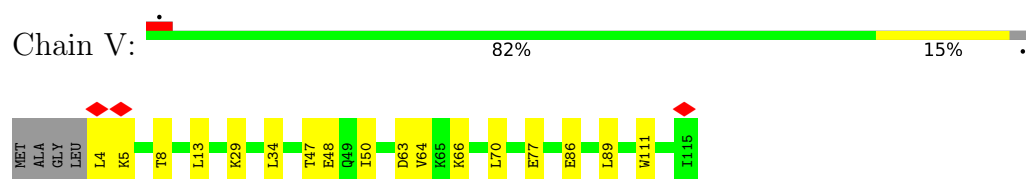
- Molecule 19: Acyl carrier protein, mitochondrial



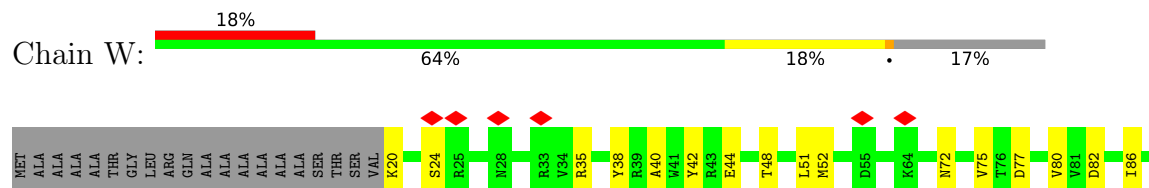
- Molecule 19: Acyl carrier protein, mitochondrial

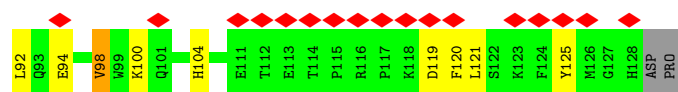


- Molecule 20: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5

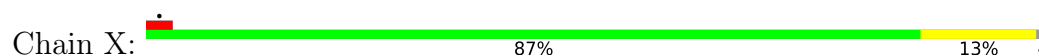


- Molecule 21: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 6

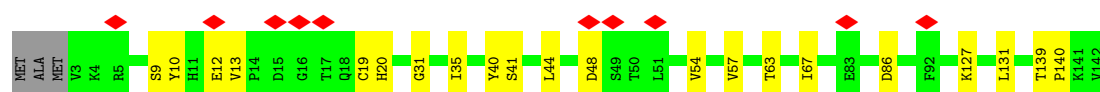
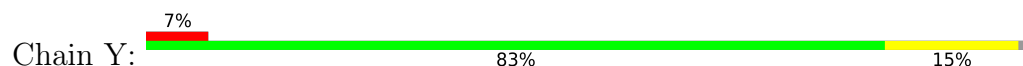




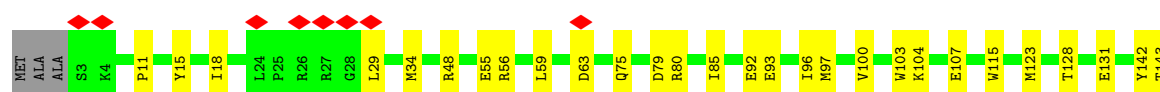
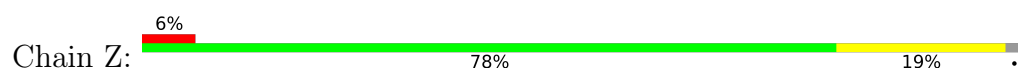
- Molecule 22: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8



- Molecule 23: MCG5603



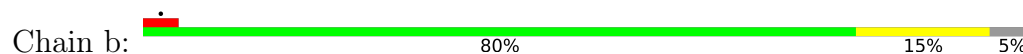
- Molecule 24: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 13



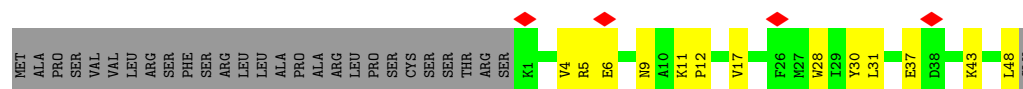
- Molecule 25: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 1



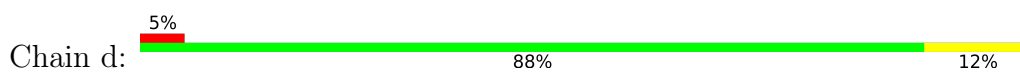
- Molecule 26: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 3



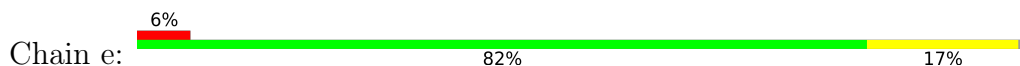
- Molecule 27: NADH dehydrogenase [ubiquinone] 1 subunit C1, mitochondrial



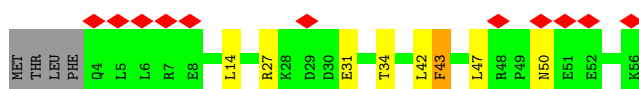
- Molecule 28: NADH dehydrogenase [ubiquinone] 1 subunit C2



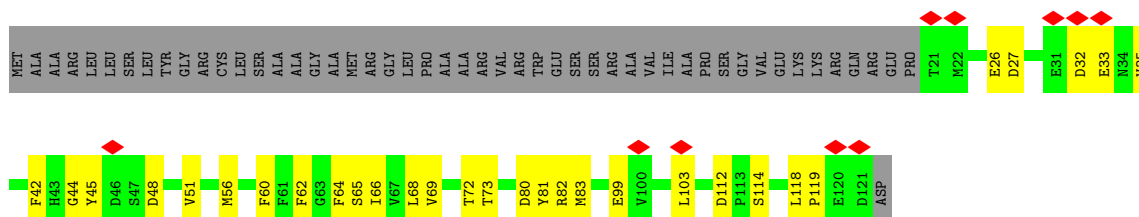
- Molecule 29: NADH dehydrogenase [ubiquinone] iron-sulfur protein 5



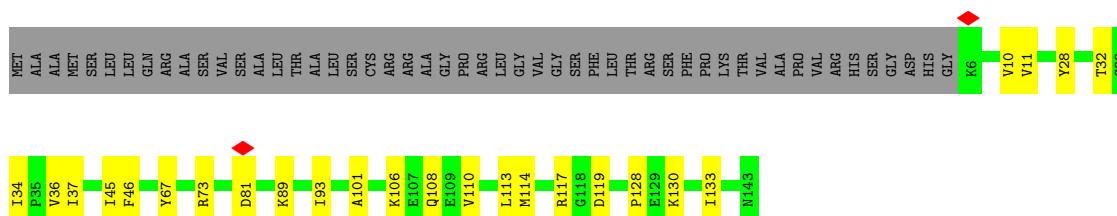
- Molecule 30: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 1



- Molecule 31: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 11, mitochondrial

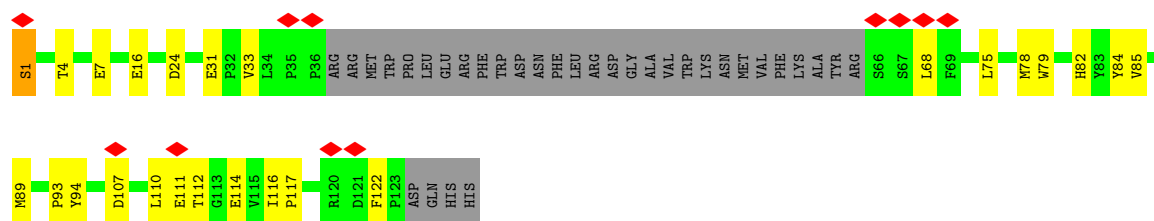


- Molecule 32: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 5, mitochondrial

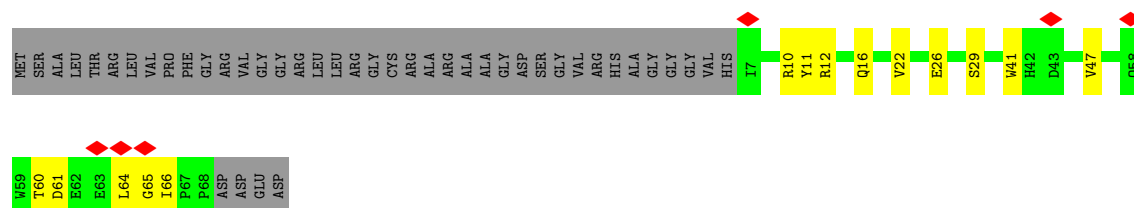


- Molecule 33: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 6





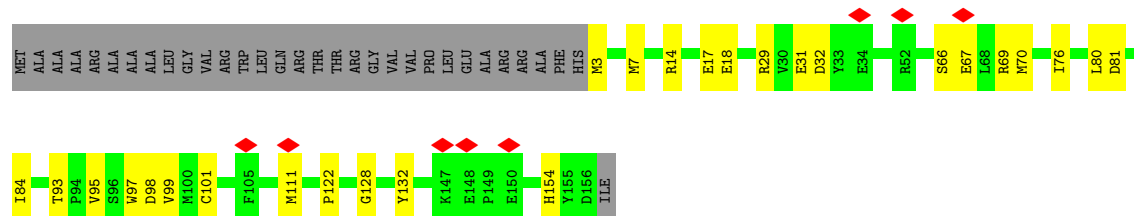
- Molecule 34: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 2, mitochondrial



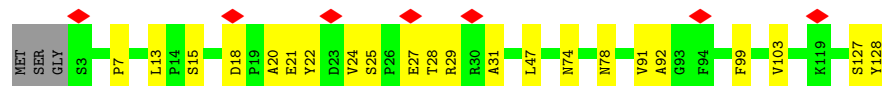
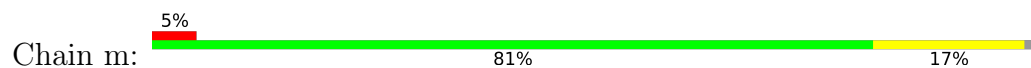
- Molecule 35: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 3



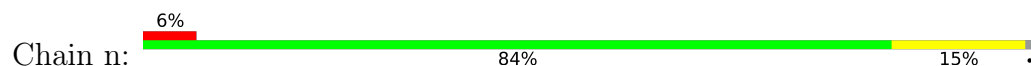
- Molecule 36: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial



- Molecule 37: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 4



- Molecule 38: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 9



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	8793	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	45	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2900	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	12.463	Depositor
Minimum map value	-3.692	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.423	Depositor
Recommended contour level	1.9	Depositor
Map size (Å)	608.4, 608.4, 608.4	wwPDB
Map dimensions	450, 450, 450	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.352, 1.352, 1.352	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3PE, EH2, PC1, SF4, FME, FMN, FES, SAC, CDL, 2MR, GTP, NAP, LMT, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.29	0/949	0.37	0/1297
2	B	0.34	0/1278	0.45	0/1730
3	C	0.29	0/1771	0.36	0/2412
4	D	0.31	0/3539	0.39	0/4793
5	E	0.19	0/1679	0.36	0/2288
6	F	0.21	0/3374	0.40	0/4557
7	G	0.21	0/5335	0.37	0/7236
8	H	0.32	0/2607	0.41	0/3564
9	I	0.30	0/1461	0.36	0/1974
10	J	0.31	0/1322	0.38	0/1799
11	K	0.32	0/738	0.42	0/1002
12	L	0.30	0/4913	0.41	0/6686
13	M	0.33	0/3709	0.42	0/5052
14	N	0.32	0/2748	0.42	0/3741
15	O	0.28	0/2674	0.34	0/3626
16	P	0.26	0/2697	0.35	0/3658
17	R	0.24	0/420	0.41	0/566
18	S	0.19	0/604	0.36	0/827
19	T	0.22	0/620	0.45	0/836
19	U	0.25	0/704	0.36	0/951
20	V	0.26	0/937	0.33	0/1270
21	W	0.26	0/957	0.38	0/1284
22	X	0.26	0/1434	0.36	0/1937
23	Y	0.25	0/1061	0.36	0/1439
24	Z	0.28	0/1198	0.39	0/1616
25	a	0.29	0/569	0.40	0/766
26	b	0.26	0/651	0.36	0/895
27	c	0.28	0/409	0.37	0/555
28	d	0.28	0/1028	0.34	0/1387
29	e	0.28	0/900	0.36	0/1199
30	f	0.26	0/468	0.41	0/630
31	g	0.27	0/878	0.36	0/1196

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	h	0.28	0/1197	0.38	0/1621
33	i	0.25	0/804	0.36	0/1094
34	j	0.24	0/561	0.31	0/768
35	k	0.24	0/629	0.35	0/851
36	l	0.25	0/1348	0.32	0/1840
37	m	0.27	0/1079	0.35	0/1463
38	n	0.26	0/1589	0.34	0/2152
39	o	0.22	0/1004	0.34	0/1348
40	p	0.27	0/1457	0.35	0/1969
41	r	0.21	0/502	0.31	0/680
42	s	0.19	0/277	0.40	0/377
All	All	0.27	0/64079	0.38	0/86932

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
13	M	0	1
19	T	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
13	M	116	ILE	Peptide
19	T	33	ASN	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	933	0	969	34	0
2	B	1247	0	1255	33	0
3	C	1721	0	1680	22	0
4	D	3463	0	3416	69	0
5	E	1639	0	1631	56	0
6	F	3300	0	3258	105	0
7	G	5248	0	5236	144	0
8	H	2540	0	2626	80	0
9	I	1431	0	1383	32	0
10	J	1300	0	1315	54	0
11	K	737	0	768	35	0
12	L	4800	0	4985	148	0
13	M	3632	0	3853	100	0
14	N	2696	0	2895	82	0
15	O	2607	0	2566	52	0
16	P	2626	0	2642	38	0
17	R	413	0	403	17	0
18	S	594	0	543	24	0
19	T	611	0	602	15	0
19	U	692	0	686	17	0
20	V	915	0	954	12	0
21	W	935	0	959	25	0
22	X	1396	0	1383	23	0
23	Y	1037	0	1024	16	0
24	Z	1167	0	1166	27	0
25	a	556	0	568	16	0
26	b	628	0	628	8	0
27	c	398	0	401	15	0
28	d	996	0	1001	18	0
29	e	877	0	871	19	0
30	f	456	0	452	11	0
31	g	850	0	783	30	0
32	h	1162	0	1163	26	0
33	i	787	0	797	29	0
34	j	537	0	495	13	0
35	k	609	0	603	13	0
36	l	1294	0	1186	23	0
37	m	1050	0	1061	15	0
38	n	1534	0	1466	19	0
39	o	979	0	963	20	0
40	p	1424	0	1393	40	0
41	r	487	0	502	4	0
42	s	269	0	256	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
43	B	8	0	0	0	0
43	F	8	0	0	1	0
43	G	16	0	0	2	0
43	I	16	0	0	2	0
44	B	41	0	59	0	0
44	H	42	0	58	3	0
45	E	4	0	0	1	0
45	G	4	0	0	1	0
46	F	31	0	18	2	0
47	H	44	0	65	1	0
47	I	51	0	82	1	0
47	J	33	0	40	0	0
47	L	49	0	75	1	0
47	M	42	0	61	1	0
47	Y	41	0	56	0	0
47	i	42	0	61	3	0
48	L	74	0	92	1	0
48	M	59	0	68	3	0
48	N	65	0	77	0	0
48	d	130	0	151	0	0
48	g	70	0	87	0	0
49	L	70	0	88	1	0
50	O	31	0	10	2	0
51	P	48	0	25	1	0
52	R	1	0	0	0	0
53	T	37	0	0	1	0
53	U	37	0	0	1	0
All	All	63667	0	63960	1335	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (1335) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:418:LYS:NZ	38:n:91:GLU:OE2	1.95	0.99
12:L:567:SER:OG	49:L:704:LMT:O2'	1.84	0.94
31:g:68:LEU:O	31:g:72:THR:OG1	1.84	0.93
6:F:303:SER:HG	6:F:354:TYR:HH	1.16	0.89
31:g:83:MET:SD	40:p:96:LYS:NZ	2.47	0.87
16:P:50:ARG:NH1	51:P:501:NAP:O3X	2.08	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:259:MET:SD	4:D:421:GLN:NE2	2.47	0.87
15:O:113:GLU:OE1	15:O:266:LYS:NZ	2.07	0.87
10:J:16:CYS:SG	11:K:15:SER:OG	2.33	0.86
15:O:25:ILE:HD12	15:O:48:MET:HE1	1.55	0.86
6:F:401:GLY:O	7:G:64:LYS:NZ	2.10	0.85
15:O:162:GLU:N	15:O:162:GLU:OE1	2.09	0.85
12:L:161:ARG:NH2	12:L:163:ASP:OD2	2.10	0.84
10:J:66:VAL:O	10:J:70:THR:HG23	1.77	0.84
13:M:346:ARG:NE	36:l:67:GLU:OE2	2.10	0.84
8:H:250:ASN:OD1	8:H:251:LEU:N	2.11	0.83
37:m:15:SER:OG	37:m:21:GLU:OE1	1.96	0.83
3:C:164:LYS:NZ	4:D:92:GLU:OE1	2.11	0.82
4:D:173:GLU:N	4:D:173:GLU:OE1	2.12	0.82
6:F:263:ASN:ND2	6:F:287:VAL:O	2.13	0.81
10:J:168:GLU:CD	14:N:5:THR:HG21	2.05	0.81
2:B:58:MET:HE1	2:B:113:VAL:HG11	1.61	0.81
6:F:224:ASN:OD1	6:F:225:VAL:N	2.13	0.80
11:K:1:FME:SD	14:N:79:LYS:NZ	2.55	0.80
13:M:244:ILE:O	40:p:148:TYR:OH	1.98	0.80
37:m:128:TYR:OH	40:p:145:LEU:O	1.99	0.80
22:X:16:GLU:N	22:X:16:GLU:OE1	2.14	0.80
2:B:79:MET:HE1	2:B:86:PHE:CD2	2.18	0.79
40:p:98:ASP:OD2	40:p:142:TYR:OH	2.01	0.79
7:G:260:GLU:OE2	7:G:397:LYS:NZ	2.15	0.79
36:l:128:GLY:O	39:o:97:ARG:NH1	2.15	0.78
14:N:312:LYS:O	14:N:316:HIS:ND1	2.17	0.78
1:A:4:TYR:HH	10:J:4:TYR:HH	1.24	0.77
28:d:119:VAL:HG11	40:p:146:GLY:HA2	1.65	0.77
5:E:183:LYS:O	5:E:187:ARG:NH1	2.17	0.77
5:E:31:ILE:HD11	5:E:53:LEU:HD23	1.67	0.77
13:M:46:ASP:OD1	13:M:47:GLU:N	2.17	0.77
12:L:65:THR:HA	33:i:89:MET:HE2	1.64	0.77
21:W:24:SER:OG	21:W:82:ASP:OD1	2.02	0.77
34:j:29:SER:OG	35:k:74:LYS:O	2.03	0.76
36:l:98:ASP:OD1	36:l:99:VAL:N	2.17	0.76
6:F:24:ASN:ND2	6:F:113:HIS:O	2.19	0.76
7:G:257:ASP:O	7:G:394:ARG:NH2	2.19	0.75
16:P:174:ARG:NH2	16:P:316:GLU:OE1	2.19	0.75
2:B:119:GLU:O	2:B:121:ARG:N	2.19	0.75
12:L:601:LEU:HD22	14:N:153:ILE:HD11	1.68	0.75
12:L:427:ILE:O	12:L:431:THR:OG1	2.05	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:116:ILE:HD11	5:E:169:ILE:HB	1.70	0.74
7:G:469:ALA:O	7:G:472:ASN:ND2	2.21	0.74
16:P:332:GLU:N	16:P:332:GLU:OE1	2.19	0.74
32:h:11:VAL:HG23	38:n:98:VAL:HG13	1.70	0.73
7:G:346:GLU:OE1	7:G:346:GLU:N	2.21	0.73
6:F:137:TYR:CD2	6:F:192:LEU:HD11	2.24	0.72
8:H:195:ARG:HD3	8:H:231:ILE:HD11	1.71	0.72
17:R:47:GLN:N	17:R:47:GLN:OE1	2.22	0.72
31:g:32:ASP:OD1	31:g:33:GLU:N	2.23	0.72
3:C:61:LEU:HD12	3:C:120:ILE:HG21	1.72	0.72
6:F:303:SER:OG	6:F:354:TYR:OH	1.97	0.72
12:L:401:THR:HG21	12:L:482:MET:HE2	1.72	0.72
24:Z:97:MET:HE1	29:e:78:ILE:HA	1.70	0.71
12:L:257:ILE:HG23	12:L:317:LEU:HD11	1.72	0.71
36:l:67:GLU:OE1	36:l:67:GLU:N	2.23	0.71
2:B:46:LEU:HD11	8:H:53:MET:HE1	1.71	0.71
4:D:259:MET:HE2	4:D:259:MET:HA	1.73	0.71
12:L:251:THR:O	12:L:254:VAL:HG22	1.90	0.71
31:g:82:ARG:O	40:p:96:LYS:NZ	2.23	0.71
5:E:16:ASN:ND2	5:E:19:THR:OG1	2.24	0.71
6:F:347:ILE:HD11	6:F:414:PRO:HB3	1.71	0.71
36:l:17:GLU:OE1	36:l:17:GLU:N	2.21	0.71
4:D:193:TYR:OH	4:D:201:GLN:O	2.05	0.70
5:E:31:ILE:CD1	5:E:53:LEU:HD23	2.20	0.70
7:G:324:ASP:HB3	7:G:571:ALA:HB1	1.72	0.70
8:H:264:LEU:HD11	25:a:16:LEU:HD12	1.73	0.70
14:N:156:MET:O	14:N:160:THR:OG1	2.10	0.70
13:M:129:THR:HG21	13:M:235:LEU:HD11	1.73	0.70
8:H:24:GLU:HA	8:H:271:LEU:HD13	1.73	0.70
14:N:27:MET:O	14:N:31:VAL:HG23	1.91	0.69
14:N:319:LYS:NZ	15:O:267:THR:HG21	2.07	0.69
14:N:319:LYS:HZ1	15:O:267:THR:HG21	1.57	0.69
15:O:138:MET:CE	50:O:401:GTP:HN21	2.05	0.69
14:N:108:LEU:HD22	14:N:191:LEU:HD22	1.75	0.69
6:F:253:THR:OG1	6:F:270:GLU:O	2.09	0.69
14:N:314:MET:CE	15:O:267:THR:HG23	2.23	0.69
4:D:371:LYS:NZ	4:D:422:ASP:OD1	2.26	0.68
1:A:33:LYS:NZ	2:B:119:GLU:OE2	2.27	0.68
12:L:89:PHE:CD2	12:L:132:THR:HG21	2.28	0.68
12:L:387:THR:OG1	12:L:461:SER:O	2.09	0.68
42:s:44:THR:O	42:s:48:LEU:HD12	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:648:LEU:O	7:G:652:VAL:HG23	1.92	0.68
6:F:347:ILE:HG21	6:F:418:LEU:HD22	1.74	0.68
29:e:40:ILE:HG21	32:h:133:ILE:HD11	1.76	0.68
6:F:276:LEU:CD1	6:F:317:MET:HE2	2.23	0.68
33:i:82:HIS:ND1	40:p:36:TYR:OH	2.26	0.68
9:I:79:CYS:HA	43:I:202:SF4:S2	2.34	0.67
19:T:43:ASP:OD1	19:T:44:SER:N	2.27	0.67
5:E:8:VAL:HG21	5:E:10:ARG:NH2	2.10	0.66
5:E:148:CYS:N	6:F:105:CYS:SG	2.67	0.66
1:A:68:GLU:HG2	10:J:159:LEU:HD13	1.75	0.66
4:D:111:MET:SD	4:D:111:MET:N	2.69	0.66
10:J:9:SER:OG	11:K:7:ASN:ND2	2.28	0.66
15:O:90:ASP:O	15:O:95:ARG:NH2	2.28	0.66
22:X:154:GLU:N	22:X:154:GLU:OE1	2.28	0.66
13:M:459:MET:O	31:g:82:ARG:NH2	2.28	0.66
7:G:597:TRP:CE3	7:G:619:VAL:HG21	2.30	0.65
14:N:128:LEU:O	14:N:132:THR:OG1	2.12	0.65
17:R:50:ASN:OD1	17:R:51:GLU:N	2.29	0.65
26:b:27:LEU:HA	26:b:30:ILE:HG22	1.77	0.65
16:P:300:LEU:HD22	16:P:305:VAL:HB	1.78	0.65
19:U:86:VAL:HG23	19:U:87:TYR:H	1.62	0.65
13:M:134:THR:OG1	14:N:302:LEU:HD13	1.97	0.65
33:i:82:HIS:NE2	40:p:41:ASP:OD1	2.30	0.65
13:M:452:LYS:NZ	47:M:501:3PE:O14	2.29	0.65
7:G:668:ILE:CG2	7:G:691:VAL:HG21	2.28	0.64
2:B:55:LEU:HD13	2:B:95:VAL:HG11	1.79	0.64
4:D:146:ARG:NH2	4:D:368:GLU:O	2.30	0.64
6:F:276:LEU:HD13	6:F:317:MET:HE2	1.80	0.64
9:I:76:GLU:OE2	17:R:44:ILE:HB	1.97	0.64
28:d:61:GLN:O	28:d:65:VAL:HG23	1.97	0.64
6:F:151:VAL:HG12	42:s:48:LEU:HD21	1.78	0.64
22:X:164:THR:HG23	22:X:164:THR:O	1.96	0.64
13:M:1:FME:HE2	13:M:111:SER:HB2	1.79	0.64
22:X:108:GLN:NE2	22:X:112:ASP:OD2	2.30	0.64
29:e:40:ILE:CG2	32:h:133:ILE:HD11	2.27	0.64
12:L:594:ASN:ND2	23:Y:40:TYR:OH	2.30	0.64
13:M:141:GLU:OE2	13:M:141:GLU:N	2.22	0.64
34:j:16:GLN:N	34:j:16:GLN:OE1	2.31	0.64
6:F:126:GLY:HA2	6:F:131:ALA:HB3	1.80	0.64
12:L:69:VAL:CG1	12:L:76:LEU:HD12	2.27	0.64
11:K:42:ILE:O	11:K:46:VAL:HG23	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:p:105:ILE:HD13	40:p:134:VAL:HG21	1.79	0.64
37:m:24:VAL:HG13	37:m:29:ARG:NH1	2.13	0.64
19:T:51:ILE:HD11	19:T:62:ILE:HG21	1.79	0.63
31:g:73:THR:HG21	32:h:36:VAL:HG11	1.80	0.63
4:D:161:ILE:HD11	4:D:416:ALA:HB2	1.80	0.63
28:d:119:VAL:HG12	40:p:144:ASP:O	1.98	0.63
22:X:18:LYS:HA	25:a:54:ILE:HD12	1.80	0.63
12:L:483:PRO:HD2	12:L:486:LEU:HD12	1.81	0.63
13:M:310:MET:SD	13:M:314:MET:HE2	2.39	0.63
15:O:237:ASP:OD1	20:V:29:LYS:NZ	2.28	0.63
6:F:21:ILE:HG23	6:F:233:THR:HG21	1.80	0.63
10:J:66:VAL:HG13	11:K:27:MET:HE2	1.80	0.63
29:e:58:GLU:N	29:e:58:GLU:OE1	2.32	0.63
13:M:207:MET:O	13:M:298:ILE:HD11	1.99	0.63
15:O:272:TYR:O	15:O:275:ILE:HG23	1.99	0.63
6:F:287:VAL:HG12	6:F:288:THR:H	1.63	0.63
1:A:95:VAL:HG11	8:H:302:MET:HG3	1.80	0.62
14:N:21:MET:HE3	29:e:3:LEU:HD22	1.81	0.62
14:N:218:ALA:HB1	14:N:244:ILE:HD11	1.81	0.62
31:g:112:ASP:OD2	31:g:114:SER:OG	2.16	0.62
2:B:122:TYR:CE2	2:B:177:ILE:HD11	2.33	0.62
10:J:32:LEU:HD21	10:J:59:TYR:CD1	2.34	0.62
12:L:124:PHE:CG	12:L:247:LEU:HD11	2.34	0.62
36:l:132:TYR:O	36:l:154:HIS:NE2	2.32	0.62
40:p:140:ASP:O	40:p:161:ARG:NH2	2.31	0.62
3:C:17:VAL:O	3:C:21:GLN:HG3	1.99	0.62
6:F:120:GLU:OE2	6:F:236:ARG:NH1	2.32	0.62
7:G:624:GLU:HB2	7:G:631:VAL:HG21	1.82	0.62
30:f:31:GLU:O	30:f:34:THR:OG1	2.17	0.62
6:F:99:GLU:OE2	6:F:106:LYS:N	2.33	0.62
27:c:31:LEU:HD21	28:d:26:LEU:HD21	1.81	0.62
28:d:100:ASP:OD2	32:h:117:ARG:NH1	2.32	0.62
32:h:106:LYS:O	32:h:110:VAL:HG12	2.00	0.62
18:S:17:GLU:OE1	18:S:67:ARG:NH2	2.33	0.62
6:F:174:ASP:OD2	42:s:57:LEU:HD12	2.00	0.62
7:G:572:THR:HG1	7:G:633:TYR:HH	1.46	0.62
33:i:1:SAC:HB3	33:i:1:SAC:H2A1	1.81	0.62
5:E:67:ASN:ND2	5:E:81:TYR:OH	2.33	0.61
24:Z:96:ILE:HD11	29:e:91:TYR:CZ	2.35	0.61
6:F:143:TYR:OH	42:s:42:THR:O	2.13	0.61
36:l:70:MET:HE3	36:l:80:LEU:HD13	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:l:122:PRO:O	39:o:8:TYR:OH	2.06	0.61
40:p:105:ILE:CD1	40:p:134:VAL:HG21	2.30	0.61
8:H:181:MET:HE1	8:H:304:HIS:NE2	2.15	0.61
14:N:34:GLU:O	14:N:38:LEU:HD12	2.01	0.61
7:G:572:THR:OG1	7:G:633:TYR:OH	2.05	0.61
15:O:318:TRP:CD2	15:O:319:LEU:HD23	2.36	0.61
16:P:81:ILE:O	16:P:85:VAL:HG22	2.00	0.61
5:E:101:GLN:NE2	5:E:155:GLN:OE1	2.34	0.61
7:G:37:ILE:O	7:G:39:ARG:NH1	2.33	0.61
9:I:76:GLU:OE1	9:I:76:GLU:HA	2.00	0.61
12:L:203:PHE:CZ	40:p:109:LEU:HD11	2.36	0.61
19:T:45:LEU:O	19:T:48:VAL:HG12	2.01	0.61
38:n:34:ASP:OD1	38:n:35:LYS:N	2.33	0.61
3:C:66:ASP:HB2	20:V:89:LEU:HD22	1.83	0.60
7:G:180:ASP:OD2	7:G:180:ASP:O	2.20	0.60
5:E:119:THR:HG21	5:E:166:PRO:HB3	1.83	0.60
7:G:453:LEU:O	7:G:493:LEU:N	2.34	0.60
2:B:92:GLN:NE2	8:H:212:ASN:O	2.34	0.60
7:G:201:ASP:OD2	7:G:268:ARG:NH2	2.33	0.60
8:H:189:THR:HG22	8:H:234:MET:CE	2.31	0.60
36:l:66:SER:O	36:l:69:ARG:N	2.33	0.60
14:N:17:PRO:O	14:N:21:MET:HG3	2.01	0.60
24:Z:142:TYR:HE1	25:a:36:LYS:HZ1	1.49	0.60
7:G:11:VAL:HG11	7:G:73:VAL:CG2	2.32	0.60
4:D:61:VAL:HG23	4:D:61:VAL:O	2.02	0.60
7:G:338:VAL:O	7:G:339:ASP:CB	2.50	0.60
31:g:118:LEU:HD11	40:p:163:LEU:HG	1.83	0.60
3:C:80:ALA:HB3	3:C:138:PHE:CD1	2.37	0.60
7:G:145:LEU:HD22	7:G:269:PHE:CD2	2.37	0.60
37:m:99:PHE:O	37:m:103:VAL:HG23	2.01	0.60
32:h:117:ARG:NH2	32:h:119:ASP:OD2	2.35	0.59
9:I:76:GLU:OE2	17:R:44:ILE:N	2.35	0.59
11:K:33:LEU:HD23	11:K:36:MET:HE2	1.84	0.59
15:O:12:ASP:OD1	15:O:17:LYS:NZ	2.32	0.59
33:i:116:ILE:HG23	33:i:117:PRO:HD2	1.84	0.59
5:E:109:MET:SD	6:F:349:ARG:HG2	2.42	0.59
2:B:126:MET:SD	2:B:166:LEU:HD22	2.42	0.59
7:G:11:VAL:HG11	7:G:73:VAL:HG22	1.83	0.59
14:N:200:LEU:HD12	14:N:269:GLU:HG3	1.84	0.59
14:N:243:MET:HE2	14:N:330:ALA:HB1	1.84	0.59
21:W:119:ASP:OD1	21:W:120:PHE:N	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:156:GLY:HA2	12:L:164:ALA:HB1	1.84	0.59
4:D:413:ASP:OD1	8:H:281:ARG:NH1	2.35	0.59
7:G:415:LEU:C	7:G:416:THR:HG1	2.07	0.59
12:L:136:ASN:ND2	12:L:197:GLU:OE2	2.36	0.59
12:L:146:GLY:O	12:L:149:ILE:HG22	2.03	0.59
7:G:294:THR:OG1	7:G:298:ASP:OD1	2.18	0.59
9:I:155:LYS:HA	9:I:155:LYS:HE2	1.84	0.59
5:E:78:MET:HE2	7:G:173:GLY:CA	2.33	0.58
4:D:73:VAL:HG21	4:D:414:VAL:HG21	1.85	0.58
15:O:170:ILE:HD11	15:O:238:ILE:HD11	1.84	0.58
14:N:218:ALA:CB	14:N:244:ILE:HD11	2.33	0.58
8:H:181:MET:HE1	8:H:304:HIS:CD2	2.38	0.58
35:k:48:ASP:O	35:k:48:ASP:OD1	2.22	0.58
4:D:151:ILE:O	4:D:155:THR:OG1	2.20	0.58
7:G:189:LYS:NZ	7:G:190:MET:O	2.37	0.58
12:L:357:ARG:O	12:L:436:ARG:NE	2.31	0.58
38:n:56:MET:HE3	38:n:56:MET:HA	1.86	0.58
7:G:498:SER:O	7:G:498:SER:OG	2.22	0.58
6:F:42:TRP:CE2	6:F:161:LEU:HD13	2.38	0.58
8:H:260:MET:SD	8:H:264:LEU:HD22	2.44	0.58
13:M:76:MET:CE	13:M:99:LEU:HD22	2.34	0.58
33:i:82:HIS:CE1	40:p:36:TYR:HH	2.21	0.58
7:G:324:ASP:OD1	7:G:324:ASP:N	2.32	0.58
7:G:415:LEU:O	7:G:416:THR:OG1	2.03	0.58
13:M:318:ALA:HB2	13:M:373:ILE:HG13	1.85	0.58
19:U:19:LEU:HB3	19:U:30:LEU:HD11	1.85	0.58
6:F:137:TYR:CG	6:F:192:LEU:HD11	2.38	0.58
6:F:347:ILE:HD11	6:F:414:PRO:CB	2.34	0.57
12:L:562:ILE:HB	12:L:563:PRO:HD3	1.85	0.57
22:X:46:ARG:NH1	24:Z:75:GLN:OE1	2.37	0.57
10:J:59:TYR:CZ	11:K:34:GLU:OE1	2.57	0.57
10:J:151:MET:HA	10:J:151:MET:HE2	1.86	0.57
13:M:324:SER:OG	13:M:440:HIS:NE2	2.37	0.57
15:O:41:GLU:OE1	15:O:45:GLN:NE2	2.36	0.57
13:M:285:LEU:HD22	13:M:410:MET:SD	2.43	0.57
14:N:25:ASN:OD1	14:N:27:MET:N	2.38	0.57
28:d:112:ILE:O	40:p:159:LYS:NZ	2.38	0.57
32:h:110:VAL:CG2	32:h:114:MET:HE3	2.34	0.57
4:D:293:CYS:SG	4:D:420:THR:HG21	2.44	0.57
15:O:170:ILE:CD1	15:O:238:ILE:HD11	2.33	0.57
3:C:102:ASN:OD1	3:C:102:ASN:O	2.23	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:W:44:GLU:OE2	21:W:48:THR:OG1	2.21	0.57
12:L:162:THR:O	12:L:166:THR:HG23	2.05	0.57
40:p:89:MET:HE3	40:p:89:MET:HA	1.86	0.57
12:L:89:PHE:HD2	12:L:132:THR:HG21	1.69	0.57
14:N:240:MET:HE3	14:N:326:PHE:CZ	2.40	0.57
19:U:11:ILE:HD13	19:U:80:ILE:HG22	1.87	0.57
40:p:128:LEU:HD23	40:p:128:LEU:O	2.05	0.57
2:B:55:LEU:HD12	2:B:55:LEU:O	2.04	0.57
7:G:55:CYS:SG	7:G:68:ALA:N	2.78	0.57
7:G:118:ASP:O	7:G:122:MET:HG2	2.04	0.57
11:K:24:SER:OG	11:K:25:HIS:N	2.37	0.57
12:L:424:MET:HA	12:L:424:MET:HE2	1.85	0.56
28:d:119:VAL:HG22	28:d:119:VAL:O	2.05	0.56
1:A:92:PHE:HA	8:H:302:MET:HE3	1.86	0.56
7:G:522:LEU:HD21	7:G:524:LEU:HD21	1.88	0.56
13:M:222:GLU:N	13:M:222:GLU:OE1	2.38	0.56
5:E:89:MET:HG2	6:F:181:ALA:HB1	1.88	0.56
7:G:194:GLU:OE1	7:G:194:GLU:N	2.26	0.56
8:H:20:LEU:HA	25:a:12:MET:CE	2.36	0.56
12:L:237:MET:HE3	12:L:303:ALA:HB2	1.86	0.56
16:P:63:GLY:HA3	16:P:68:LEU:HD13	1.88	0.56
36:l:31:GLU:N	36:l:31:GLU:OE1	2.36	0.56
9:I:15:ASP:OD1	9:I:16:MET:N	2.38	0.56
14:N:137:ALA:HB3	14:N:138:PRO:HD3	1.86	0.56
19:U:86:VAL:HG23	19:U:87:TYR:N	2.20	0.56
33:i:68:LEU:HD12	33:i:68:LEU:H	1.70	0.56
5:E:156:ILE:O	5:E:158:ASP:N	2.38	0.56
1:A:61:THR:HG21	1:A:105:GLU:OE1	2.05	0.56
14:N:53:THR:O	14:N:57:THR:HG23	2.05	0.56
2:B:50:ALA:HB2	8:H:53:MET:HE2	1.86	0.56
6:F:297:VAL:HG12	6:F:336:VAL:HG12	1.86	0.56
12:L:22:SER:HA	12:L:27:ILE:HD12	1.85	0.56
12:L:83:ASP:OD1	12:L:83:ASP:N	2.38	0.56
12:L:377:SER:OG	12:L:427:ILE:HD11	2.05	0.56
25:a:14:VAL:O	25:a:18:ILE:HG13	2.04	0.56
25:a:54:ILE:HD11	29:e:105:PRO:HB3	1.86	0.56
36:l:80:LEU:O	36:l:84:ILE:HG23	2.05	0.56
5:E:82:GLU:OE2	7:G:174:THR:N	2.37	0.56
40:p:47:VAL:O	40:p:51:ILE:HG13	2.05	0.56
8:H:318:MET:HE2	8:H:318:MET:HA	1.88	0.56
7:G:197:GLY:N	7:G:265:ASP:OD2	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:366:THR:OG1	7:G:491:ASN:ND2	2.39	0.55
16:P:90:VAL:HG11	16:P:218:THR:HG22	1.87	0.55
16:P:301:GLU:OE1	16:P:301:GLU:N	2.37	0.55
37:m:27:GLU:OE2	37:m:27:GLU:N	2.33	0.55
5:E:48:LEU:HB2	5:E:49:PRO:HD3	1.88	0.55
18:S:95:SER:O	18:S:95:SER:OG	2.21	0.55
7:G:274:LEU:O	7:G:278:ARG:NH1	2.37	0.55
10:J:63:MET:HE2	10:J:63:MET:N	2.21	0.55
1:A:112:GLU:OE2	8:H:291:LYS:NZ	2.37	0.55
5:E:199:LEU:HD12	6:F:26:TYR:CD2	2.41	0.55
31:g:65:SER:OG	32:h:28:TYR:O	2.18	0.55
2:B:180:GLU:O	2:B:181:GLN:HG2	2.06	0.55
6:F:111:MET:HE1	6:F:138:ILE:HD11	1.88	0.55
7:G:334:LEU:HD12	7:G:609:ILE:HG21	1.87	0.55
13:M:173:THR:HB	32:h:101:ALA:HB2	1.88	0.55
13:M:278:ARG:NH2	36:l:81:ASP:OD1	2.35	0.55
31:g:48:ASP:OD2	31:g:51:VAL:HG23	2.07	0.55
2:B:57:PRO:HD3	2:B:84:VAL:HG13	1.88	0.55
12:L:304:PHE:CZ	12:L:526:LEU:HD22	2.41	0.55
13:M:23:THR:O	13:M:27:VAL:HG23	2.06	0.55
16:P:143:SER:OG	16:P:282:ASP:OD1	2.22	0.55
23:Y:31:GLY:O	23:Y:35:ILE:HD12	2.06	0.55
2:B:46:LEU:CD1	8:H:53:MET:HE1	2.37	0.55
6:F:268:VAL:HG21	6:F:283:HIS:CD2	2.42	0.55
7:G:82:ASN:OD1	7:G:82:ASN:O	2.24	0.55
14:N:218:ALA:CB	14:N:240:MET:HE1	2.35	0.55
7:G:665:GLN:N	7:G:665:GLN:OE1	2.40	0.55
11:K:57:MET:N	11:K:58:PRO:CD	2.70	0.55
5:E:46:ALA:HB1	5:E:50:VAL:HG23	1.87	0.55
6:F:66:ARG:O	6:F:68:ARG:NH1	2.40	0.55
6:F:97:ALA:CB	6:F:111:MET:HE2	2.37	0.55
9:I:82:CYS:SG	9:I:83:LYS:N	2.80	0.54
10:J:33:ILE:HD12	10:J:60:LEU:HD23	1.90	0.54
4:D:326:ASP:OD1	7:G:127:ARG:NH2	2.41	0.54
7:G:371:VAL:HG22	7:G:450:MET:HE1	1.88	0.54
12:L:132:THR:HG22	12:L:258:PHE:CE2	2.42	0.54
19:U:11:ILE:HD11	19:U:86:VAL:HG11	1.90	0.54
24:Z:92:GLU:O	24:Z:96:ILE:HG12	2.08	0.54
6:F:329:LEU:HD23	6:F:329:LEU:O	2.08	0.54
8:H:283:ASP:OD1	8:H:284:GLN:N	2.39	0.54
9:I:101:ARG:HD2	17:R:41:ILE:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:H:402:PC1:H242	10:J:30:LEU:HD21	1.90	0.54
12:L:378:LEU:HD22	12:L:383:MET:HE2	1.90	0.54
13:M:210:TYR:O	13:M:213:HIS:ND1	2.40	0.54
19:T:57:GLU:C	19:T:57:GLU:OE2	2.50	0.54
39:o:48:ALA:HB3	39:o:50:LEU:HD13	1.88	0.54
4:D:405:MET:HE1	4:D:421:GLN:CD	2.33	0.54
5:E:204:GLU:OE1	6:F:16:LYS:NZ	2.26	0.54
6:F:372:MET:HG2	6:F:392:LEU:HD11	1.89	0.54
44:H:402:PC1:H371	10:J:92:LEU:HD21	1.88	0.54
12:L:3:ILE:HD13	40:p:36:TYR:CD2	2.43	0.54
14:N:275:CYS:SG	14:N:278:MET:HB2	2.48	0.54
29:e:14:ASP:OD1	29:e:14:ASP:C	2.51	0.54
31:g:65:SER:O	32:h:32:THR:HG21	2.07	0.54
33:i:1:SAC:H2A1	33:i:1:SAC:CB	2.38	0.54
36:l:97:TRP:CD1	36:l:101:CYS:HG	2.26	0.54
12:L:149:ILE:HD11	13:M:364:LEU:HD21	1.90	0.54
12:L:257:ILE:HB	12:L:329:ILE:HD11	1.89	0.54
15:O:175:PRO:O	15:O:179:VAL:HG23	2.08	0.54
15:O:219:GLU:OE1	15:O:241:LEU:HD22	2.08	0.54
7:G:73:VAL:O	7:G:74:MET:HE2	2.08	0.54
7:G:684:MET:HE3	7:G:684:MET:HA	1.88	0.54
11:K:26:LEU:HD23	11:K:78:LEU:HD12	1.90	0.54
25:a:40:ARG:O	25:a:41:VAL:HG13	2.08	0.54
4:D:208:LEU:HD22	4:D:315:LEU:CD2	2.38	0.54
7:G:663:PRO:O	7:G:665:GLN:N	2.41	0.54
12:L:193:MET:HE2	12:L:200:GLN:O	2.08	0.54
14:N:135:LYS:CE	14:N:187:MET:HE2	2.38	0.54
31:g:69:VAL:O	31:g:73:THR:HG23	2.07	0.54
36:l:14:ARG:N	36:l:18:GLU:OE1	2.35	0.54
7:G:174:THR:OG1	7:G:181:MET:HE3	2.08	0.53
14:N:214:PRO:HB2	14:N:247:MET:SD	2.48	0.53
6:F:108:ARG:O	6:F:112:ARG:HG2	2.08	0.53
10:J:62:GLY:C	10:J:63:MET:HE2	2.33	0.53
9:I:59:LEU:HD22	9:I:63:PHE:CD1	2.43	0.53
12:L:122:THR:O	12:L:126:ILE:HD12	2.08	0.53
23:Y:139:THR:HG22	23:Y:140:PRO:HD2	1.89	0.53
36:l:81:ASP:OD1	36:l:81:ASP:N	2.41	0.53
10:J:77:GLU:OE2	16:P:324:TYR:OH	2.24	0.53
12:L:180:ILE:HD12	13:M:397:GLY:HA2	1.89	0.53
14:N:44:LEU:HG	14:N:122:ILE:HG21	1.91	0.53
17:R:53:GLU:N	17:R:53:GLU:OE1	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:S:22:LEU:CD2	18:S:54:ILE:HD11	2.39	0.53
4:D:343:GLU:O	4:D:347:HIS:ND1	2.42	0.53
7:G:111:GLY:O	7:G:117:GLN:NE2	2.42	0.53
42:s:41:ASN:OD1	42:s:43:TYR:N	2.42	0.53
8:H:189:THR:HG22	8:H:234:MET:HE3	1.89	0.53
13:M:360:LEU:O	13:M:364:LEU:HG	2.09	0.53
8:H:73:ILE:O	8:H:76:THR:OG1	2.26	0.53
8:H:144:VAL:HG23	8:H:145:THR:HG23	1.90	0.53
15:O:221:LEU:HD21	15:O:241:LEU:HD21	1.90	0.53
53:U:201:EHZ:O5	53:U:201:EHZ:O6	2.26	0.53
27:c:4:VAL:HG23	27:c:5:ARG:H	1.73	0.53
4:D:106:LEU:O	4:D:427:GLU:N	2.41	0.53
4:D:337:GLU:O	4:D:344:SER:OG	2.27	0.53
7:G:191:PHE:C	7:G:192:MET:HE2	2.34	0.53
8:H:31:MET:HG3	9:I:43:LEU:HD13	1.91	0.53
12:L:407:TRP:O	12:L:411:ILE:HG12	2.09	0.53
18:S:14:GLY:O	18:S:69:ALA:N	2.41	0.53
19:T:61:GLU:OE1	19:T:62:ILE:N	2.42	0.53
1:A:79:ILE:HG12	1:A:87:MET:HE2	1.91	0.53
1:A:88:MET:SD	8:H:309:ILE:HG21	2.49	0.53
38:n:116:ASP:O	38:n:119:SER:OG	2.24	0.53
39:o:73:PHE:CD1	39:o:74:PRO:HA	2.44	0.53
2:B:41:THR:O	2:B:44:ASP:OD1	2.26	0.52
6:F:245:PHE:O	6:F:252:GLY:N	2.42	0.52
12:L:145:GLU:OE1	13:M:369:LEU:HD12	2.09	0.52
12:L:173:LEU:HD12	13:M:405:MET:CE	2.40	0.52
16:P:175:GLU:OE1	16:P:175:GLU:N	2.40	0.52
7:G:205:VAL:O	7:G:207:ALA:N	2.37	0.52
7:G:568:GLU:O	7:G:585:VAL:HG13	2.08	0.52
11:K:66:PHE:CZ	14:N:31:VAL:HG13	2.43	0.52
9:I:76:GLU:OE2	17:R:40:ALA:O	2.27	0.52
48:M:502:CDL:C52	28:d:53:VAL:HG11	2.40	0.52
13:M:166:LEU:HD23	13:M:166:LEU:O	2.09	0.52
15:O:174:VAL:HG23	15:O:179:VAL:HG22	1.90	0.52
19:U:15:VAL:HG23	19:U:54:MET:SD	2.50	0.52
4:D:161:ILE:HD11	4:D:416:ALA:CB	2.39	0.52
7:G:331:LEU:HD13	7:G:525:LEU:HD22	1.91	0.52
15:O:221:LEU:HD11	15:O:241:LEU:HD11	1.91	0.52
37:m:91:VAL:HG13	37:m:92:ALA:N	2.24	0.52
1:A:15:LEU:HD23	1:A:19:LEU:HD13	1.92	0.52
2:B:55:LEU:HD13	2:B:95:VAL:CG1	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:297:GLU:OE1	21:W:125:TYR:CE1	2.62	0.52
8:H:63:PRO:HB2	8:H:66:THR:HG23	1.92	0.52
15:O:198:TYR:CE2	15:O:202:ILE:HD11	2.45	0.52
15:O:315:LYS:O	15:O:316:TRP:HB2	2.10	0.52
24:Z:128:THR:HG22	24:Z:131:GLU:HG3	1.91	0.52
6:F:148:ASN:O	6:F:151:VAL:HG22	2.09	0.52
12:L:223:LYS:HG2	12:L:255:ALA:HB3	1.91	0.52
13:M:14:LEU:HD21	13:M:26:ASN:HB3	1.91	0.52
19:U:55:GLU:HG2	19:U:61:GLU:HA	1.92	0.52
37:m:24:VAL:HG12	37:m:24:VAL:O	2.08	0.52
5:E:53:LEU:HD22	42:s:46:LEU:HD13	1.92	0.52
12:L:99:SER:OG	12:L:457:LEU:HD11	2.10	0.52
13:M:195:LEU:O	13:M:196:TRP:C	2.53	0.52
13:M:369:LEU:O	13:M:372:SER:OG	2.14	0.52
34:j:61:ASP:HA	34:j:66:ILE:HD12	1.90	0.52
3:C:69:ASN:CG	20:V:50:ILE:HD11	2.35	0.52
6:F:42:TRP:CD2	6:F:161:LEU:HD13	2.45	0.52
8:H:4:ILE:O	8:H:8:THR:HG23	2.09	0.52
8:H:148:ILE:HG21	8:H:297:THR:HG22	1.92	0.52
10:J:124:LEU:HD23	10:J:124:LEU:C	2.35	0.52
23:Y:86:ASP:C	23:Y:86:ASP:OD1	2.52	0.52
27:c:37:GLU:OE2	27:c:37:GLU:HA	2.10	0.52
33:i:75:LEU:HD21	33:i:79:TRP:CZ2	2.45	0.52
4:D:19:MET:HG2	14:N:173:THR:HG23	1.92	0.52
5:E:78:MET:HE2	7:G:173:GLY:HA3	1.90	0.52
21:W:121:LEU:HD23	21:W:121:LEU:O	2.10	0.52
7:G:296:TRP:HB3	21:W:125:TYR:HE2	1.75	0.51
8:H:260:MET:SD	8:H:260:MET:C	2.92	0.51
10:J:168:GLU:OE2	14:N:5:THR:HG21	2.11	0.51
12:L:3:ILE:HD13	40:p:36:TYR:CG	2.45	0.51
12:L:350:LEU:O	12:L:352:ASP:N	2.42	0.51
13:M:414:THR:HG22	13:M:415:GLN:OE1	2.10	0.51
14:N:296:LEU:O	14:N:300:THR:HG23	2.10	0.51
33:i:31:GLU:O	33:i:33:VAL:N	2.43	0.51
8:H:85:LEU:HB3	8:H:233:LEU:HD21	1.92	0.51
12:L:430:VAL:O	12:L:430:VAL:HG12	2.09	0.51
27:c:4:VAL:HG23	27:c:5:ARG:N	2.26	0.51
32:h:10:VAL:HG23	32:h:10:VAL:O	2.10	0.51
39:o:102:GLU:OE2	39:o:106:ARG:NE	2.43	0.51
10:J:27:TYR:CE2	10:J:81:THR:HG22	2.46	0.51
19:T:74:GLN:O	19:T:77:VAL:HG22	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:6:LEU:O	5:E:92:ARG:NH2	2.39	0.51
6:F:30:ASP:O	6:F:39:ARG:NH2	2.44	0.51
21:W:72:ASN:O	21:W:75:VAL:HG22	2.11	0.51
22:X:164:THR:O	22:X:166:PHE:N	2.43	0.51
33:i:4:THR:OG1	33:i:7:GLU:OE2	2.28	0.51
35:k:34:LEU:HD23	35:k:34:LEU:O	2.10	0.51
5:E:44:ALA:O	5:E:47:VAL:HG23	2.10	0.51
11:K:59:ILE:HG23	14:N:27:MET:HE2	1.93	0.51
12:L:15:LEU:CD1	12:L:129:LEU:HD12	2.39	0.51
15:O:36:ASN:ND2	50:O:401:GTP:O1B	2.44	0.51
15:O:320:LYS:HD2	27:c:9:ASN:OD1	2.10	0.51
4:D:222:ILE:HD11	4:D:301:VAL:HG13	1.92	0.51
12:L:69:VAL:HG12	12:L:76:LEU:HD12	1.92	0.51
12:L:140:LEU:O	12:L:140:LEU:HD23	2.10	0.51
14:N:221:LEU:HD11	14:N:319:LYS:HD3	1.93	0.51
6:F:295:LEU:HD22	6:F:343:ILE:HD13	1.92	0.51
23:Y:54:VAL:O	23:Y:57:VAL:HG12	2.10	0.51
29:e:11:ILE:HD12	29:e:13:LEU:HD21	1.92	0.51
35:k:48:ASP:OD1	35:k:48:ASP:C	2.52	0.51
6:F:277:LYS:O	6:F:281:GLU:OE1	2.29	0.51
12:L:433:THR:HG22	12:L:434:LYS:N	2.25	0.51
15:O:62:THR:HG23	15:O:63:THR:HG23	1.93	0.51
23:Y:127:LYS:O	23:Y:131:LEU:HD13	2.11	0.51
1:A:63:LEU:HD21	10:J:62:GLY:HA2	1.91	0.51
8:H:250:ASN:OD1	8:H:250:ASN:C	2.54	0.51
7:G:673:MET:O	7:G:674:THR:HG23	2.10	0.51
8:H:207:LEU:O	8:H:208:VAL:C	2.54	0.51
10:J:67:PHE:O	10:J:71:THR:HG23	2.11	0.51
16:P:121:SER:HB2	16:P:126:VAL:HG11	1.92	0.51
31:g:26:GLU:OE1	31:g:27:ASP:N	2.39	0.51
6:F:24:ASN:OD1	6:F:24:ASN:O	2.28	0.50
7:G:668:ILE:HG23	7:G:691:VAL:HG21	1.92	0.50
12:L:536:ILE:HG23	12:L:537:THR:N	2.26	0.50
13:M:367:LEU:O	13:M:368:ALA:HB3	2.10	0.50
15:O:171:TYR:CD2	15:O:211:LEU:HD11	2.46	0.50
1:A:95:VAL:HG11	8:H:302:MET:CG	2.41	0.50
4:D:223:ASP:OD1	4:D:305:ARG:NH2	2.35	0.50
6:F:136:ILE:HG22	6:F:138:ILE:CD1	2.41	0.50
7:G:156:CYS:SG	7:G:158:ARG:HD3	2.51	0.50
8:H:171:HIS:O	24:Z:48:ARG:NH2	2.44	0.50
9:I:27:LEU:HD23	47:I:201:3PE:O32	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:112:ALA:HB1	13:M:117:MET:HB3	1.91	0.50
13:M:207:MET:O	13:M:208:PRO:C	2.53	0.50
13:M:221:VAL:HG11	13:M:284:SER:OG	2.11	0.50
5:E:202:LEU:HD21	6:F:20:ARG:HD2	1.93	0.50
6:F:33:LEU:HD12	6:F:33:LEU:O	2.11	0.50
7:G:294:THR:HG1	7:G:298:ASP:CG	2.19	0.50
7:G:315:VAL:HG23	7:G:315:VAL:O	2.11	0.50
12:L:101:MET:SD	12:L:118:ILE:HG23	2.52	0.50
13:M:347:GLY:N	13:M:414:THR:O	2.44	0.50
14:N:108:LEU:CD1	14:N:187:MET:HB3	2.40	0.50
5:E:100:ILE:HG21	5:E:120:LEU:HD13	1.93	0.50
6:F:270:GLU:OE1	6:F:283:HIS:NE2	2.44	0.50
12:L:260:LEU:HD11	12:L:314:MET:HE1	1.94	0.50
31:g:56:MET:SD	31:g:56:MET:C	2.95	0.50
36:l:93:THR:OG1	36:l:95:VAL:O	2.26	0.50
4:D:38:ILE:HD11	15:O:157:ARG:CZ	2.42	0.50
4:D:342:MET:HE1	7:G:103:LEU:HA	1.93	0.50
6:F:264:HIS:O	6:F:284:ALA:HB1	2.12	0.50
8:H:113:VAL:HG21	8:H:139:THR:OG1	2.12	0.50
11:K:17:LEU:CD2	11:K:21:MET:SD	3.00	0.50
19:U:9:ASP:N	19:U:9:ASP:OD1	2.43	0.50
30:f:42:LEU:HB2	40:p:70:VAL:HG11	1.92	0.50
33:i:94:TYR:OH	40:p:10:PRO:O	2.29	0.50
6:F:174:ASP:CG	42:s:57:LEU:HD12	2.37	0.50
7:G:296:TRP:CZ3	7:G:561:LEU:HD22	2.47	0.50
9:I:154:GLU:O	9:I:158:ASN:OD1	2.29	0.50
11:K:91:GLN:OE1	11:K:91:GLN:N	2.37	0.50
12:L:290:ILE:HG13	12:L:418:MET:HE1	1.94	0.50
12:L:601:LEU:C	12:L:601:LEU:HD23	2.37	0.50
13:M:221:VAL:HG23	13:M:222:GLU:OE1	2.11	0.50
13:M:310:MET:HE1	13:M:380:PHE:CG	2.47	0.50
5:E:52:ASP:OD2	6:F:179:ARG:NH1	2.43	0.50
6:F:60:MET:HE2	6:F:60:MET:HA	1.94	0.50
7:G:575:ASN:OD1	7:G:579:ARG:O	2.30	0.50
13:M:114:GLU:OE2	13:M:174:LEU:HB2	2.12	0.50
15:O:319:LEU:HD12	27:c:17:VAL:HG11	1.92	0.50
16:P:219:LYS:O	16:P:220:ASP:C	2.55	0.50
26:b:78:ASP:N	26:b:78:ASP:OD2	2.45	0.50
27:c:9:ASN:OD1	27:c:9:ASN:O	2.30	0.50
1:A:69:ILE:CD1	8:H:148:ILE:HD11	2.41	0.50
4:D:8:ILE:O	4:D:12:GLU:HG2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:97:ALA:HB2	6:F:225:VAL:CG2	2.41	0.50
12:L:233:LEU:HD23	12:L:233:LEU:C	2.37	0.50
4:D:161:ILE:HG23	4:D:238:ARG:HG2	1.94	0.50
6:F:24:ASN:O	6:F:24:ASN:CG	2.55	0.50
34:j:12:ARG:NH1	35:k:53:ASN:O	2.40	0.50
7:G:222:THR:HG22	7:G:223:ARG:O	2.12	0.49
13:M:1:FME:HE2	13:M:111:SER:CB	2.42	0.49
19:U:19:LEU:CB	19:U:30:LEU:HD11	2.41	0.49
37:m:18:ASP:OD1	37:m:20:ALA:N	2.45	0.49
5:E:123:LYS:HG2	5:E:173:ILE:HG21	1.94	0.49
12:L:273:ILE:O	12:L:277:MET:HG3	2.12	0.49
12:L:546:LEU:HB2	13:M:278:ARG:HD3	1.94	0.49
14:N:148:LEU:HD12	14:N:148:LEU:N	2.27	0.49
15:O:319:LEU:CD1	27:c:17:VAL:HG11	2.42	0.49
16:P:295:THR:C	16:P:296:ASP:OD1	2.54	0.49
28:d:16:ASP:O	28:d:16:ASP:OD2	2.30	0.49
8:H:260:MET:O	8:H:264:LEU:HD13	2.11	0.49
21:W:77:ASP:HB3	21:W:80:VAL:HG23	1.94	0.49
4:D:44:VAL:O	4:D:44:VAL:HG13	2.11	0.49
12:L:230:HIS:N	12:L:231:PRO:CD	2.76	0.49
13:M:195:LEU:O	13:M:198:ALA:N	2.45	0.49
36:l:29:ARG:HH12	37:m:31:ALA:HB1	1.77	0.49
2:B:143:VAL:O	2:B:145:GLY:N	2.45	0.49
6:F:403:THR:HG21	6:F:408:GLY:HA3	1.95	0.49
7:G:340:SER:OG	7:G:341:ASP:N	2.43	0.49
12:L:64:THR:HG21	12:L:77:LYS:HE3	1.94	0.49
18:S:19:ARG:HG2	18:S:65:TRP:HB2	1.92	0.49
19:T:51:ILE:HD13	19:T:67:ALA:HB1	1.93	0.49
19:T:73:PRO:O	19:T:77:VAL:HG13	2.12	0.49
27:c:28:TRP:NE1	28:d:66:THR:OG1	2.44	0.49
30:f:42:LEU:O	30:f:43:PHE:CB	2.60	0.49
34:j:65:GLY:O	39:o:103:ARG:NH1	2.45	0.49
8:H:200:LEU:N	8:H:200:LEU:HD12	2.28	0.49
10:J:32:LEU:HD22	10:J:63:MET:SD	2.52	0.49
12:L:160:GLY:N	38:n:95:CYS:SG	2.83	0.49
12:L:381:THR:O	12:L:381:THR:OG1	2.30	0.49
19:T:49:GLU:OE2	21:W:38:TYR:OH	2.16	0.49
26:b:40:TYR:CD2	26:b:83:LEU:HD11	2.47	0.49
34:j:60:THR:O	34:j:64:LEU:HD12	2.13	0.49
4:D:87:THR:HG21	4:D:106:LEU:HD21	1.95	0.49
7:G:160:ILE:HD11	7:G:183:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:597:TRP:CD2	7:G:619:VAL:HG21	2.48	0.49
12:L:63:ILE:CG2	33:i:89:MET:HE1	2.42	0.49
38:n:31:ILE:HD12	38:n:31:ILE:H	1.78	0.49
15:O:214:MET:HE2	15:O:214:MET:HA	1.93	0.49
18:S:88:ARG:O	18:S:89:ALA:C	2.56	0.49
22:X:17:VAL:HG23	22:X:18:LYS:N	2.28	0.49
33:i:82:HIS:CG	40:p:36:TYR:HH	2.28	0.49
39:o:104:GLU:OE1	39:o:104:GLU:C	2.56	0.49
4:D:412:ALA:HA	8:H:206:GLU:OE1	2.13	0.49
12:L:353:GLU:HA	12:L:353:GLU:OE1	2.12	0.49
13:M:23:THR:HG23	13:M:24:TRP:N	2.28	0.49
16:P:166:ILE:HD13	16:P:228:PHE:HB2	1.94	0.49
29:e:40:ILE:HD11	32:h:128:PRO:HB2	1.93	0.49
30:f:42:LEU:O	30:f:43:PHE:HB2	2.12	0.49
30:f:47:LEU:H	30:f:47:LEU:HD12	1.78	0.49
5:E:26:GLU:N	5:E:26:GLU:OE1	2.45	0.49
7:G:315:VAL:HG23	7:G:340:SER:HB3	1.95	0.49
10:J:117:LEU:O	10:J:117:LEU:HD23	2.13	0.49
11:K:20:LEU:HD12	12:L:592:LEU:CD1	2.43	0.49
19:U:57:GLU:OE1	34:j:11:TYR:OH	2.21	0.49
22:X:43:MET:HE1	24:Z:75:GLN:HG3	1.94	0.49
10:J:44:LEU:HD23	10:J:49:SER:HA	1.95	0.48
13:M:119:TYR:CZ	13:M:123:GLU:OE1	2.66	0.48
4:D:9:GLU:HA	4:D:9:GLU:OE2	2.12	0.48
4:D:134:ALA:HB2	4:D:323:ILE:O	2.12	0.48
6:F:257:ASN:OD1	6:F:257:ASN:N	2.47	0.48
6:F:347:ILE:HD12	6:F:350:LEU:HD12	1.94	0.48
13:M:207:MET:HB3	13:M:298:ILE:HD11	1.94	0.48
34:j:41:TRP:CE2	35:k:87:LEU:HD21	2.48	0.48
1:A:69:ILE:HD12	8:H:148:ILE:HD11	1.96	0.48
2:B:153:ASP:OD1	2:B:153:ASP:N	2.47	0.48
4:D:335:ARG:NH2	9:I:128:CYS:O	2.41	0.48
10:J:168:GLU:OE2	14:N:5:THR:OG1	2.31	0.48
12:L:319:MET:O	12:L:321:GLN:N	2.45	0.48
12:L:424:MET:HE1	35:k:68:PHE:CE1	2.48	0.48
16:P:215:VAL:O	16:P:218:THR:OG1	2.30	0.48
31:g:35:VAL:HG11	31:g:42:PHE:CZ	2.47	0.48
5:E:105:THR:HG22	5:E:106:THR:H	1.78	0.48
7:G:569:LYS:HG3	7:G:571:ALA:HB2	1.96	0.48
39:o:14:VAL:HG12	39:o:109:GLN:HA	1.96	0.48
7:G:426:PRO:HB3	7:G:661:LEU:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:528:ASP:OD1	7:G:545:TYR:OH	2.31	0.48
9:I:176:LEU:H	9:I:176:LEU:HD12	1.78	0.48
14:N:70:ILE:HG13	14:N:101:ALA:HB3	1.96	0.48
22:X:164:THR:OG1	28:d:83:TYR:CZ	2.66	0.48
31:g:69:VAL:HG11	32:h:32:THR:OG1	2.14	0.48
33:i:82:HIS:CE1	40:p:41:ASP:OD1	2.67	0.48
40:p:42:TRP:CD1	40:p:46:LEU:HD21	2.49	0.48
5:E:109:MET:O	5:E:112:ASP:N	2.42	0.48
6:F:82:MET:HE3	6:F:129:MET:HE3	1.95	0.48
9:I:106:ARG:O	9:I:107:ARG:HD3	2.13	0.48
22:X:164:THR:O	22:X:164:THR:CG2	2.62	0.48
29:e:67:LEU:O	29:e:68:ARG:HB2	2.13	0.48
8:H:111:LEU:HD13	10:J:60:LEU:HD21	1.95	0.48
8:H:293:PHE:O	8:H:297:THR:N	2.44	0.48
14:N:5:THR:HG22	14:N:9:ILE:HD12	1.95	0.48
17:R:57:ILE:HD12	17:R:58:ALA:H	1.78	0.48
22:X:150:ASN:OD1	29:e:50:ARG:NH1	2.44	0.48
1:A:61:THR:HG22	1:A:65:PHE:CE2	2.49	0.48
9:I:147:GLU:OE2	16:P:66:GLY:N	2.45	0.48
12:L:26:LEU:HD12	12:L:26:LEU:N	2.28	0.48
13:M:439:LEU:HD21	31:g:64:PHE:CE1	2.49	0.48
14:N:122:ILE:HD11	14:N:127:GLY:HA2	1.96	0.48
37:m:22:TYR:CD1	37:m:22:TYR:N	2.82	0.48
39:o:83:GLN:O	39:o:87:ASP:OD1	2.32	0.48
4:D:371:LYS:HE3	4:D:424:VAL:HG22	1.95	0.48
6:F:65:LEU:HD23	6:F:75:THR:HG21	1.96	0.48
10:J:54:MET:O	10:J:58:ILE:HG12	2.14	0.48
10:J:144:MET:HG2	11:K:58:PRO:HB3	1.95	0.48
11:K:16:LEU:O	11:K:20:LEU:HG	2.14	0.48
12:L:88:LEU:HD23	12:L:326:PHE:CE1	2.48	0.48
13:M:375:LEU:O	13:M:376:MET:C	2.56	0.48
14:N:314:MET:HE3	15:O:267:THR:HG23	1.94	0.48
33:i:122:PHE:HD2	39:o:39:VAL:HG11	1.79	0.48
7:G:57:VAL:HG22	7:G:58:GLU:H	1.78	0.48
8:H:31:MET:SD	9:I:39:THR:HG22	2.54	0.48
1:A:115:GLU:N	1:A:115:GLU:OE2	2.47	0.47
3:C:17:VAL:HG23	3:C:18:THR:N	2.28	0.47
12:L:96:VAL:HG12	12:L:100:ILE:HD12	1.95	0.47
18:S:57:CYS:O	18:S:60:VAL:HG12	2.14	0.47
18:S:90:MET:O	18:S:93:VAL:HG12	2.14	0.47
39:o:43:GLN:NE2	39:o:47:ASP:OD1	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:170:GLU:O	5:E:173:ILE:HG22	2.14	0.47
12:L:5:THR:HG23	12:L:63:ILE:HD13	1.95	0.47
31:g:83:MET:SD	40:p:96:LYS:HD3	2.54	0.47
36:l:32:ASP:N	36:l:32:ASP:OD1	2.47	0.47
4:D:259:MET:HG3	4:D:398:HIS:CE1	2.49	0.47
8:H:17:MET:CE	8:H:225:MET:HG2	2.44	0.47
12:L:10:LEU:HD22	33:i:78:MET:CG	2.44	0.47
22:X:90:TYR:O	25:a:35:GLU:OE2	2.32	0.47
39:o:68:CYS:SG	39:o:79:CYS:SG	3.12	0.47
2:B:180:GLU:O	2:B:181:GLN:CG	2.61	0.47
3:C:199:LEU:HD21	4:D:385:ARG:NH1	2.30	0.47
4:D:342:MET:HE2	4:D:342:MET:HA	1.96	0.47
7:G:230:VAL:HG11	7:G:573:TYR:CD2	2.48	0.47
7:G:384:PRO:HA	7:G:387:GLU:O	2.14	0.47
12:L:152:PHE:CD2	12:L:168:ALA:HB1	2.50	0.47
13:M:36:LEU:HD12	31:g:66:ILE:HG22	1.95	0.47
13:M:141:GLU:H	13:M:141:GLU:CD	2.18	0.47
16:P:188:PHE:O	16:P:190:ALA:N	2.47	0.47
4:D:343:GLU:HG3	7:G:98:LEU:HD13	1.97	0.47
7:G:216:THR:O	7:G:217:ALA:HB2	2.14	0.47
7:G:343:LEU:HD12	7:G:507:TYR:CZ	2.50	0.47
13:M:285:LEU:HD23	13:M:285:LEU:C	2.39	0.47
15:O:70:ASP:OD1	15:O:71:ILE:N	2.46	0.47
18:S:86:VAL:O	18:S:87:THR:C	2.57	0.47
28:d:36:LEU:HD23	28:d:36:LEU:O	2.15	0.47
8:H:179:TRP:CZ3	8:H:183:MET:HE2	2.50	0.47
14:N:40:ILE:CG2	14:N:44:LEU:HD13	2.44	0.47
20:V:47:THR:HG23	20:V:86:GLU:OE2	2.15	0.47
2:B:77:TYR:O	2:B:78:ASP:C	2.57	0.47
6:F:84:LYS:HB3	6:F:85:PRO:HD2	1.96	0.47
8:H:9:LEU:HD13	8:H:95:LEU:HD22	1.97	0.47
12:L:78:MET:HE1	47:i:201:3PE:H362	1.96	0.47
12:L:183:ILE:O	12:L:187:VAL:HG13	2.15	0.47
13:M:127:ILE:HB	13:M:128:PRO:HD3	1.96	0.47
14:N:246:LEU:HD22	14:N:257:LEU:HD21	1.97	0.47
16:P:64:ASP:N	16:P:64:ASP:OD1	2.47	0.47
22:X:52:PRO:HB3	24:Z:79:ASP:OD2	2.14	0.47
23:Y:12:GLU:HG3	23:Y:13:VAL:N	2.30	0.47
29:e:67:LEU:N	29:e:67:LEU:HD23	2.30	0.47
2:B:90:PRO:HB3	2:B:113:VAL:HG13	1.96	0.47
6:F:294:LEU:HD11	6:F:297:VAL:HG13	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:366:THR:O	7:G:367:THR:OG1	2.25	0.47
8:H:118:TRP:CH2	10:J:30:LEU:HD13	2.50	0.47
12:L:131:LEU:HD23	12:L:131:LEU:C	2.40	0.47
12:L:284:THR:O	12:L:288:THR:HG23	2.15	0.47
16:P:220:ASP:OD2	16:P:223:ALA:HB2	2.15	0.47
18:S:32:VAL:HG13	18:S:78:LEU:HD21	1.96	0.47
6:F:435:GLN:O	6:F:436:GLN:C	2.58	0.47
10:J:19:LEU:O	10:J:19:LEU:HD23	2.15	0.47
10:J:101:PHE:O	10:J:105:VAL:HG23	2.15	0.47
4:D:106:LEU:HD11	4:D:391:ILE:HD13	1.97	0.47
10:J:145:TYR:CE1	10:J:152:MET:HE3	2.50	0.47
16:P:333:GLU:HA	16:P:333:GLU:OE1	2.14	0.47
6:F:107:ASP:O	6:F:111:MET:HG2	2.15	0.46
12:L:295:GLN:HE21	12:L:526:LEU:HD21	1.80	0.46
13:M:36:LEU:HD23	13:M:36:LEU:O	2.15	0.46
15:O:305:ALA:O	15:O:309:ASN:ND2	2.48	0.46
7:G:615:THR:O	7:G:619:VAL:HG23	2.14	0.46
12:L:293:LEU:HD11	12:L:418:MET:HG2	1.96	0.46
20:V:34:LEU:HD13	20:V:48:GLU:HG3	1.97	0.46
36:I:97:TRP:CD1	36:I:97:TRP:C	2.93	0.46
3:C:57:VAL:HG23	3:C:58:ILE:N	2.30	0.46
4:D:159:LEU:HD12	4:D:164:MET:HA	1.98	0.46
5:E:95:VAL:HG22	5:E:96:GLY:N	2.30	0.46
6:F:365:CYS:O	6:F:366:ARG:C	2.58	0.46
8:H:179:TRP:HZ3	8:H:183:MET:HE2	1.80	0.46
12:L:486:LEU:O	12:L:489:THR:HG22	2.15	0.46
18:S:22:LEU:HD11	18:S:24:GLN:OE1	2.15	0.46
24:Z:85:ILE:HG21	24:Z:123:MET:HG3	1.98	0.46
4:D:347:HIS:CG	7:G:121:MET:HE1	2.50	0.46
5:E:7:PHE:CE2	7:G:180:ASP:OD1	2.69	0.46
5:E:22:ASP:OD1	5:E:23:PHE:N	2.49	0.46
6:F:20:ARG:NE	6:F:269:GLU:O	2.49	0.46
7:G:331:LEU:C	7:G:331:LEU:HD23	2.40	0.46
9:I:101:ARG:CD	17:R:41:ILE:HD11	2.45	0.46
12:L:187:VAL:HG12	13:M:383:MET:HG3	1.98	0.46
22:X:21:SER:OG	25:a:51:ASP:OD2	2.33	0.46
39:o:45:MET:HG3	39:o:50:LEU:HD22	1.97	0.46
1:A:52:SER:OG	4:D:70:GLY:N	2.48	0.46
1:A:90:MET:HG3	10:J:149:THR:HG22	1.97	0.46
2:B:47:ILE:HG23	8:H:57:MET:HE1	1.97	0.46
4:D:393:ALA:HB1	4:D:427:GLU:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:186:CYS:O	6:F:192:LEU:HD13	2.15	0.46
7:G:380:VAL:HG23	7:G:380:VAL:O	2.15	0.46
7:G:628:PRO:HB2	18:S:57:CYS:SG	2.56	0.46
11:K:1:FME:N	11:K:2:PRO:HD2	2.30	0.46
11:K:70:GLU:OE1	14:N:68:MET:HE2	2.15	0.46
12:L:74:MET:HE1	13:M:307:TRP:HZ2	1.81	0.46
14:N:109:ALA:N	14:N:161:SER:OG	2.46	0.46
15:O:46:LEU:CD1	15:O:48:MET:HE2	2.46	0.46
20:V:66:LYS:O	20:V:70:LEU:HD13	2.15	0.46
42:s:40:TYR:CD1	42:s:44:THR:HG21	2.50	0.46
1:A:87:MET:SD	1:A:88:MET:N	2.88	0.46
7:G:94:MET:HE3	7:G:119:GLN:C	2.40	0.46
8:H:85:LEU:HD23	8:H:105:ILE:HD13	1.97	0.46
10:J:63:MET:HE2	10:J:63:MET:CA	2.46	0.46
12:L:124:PHE:CD2	12:L:124:PHE:C	2.92	0.46
13:M:232:ALA:HA	13:M:236:LEU:HD12	1.98	0.46
16:P:169:SER:OG	16:P:170:ASP:N	2.45	0.46
19:U:46:ASP:O	19:U:50:ILE:HG12	2.16	0.46
24:Z:11:PRO:HD3	24:Z:15:TYR:CZ	2.51	0.46
35:k:29:ILE:HG21	35:k:34:LEU:HD22	1.97	0.46
42:s:34:ASN:OD1	42:s:37:HIS:HD2	1.99	0.46
5:E:126:ILE:HD12	5:E:130:GLU:O	2.15	0.46
6:F:176:PHE:CD1	6:F:176:PHE:N	2.84	0.46
7:G:609:ILE:HG22	7:G:610:THR:N	2.30	0.46
14:N:213:ALA:HB3	14:N:214:PRO:HD3	1.97	0.46
20:V:13:LEU:HD21	20:V:77:GLU:HB2	1.97	0.46
3:C:80:ALA:HB3	3:C:138:PHE:CE1	2.50	0.46
5:E:126:ILE:HG21	5:E:132:THR:HA	1.97	0.46
7:G:106:PRO:O	7:G:218:ARG:NH2	2.48	0.46
24:Z:128:THR:CG2	24:Z:131:GLU:HG3	2.45	0.46
28:d:16:ASP:OD2	28:d:16:ASP:C	2.57	0.46
5:E:113:SER:O	5:E:116:ILE:HG22	2.16	0.46
7:G:35:MET:HE3	7:G:36:GLN:C	2.41	0.46
7:G:296:TRP:CZ2	7:G:562:PRO:HD2	2.51	0.46
17:R:57:ILE:HD12	17:R:58:ALA:N	2.29	0.46
1:A:34:ALA:HB1	8:H:64:LEU:HG	1.97	0.46
3:C:91:GLU:OE2	3:C:108:LYS:NZ	2.41	0.46
5:E:128:VAL:HA	5:E:139:LEU:CD2	2.46	0.46
6:F:365:CYS:SG	6:F:366:ARG:N	2.88	0.46
14:N:149:LEU:HD13	14:N:154:ILE:HD11	1.97	0.46
18:S:15:LEU:HD12	18:S:68:TYR:CE1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:T:66:ASP:O	19:T:70:LEU:HD23	2.16	0.46
29:e:46:ILE:HG22	29:e:50:ARG:NH1	2.31	0.46
40:p:42:TRP:HB3	40:p:43:PRO:HD3	1.98	0.46
1:A:87:MET:SD	1:A:87:MET:C	2.99	0.45
6:F:215:VAL:HG22	6:F:216:PHE:CE2	2.51	0.45
7:G:297:GLU:OE1	21:W:125:TYR:CZ	2.69	0.45
8:H:90:PRO:HB2	8:H:166:ILE:HD11	1.98	0.45
11:K:9:THR:HA	11:K:43:MET:HE1	1.98	0.45
14:N:135:LYS:HE3	14:N:187:MET:HE2	1.97	0.45
14:N:292:PHE:CD1	14:N:292:PHE:C	2.94	0.45
12:L:489:THR:HG23	12:L:490:ALA:N	2.31	0.45
16:P:194:VAL:HG22	16:P:288:HIS:ND1	2.30	0.45
16:P:285:GLU:O	16:P:289:ILE:HG23	2.15	0.45
20:V:13:LEU:O	20:V:13:LEU:HD23	2.16	0.45
24:Z:97:MET:SD	24:Z:100:VAL:HG21	2.56	0.45
32:h:89:LYS:O	32:h:93:ILE:HG12	2.16	0.45
5:E:85:THR:HG21	7:G:175:THR:HB	1.97	0.45
6:F:200:GLN:HE22	7:G:174:THR:HG23	1.81	0.45
8:H:38:ASN:OD1	8:H:38:ASN:C	2.58	0.45
11:K:27:MET:HE1	11:K:75:LEU:HD21	1.97	0.45
12:L:126:ILE:HD12	12:L:126:ILE:H	1.82	0.45
12:L:147:VAL:HG11	12:L:252:MET:HG3	1.98	0.45
13:M:337:ILE:HD12	13:M:339:SER:HB3	1.99	0.45
14:N:108:LEU:HG	14:N:161:SER:OG	2.16	0.45
33:i:112:THR:HG23	33:i:114:GLU:H	1.82	0.45
2:B:107:ALA:HB3	2:B:108:PRO:HD3	1.99	0.45
6:F:212:ASP:O	6:F:213:VAL:HG13	2.16	0.45
7:G:103:LEU:HD13	17:R:70:LYS:O	2.16	0.45
7:G:609:ILE:O	7:G:610:THR:OG1	2.32	0.45
9:I:83:LYS:HE2	9:I:96:ILE:HB	1.97	0.45
10:J:102:LEU:O	10:J:102:LEU:HD23	2.16	0.45
12:L:97:THR:HG21	12:L:125:LEU:HD22	1.98	0.45
27:c:43:LYS:HD3	27:c:48:LEU:HD12	1.98	0.45
32:h:67:TYR:O	32:h:73:ARG:NH1	2.49	0.45
6:F:268:VAL:HG21	6:F:283:HIS:CE1	2.51	0.45
6:F:317:MET:HE1	6:F:334:VAL:HG22	1.99	0.45
7:G:162:PHE:CE2	7:G:166:ILE:HG21	2.52	0.45
8:H:209:SER:O	8:H:213:VAL:HB	2.17	0.45
13:M:297:VAL:HG12	13:M:301:ILE:HD12	1.99	0.45
32:h:108:GLN:OE1	32:h:108:GLN:HA	2.15	0.45
4:D:101:PRO:O	4:D:105:ARG:NH1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:236:SER:HB2	7:G:259:ASN:HD21	1.81	0.45
10:J:23:PRO:HG3	10:J:82:TRP:CE2	2.52	0.45
12:L:121:LEU:HD22	12:L:246:LEU:HD23	1.98	0.45
12:L:301:ILE:CG2	12:L:426:ILE:HD11	2.47	0.45
12:L:574:THR:OG1	14:N:167:TRP:O	2.33	0.45
16:P:26:ALA:O	16:P:28:GLY:N	2.47	0.45
19:U:30:LEU:HD23	19:U:31:SER:N	2.32	0.45
2:B:44:ASP:HB3	2:B:183:LEU:HB2	1.99	0.45
6:F:202:LYS:NZ	7:G:174:THR:HG21	2.31	0.45
6:F:295:LEU:CD2	6:F:343:ILE:HD13	2.46	0.45
6:F:307:ILE:HG13	6:F:327:THR:HG21	1.99	0.45
18:S:22:LEU:HD23	18:S:54:ILE:HD11	1.98	0.45
21:W:100:LYS:HG3	21:W:104:HIS:HB2	1.98	0.45
31:g:33:GLU:N	31:g:33:GLU:OE1	2.50	0.45
33:i:107:ASP:OD1	40:p:17:PRO:HB3	2.17	0.45
37:m:127:SER:O	37:m:128:TYR:C	2.60	0.45
44:H:402:PC1:C24	10:J:30:LEU:HD21	2.47	0.45
9:I:114:ASP:OD1	9:I:117:LYS:N	2.49	0.45
19:U:55:GLU:HG2	19:U:60:PHE:O	2.16	0.45
23:Y:127:LYS:HE3	23:Y:131:LEU:HD11	1.99	0.45
24:Z:55:GLU:OE2	24:Z:55:GLU:HA	2.17	0.45
33:i:31:GLU:O	38:n:113:MET:O	2.35	0.45
4:D:38:ILE:HD12	4:D:38:ILE:H	1.82	0.45
4:D:117:ALA:HB2	4:D:367:ILE:HG12	1.99	0.45
5:E:31:ILE:O	5:E:35:VAL:HG23	2.17	0.45
7:G:216:THR:HG22	9:I:99:GLU:OE2	2.17	0.45
8:H:154:LEU:HD13	8:H:157:ASN:HB3	1.99	0.45
11:K:41:PHE:CD2	11:K:64:LEU:HG	2.52	0.45
15:O:130:SER:O	15:O:133:VAL:HG22	2.16	0.45
17:R:49:VAL:HG11	17:R:81:THR:HG21	1.99	0.45
24:Z:97:MET:HE3	24:Z:103:TRP:CZ2	2.52	0.45
24:Z:142:TYR:CD1	24:Z:142:TYR:C	2.95	0.45
26:b:18:LEU:O	26:b:22:PHE:HD1	2.00	0.45
37:m:7:PRO:HG3	37:m:13:LEU:HD13	1.99	0.45
2:B:119:GLU:O	16:P:54:TYR:OH	2.34	0.45
4:D:67:GLU:HB3	4:D:75:LYS:HB3	1.99	0.45
6:F:215:VAL:HG22	6:F:216:PHE:CD2	2.52	0.45
12:L:295:GLN:O	12:L:301:ILE:HD11	2.17	0.45
13:M:123:GLU:HB2	14:N:255:PRO:HG3	1.99	0.45
16:P:263:TYR:O	16:P:266:ILE:HG22	2.17	0.45
18:S:15:LEU:HD12	18:S:68:TYR:CZ	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:c:11:LYS:HG2	27:c:12:PRO:HD2	1.98	0.45
33:i:111:GLU:OE1	33:i:111:GLU:N	2.50	0.45
38:n:133:ASP:O	38:n:136:VAL:HG12	2.16	0.45
4:D:51:PHE:CZ	4:D:415:VAL:HG22	2.52	0.44
5:E:68:LYS:O	5:E:72:VAL:HG13	2.17	0.44
7:G:9:ILE:HG23	7:G:75:LYS:HB2	2.00	0.44
13:M:57:SER:HB3	13:M:113:THR:HG22	1.99	0.44
23:Y:19:CYS:SG	23:Y:20:HIS:N	2.90	0.44
25:a:22:SER:O	25:a:26:ILE:HG12	2.17	0.44
39:o:65:LEU:HA	39:o:68:CYS:SG	2.56	0.44
41:r:69:MET:SD	41:r:70:SER:N	2.90	0.44
1:A:73:LEU:N	1:A:74:PRO:CD	2.80	0.44
2:B:88:ALA:HB1	8:H:208:VAL:HG13	1.98	0.44
3:C:79:THR:CG2	3:C:80:ALA:N	2.79	0.44
5:E:192:CYS:O	5:E:193:CYS:SG	2.75	0.44
7:G:187:ILE:HD11	7:G:189:LYS:HB2	1.99	0.44
11:K:12:PHE:CE1	14:N:72:LEU:HD22	2.52	0.44
12:L:131:LEU:HD23	12:L:131:LEU:O	2.17	0.44
14:N:170:LEU:HD11	14:N:288:LEU:HD12	2.00	0.44
14:N:218:ALA:HB1	14:N:240:MET:HE1	1.99	0.44
14:N:242:THR:HG23	14:N:243:MET:N	2.32	0.44
21:W:52:MET:CE	21:W:100:LYS:O	2.65	0.44
38:n:19:TYR:CZ	38:n:23:LEU:HD11	2.52	0.44
6:F:125:GLY:O	6:F:129:MET:HG3	2.18	0.44
8:H:291:LYS:O	8:H:292:ASN:ND2	2.51	0.44
8:H:317:TYR:O	8:H:318:MET:HB3	2.18	0.44
12:L:96:VAL:HG12	12:L:100:ILE:CD1	2.48	0.44
13:M:76:MET:HE2	13:M:99:LEU:HD22	1.99	0.44
14:N:200:LEU:CD1	14:N:269:GLU:HG3	2.48	0.44
16:P:159:VAL:O	16:P:159:VAL:HG22	2.18	0.44
53:T:201:EHZ:O2	53:T:201:EHZ:O1	2.31	0.44
23:Y:86:ASP:OD1	23:Y:86:ASP:O	2.34	0.44
31:g:80:ASP:O	31:g:83:MET:N	2.49	0.44
33:i:16:GLU:HA	33:i:16:GLU:OE1	2.16	0.44
34:j:10:ARG:NH2	34:j:16:GLN:OE1	2.49	0.44
7:G:140:LYS:O	7:G:148:THR:OG1	2.35	0.44
10:J:46:PHE:CG	11:K:46:VAL:HG13	2.52	0.44
10:J:99:GLU:OE1	10:J:99:GLU:HA	2.17	0.44
11:K:17:LEU:HD21	11:K:21:MET:SD	2.58	0.44
12:L:468:ILE:O	12:L:472:ILE:HG23	2.17	0.44
18:S:77:SER:OG	18:S:78:LEU:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:W:100:LYS:CG	21:W:104:HIS:HB2	2.47	0.44
47:i:201:3PE:C21	47:i:201:3PE:O31	2.64	0.44
3:C:137:MET:O	3:C:160:HIS:HB3	2.17	0.44
7:G:222:THR:HG21	7:G:241:SER:HB3	1.99	0.44
7:G:575:ASN:ND2	7:G:577:GLU:HG2	2.33	0.44
11:K:54:MET:HA	11:K:57:MET:HE3	2.00	0.44
12:L:332:HIS:HA	12:L:335:PHE:CE2	2.52	0.44
12:L:539:MET:HE2	12:L:540:LYS:N	2.32	0.44
13:M:220:HIS:NE2	13:M:231:LEU:HD22	2.33	0.44
13:M:221:VAL:CG2	13:M:222:GLU:OE1	2.65	0.44
14:N:203:LEU:C	14:N:203:LEU:HD23	2.42	0.44
15:O:46:LEU:HD12	15:O:48:MET:HE2	1.99	0.44
26:b:56:VAL:HG23	26:b:62:MET:HE1	2.00	0.44
30:f:27:ARG:O	30:f:31:GLU:HG2	2.18	0.44
38:n:170:VAL:HG12	38:n:171:THR:N	2.31	0.44
5:E:148:CYS:O	5:E:151:ALA:HB2	2.17	0.44
8:H:85:LEU:HD21	8:H:108:THR:HB	1.98	0.44
12:L:149:ILE:HD11	13:M:364:LEU:CD2	2.47	0.44
23:Y:127:LYS:O	23:Y:131:LEU:CD1	2.66	0.44
37:m:47:LEU:HB3	38:n:165:LEU:HD13	1.98	0.44
6:F:87:ASP:OD1	6:F:87:ASP:O	2.35	0.44
7:G:324:ASP:CB	7:G:571:ALA:HB1	2.45	0.44
7:G:476:LYS:O	7:G:480:THR:HG23	2.18	0.44
9:I:115:MET:HE3	9:I:120:TYR:OH	2.17	0.44
13:M:119:TYR:CZ	13:M:161:LEU:HB2	2.53	0.44
13:M:220:HIS:CD2	13:M:231:LEU:HD22	2.53	0.44
14:N:85:MET:SD	14:N:86:PHE:N	2.89	0.44
15:O:320:LYS:CD	27:c:9:ASN:OD1	2.66	0.44
19:T:25:ILE:HD12	19:T:25:ILE:C	2.42	0.44
39:o:109:GLN:O	39:o:109:GLN:NE2	2.51	0.44
1:A:36:PRO:O	2:B:116:GLN:NE2	2.50	0.44
6:F:309:LYS:HA	6:F:312:CYS:SG	2.58	0.44
6:F:363:THR:O	6:F:364:PRO:C	2.61	0.44
8:H:17:MET:SD	8:H:17:MET:C	3.01	0.44
10:J:32:LEU:HD21	10:J:59:TYR:HD1	1.82	0.44
11:K:39:SER:O	11:K:43:MET:HG3	2.18	0.44
14:N:95:LEU:HA	14:N:98:THR:OG1	2.17	0.44
18:S:88:ARG:C	18:S:92:ASN:OD1	2.60	0.44
42:s:37:HIS:HA	42:s:40:TYR:CD2	2.53	0.44
6:F:202:LYS:HZ2	7:G:174:THR:HG21	1.83	0.44
7:G:291:LEU:HD21	7:G:560:ILE:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:325:ALA:HA	7:G:328:LEU:HB2	2.00	0.44
12:L:147:VAL:HG12	12:L:252:MET:HE3	2.00	0.44
13:M:270:ILE:HD12	13:M:395:LEU:HD22	1.99	0.44
19:U:17:TYR:OH	35:k:22:PRO:O	2.21	0.44
22:X:8:THR:OG1	22:X:11:GLU:HG2	2.18	0.44
4:D:212:TYR:CE1	24:Z:18:ILE:HD11	2.53	0.43
7:G:255:HIS:O	7:G:260:GLU:N	2.42	0.43
8:H:157:ASN:ND2	8:H:159:SER:O	2.51	0.43
12:L:4:PHE:HZ	12:L:87:ILE:HD11	1.83	0.43
12:L:79:SER:HB2	12:L:135:ASN:HB3	2.00	0.43
12:L:233:LEU:HB3	12:L:234:PRO:HD3	2.00	0.43
12:L:562:ILE:CB	12:L:563:PRO:HD3	2.48	0.43
13:M:200:MET:HE1	13:M:243:MET:HG2	2.00	0.43
20:V:63:ASP:OD1	20:V:64:VAL:N	2.51	0.43
28:d:90:MET:SD	32:h:110:VAL:HG11	2.59	0.43
34:j:22:VAL:O	34:j:26:GLU:HG2	2.18	0.43
5:E:136:LEU:O	5:E:137:PHE:HD1	2.01	0.43
8:H:316:PRO:HG3	24:Z:56:ARG:HB3	1.99	0.43
10:J:5:ILE:CG2	11:K:4:THR:CG2	2.96	0.43
12:L:315:VAL:HG12	12:L:319:MET:HE2	2.00	0.43
13:M:36:LEU:HD23	13:M:36:LEU:C	2.43	0.43
5:E:99:HIS:NE2	5:E:140:ILE:HG21	2.33	0.43
7:G:296:TRP:CB	21:W:125:TYR:HE2	2.31	0.43
7:G:443:LEU:HD13	7:G:477:ILE:HD12	2.00	0.43
13:M:59:ASP:N	13:M:62:SER:OG	2.48	0.43
3:C:96:LEU:HB2	3:C:105:ILE:HG22	2.00	0.43
9:I:158:ASN:OD1	9:I:158:ASN:N	2.52	0.43
12:L:72:ASN:O	12:L:73:SER:CB	2.65	0.43
13:M:115:LEU:HB2	13:M:174:LEU:HD22	1.99	0.43
13:M:235:LEU:HA	13:M:238:LEU:HG	2.00	0.43
18:S:20:VAL:HG22	18:S:64:LEU:HD23	2.00	0.43
40:p:132:THR:O	40:p:135:THR:HG22	2.19	0.43
3:C:51:CYS:HB3	20:V:111:TRP:CE2	2.53	0.43
4:D:62:LEU:HB2	4:D:425:PHE:CZ	2.54	0.43
6:F:192:LEU:O	6:F:193:ILE:C	2.59	0.43
7:G:263:ILE:HD11	7:G:268:ARG:CB	2.49	0.43
12:L:221:THR:HG23	12:L:226:GLN:HB2	2.00	0.43
13:M:151:PHE:O	13:M:155:ILE:HG12	2.19	0.43
13:M:300:SER:HB3	13:M:381:ILE:HG23	2.00	0.43
13:M:403:THR:HA	13:M:406:TYR:CE2	2.53	0.43
23:Y:41:SER:O	23:Y:44:LEU:O	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:c:6:GLU:OE2	27:c:6:GLU:HA	2.17	0.43
31:g:81:TYR:CE1	31:g:82:ARG:HG3	2.53	0.43
33:i:24:ASP:OD1	38:n:120:LYS:HD3	2.18	0.43
33:i:93:PRO:HB3	40:p:4:TRP:CD1	2.53	0.43
39:o:111:LYS:O	39:o:113:ARG:N	2.52	0.43
6:F:147:SER:O	6:F:151:VAL:HG13	2.19	0.43
7:G:429:LEU:HD21	7:G:470:VAL:CG2	2.49	0.43
15:O:189:PRO:HA	15:O:192:MET:HE3	2.00	0.43
17:R:47:GLN:HB2	17:R:48:PRO:HD2	2.01	0.43
28:d:81:TYR:O	28:d:85:VAL:HG23	2.18	0.43
38:n:55:ASP:OD1	38:n:58:ARG:HB3	2.17	0.43
4:D:116:GLN:NE2	4:D:276:ASP:OD2	2.52	0.43
6:F:68:ARG:HD2	6:F:254:LYS:HD2	2.01	0.43
6:F:184:TYR:CE1	46:F:502:FMN:H6	2.52	0.43
10:J:55:VAL:CG2	11:K:41:PHE:CE1	3.02	0.43
10:J:63:MET:HE2	10:J:63:MET:HA	1.98	0.43
13:M:91:LEU:HD21	13:M:135:ARG:NH2	2.34	0.43
13:M:118:PHE:O	13:M:122:PHE:HB3	2.19	0.43
15:O:128:ILE:HD12	15:O:160:LEU:HD12	2.01	0.43
18:S:32:VAL:O	18:S:36:ILE:HG12	2.18	0.43
5:E:105:THR:HB	45:E:301:FES:S2	2.58	0.43
6:F:406:ALA:HB3	46:F:502:FMN:HM81	2.01	0.43
7:G:36:GLN:O	7:G:36:GLN:HG2	2.18	0.43
7:G:126:ASP:OD1	41:r:56:ARG:NH2	2.37	0.43
7:G:685:ALA:O	7:G:689:LYS:HG3	2.19	0.43
8:H:151:LEU:HD23	8:H:151:LEU:C	2.43	0.43
8:H:153:VAL:HG13	8:H:154:LEU:HD22	2.00	0.43
9:I:177:TYR:CZ	41:r:38:PRO:HG3	2.54	0.43
12:L:407:TRP:CZ3	12:L:410:LEU:HD23	2.54	0.43
13:M:125:THR:O	13:M:129:THR:HG23	2.19	0.43
14:N:95:LEU:HD12	14:N:95:LEU:H	1.82	0.43
14:N:300:THR:OG1	14:N:301:SER:N	2.51	0.43
15:O:308:TYR:OH	27:c:9:ASN:ND2	2.51	0.43
16:P:113:ILE:HB	16:P:114:PRO:HD3	2.01	0.43
16:P:166:ILE:HD12	16:P:230:PHE:CE2	2.54	0.43
19:T:64:ASP:OD1	21:W:35:ARG:NH1	2.52	0.43
22:X:81:PHE:HD2	25:a:66:LEU:HD11	1.83	0.43
30:f:42:LEU:HD13	40:p:93:ARG:NH2	2.33	0.43
35:k:20:GLU:HA	35:k:20:GLU:OE2	2.19	0.43
42:s:35:LEU:N	42:s:35:LEU:HD22	2.33	0.43
1:A:51:PHE:CD2	8:H:133:LEU:HD22	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:16:ASN:OD1	5:E:18:ASP:N	2.47	0.43
6:F:126:GLY:CA	6:F:131:ALA:HB3	2.48	0.43
6:F:256:PHE:N	6:F:256:PHE:CD1	2.86	0.43
7:G:292:THR:OG1	7:G:293:TYR:N	2.51	0.43
7:G:402:ASN:O	7:G:403:ASP:C	2.62	0.43
10:J:127:GLU:HG2	10:J:128:VAL:N	2.34	0.43
12:L:63:ILE:HG22	33:i:89:MET:HE1	2.00	0.43
12:L:313:MET:HE1	12:L:332:HIS:CE1	2.53	0.43
13:M:229:MET:SD	13:M:328:CYS:SG	3.17	0.43
20:V:8:THR:OG1	20:V:13:LEU:O	2.37	0.43
22:X:4:VAL:HG22	22:X:5:GLU:N	2.33	0.43
29:e:40:ILE:HD11	32:h:128:PRO:CB	2.48	0.43
1:A:94:LEU:HD12	10:J:152:MET:SD	2.59	0.43
4:D:405:MET:HE1	4:D:421:GLN:NE2	2.34	0.43
5:E:53:LEU:HD12	5:E:53:LEU:O	2.19	0.43
7:G:137:VAL:HG12	7:G:138:GLU:N	2.33	0.43
7:G:208:LEU:HD13	43:G:802:SF4:S1	2.59	0.43
7:G:296:TRP:CH2	7:G:561:LEU:HD22	2.54	0.43
7:G:611:LEU:HD22	7:G:611:LEU:N	2.33	0.43
8:H:93:HIS:O	25:a:23:THR:HG21	2.19	0.43
10:J:6:PHE:CB	10:J:120:LEU:HD21	2.49	0.43
12:L:127:THR:HG21	12:L:146:GLY:C	2.44	0.43
12:L:149:ILE:CD1	13:M:364:LEU:CD2	2.97	0.43
14:N:21:MET:CE	29:e:3:LEU:HD22	2.49	0.43
15:O:211:LEU:HB2	15:O:212:PRO:HD3	2.01	0.43
31:g:60:PHE:HB3	32:h:28:TYR:CD2	2.54	0.43
33:i:110:LEU:HD21	40:p:18:ALA:HA	2.01	0.43
37:m:25:SER:O	37:m:28:THR:OG1	2.36	0.43
1:A:57:LEU:HD13	10:J:170:THR:OG1	2.19	0.42
4:D:143:GLU:OE1	4:D:143:GLU:HA	2.19	0.42
7:G:431:ASP:OD1	7:G:436:ARG:NH2	2.52	0.42
9:I:82:CYS:HB2	9:I:84:LEU:HD13	2.00	0.42
11:K:26:LEU:N	11:K:88:ASP:OD2	2.52	0.42
13:M:166:LEU:HD23	13:M:166:LEU:C	2.44	0.42
14:N:289:ASN:HA	14:N:292:PHE:CE2	2.53	0.42
18:S:64:LEU:HD11	18:S:78:LEU:HD12	2.01	0.42
20:V:4:LEU:HD23	20:V:5:LYS:N	2.34	0.42
30:f:42:LEU:O	30:f:42:LEU:HD23	2.19	0.42
32:h:45:ILE:HG23	32:h:46:PHE:CD2	2.54	0.42
39:o:17:ASP:C	39:o:17:ASP:OD1	2.61	0.42
4:D:151:ILE:HG21	4:D:174:ARG:HG3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:59:ILE:O	7:G:62:ALA:HB3	2.18	0.42
7:G:357:ASP:HA	18:S:53:LEU:HD13	2.01	0.42
15:O:141:GLN:HE22	15:O:197:ALA:HB3	1.84	0.42
16:P:185:TYR:HD2	16:P:191:VAL:HG13	1.83	0.42
19:U:19:LEU:HD13	19:U:73:PRO:HG3	2.01	0.42
6:F:60:MET:HE2	6:F:60:MET:CA	2.48	0.42
7:G:13:VAL:HG11	7:G:33:VAL:HG11	2.01	0.42
7:G:213:TYR:CG	7:G:213:TYR:O	2.72	0.42
8:H:175:LEU:O	8:H:179:TRP:N	2.52	0.42
13:M:57:SER:OG	13:M:58:SER:N	2.51	0.42
14:N:73:ALA:CB	14:N:98:THR:HG23	2.50	0.42
19:T:52:MET:HE1	21:W:42:TYR:CG	2.54	0.42
24:Z:93:GLU:OE2	29:e:74:ARG:NH1	2.52	0.42
30:f:50:ASN:OD1	30:f:50:ASN:C	2.61	0.42
31:g:99:GLU:OE2	40:p:108:ARG:NH1	2.51	0.42
1:A:3:LEU:HD23	47:H:401:3PE:H221	2.01	0.42
12:L:204:SER:OG	12:L:207:ASN:ND2	2.52	0.42
12:L:542:LEU:HB3	13:M:278:ARG:HD2	2.01	0.42
14:N:126:MET:HE2	14:N:126:MET:CA	2.50	0.42
23:Y:48:ASP:OD1	23:Y:48:ASP:N	2.52	0.42
25:a:6:LEU:N	25:a:7:PRO:CD	2.83	0.42
6:F:325:ALA:O	6:F:326:GLN:HB2	2.18	0.42
7:G:511:VAL:HG22	7:G:514:ILE:HB	2.01	0.42
9:I:129:PRO:HD2	43:I:202:SF4:S3	2.59	0.42
13:M:204:LEU:HD22	13:M:209:LEU:HD13	2.00	0.42
15:O:147:GLN:N	15:O:147:GLN:OE1	2.52	0.42
16:P:136:ASN:ND2	16:P:291:ASP:OD1	2.47	0.42
23:Y:10:TYR:CD1	23:Y:10:TYR:C	2.97	0.42
26:b:33:MET:C	26:b:34:ILE:HD13	2.45	0.42
37:m:74:ASN:ND2	37:m:78:ASN:OD1	2.52	0.42
3:C:79:THR:OG1	4:D:390:LYS:HG2	2.19	0.42
3:C:151:ILE:HG23	3:C:152:LEU:HG	2.01	0.42
5:E:34:ILE:HD11	42:s:46:LEU:HD22	2.02	0.42
8:H:290:TRP:O	8:H:295:PRO:HD3	2.19	0.42
12:L:10:LEU:HD22	33:i:78:MET:HG2	2.01	0.42
12:L:604:ILE:HG13	12:L:605:ASN:N	2.34	0.42
21:W:121:LEU:O	21:W:125:TYR:CD1	2.72	0.42
38:n:158:ARG:NH2	38:n:161:ASP:OD2	2.52	0.42
6:F:151:VAL:HG12	42:s:48:LEU:CD2	2.47	0.42
7:G:359:ARG:HB3	7:G:363:LEU:HD12	2.01	0.42
12:L:122:THR:O	12:L:126:ILE:CD1	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:147:VAL:CG1	12:L:252:MET:HG3	2.50	0.42
13:M:13:PRO:HB2	30:f:14:LEU:HD13	2.02	0.42
13:M:240:SER:OG	13:M:316:MET:SD	2.50	0.42
14:N:314:MET:HE1	15:O:267:THR:HG23	2.00	0.42
16:P:280:THR:HG22	16:P:281:LYS:N	2.35	0.42
28:d:17:GLU:OE1	28:d:17:GLU:N	2.34	0.42
1:A:77:TRP:CE2	10:J:136:GLU:OE2	2.73	0.42
1:A:114:THR:O	1:A:114:THR:HG23	2.19	0.42
3:C:109:THR:OG1	3:C:110:TYR:N	2.53	0.42
4:D:45:THR:O	4:D:45:THR:HG22	2.20	0.42
8:H:110:SER:O	8:H:113:VAL:HG12	2.20	0.42
12:L:3:ILE:HG23	12:L:4:PHE:N	2.34	0.42
16:P:179:LEU:CD2	16:P:245:ILE:HD11	2.49	0.42
22:X:92:ASN:OD1	22:X:92:ASN:C	2.62	0.42
32:h:113:LEU:HD22	32:h:117:ARG:NH1	2.35	0.42
1:A:79:ILE:CD1	8:H:155:LEU:CD2	2.98	0.42
2:B:35:ARG:O	2:B:38:TYR:HB3	2.19	0.42
3:C:155:TYR:HD1	21:W:94:GLU:OE1	2.02	0.42
4:D:19:MET:HE2	14:N:295:ARG:NE	2.35	0.42
4:D:185:SER:CB	4:D:191:ALA:HB1	2.50	0.42
4:D:259:MET:CE	4:D:296:ARG:CD	2.98	0.42
5:E:43:GLN:CD	5:E:43:GLN:H	2.27	0.42
6:F:30:ASP:OD1	6:F:31:TRP:N	2.53	0.42
6:F:137:TYR:C	6:F:138:ILE:HD12	2.44	0.42
7:G:429:LEU:HD21	7:G:470:VAL:HG23	2.00	0.42
12:L:594:ASN:OD1	14:N:110:PRO:HB2	2.19	0.42
13:M:63:THR:OG1	13:M:64:PRO:HD3	2.20	0.42
16:P:26:ALA:O	16:P:31:GLY:HA3	2.20	0.42
21:W:98:VAL:HG12	21:W:98:VAL:O	2.19	0.42
36:l:76:ILE:HG23	36:l:80:LEU:HD22	2.00	0.42
40:p:70:VAL:HB	40:p:71:PRO:HD2	2.02	0.42
3:C:78:LEU:HD23	3:C:78:LEU:O	2.20	0.42
5:E:31:ILE:HD13	5:E:53:LEU:HD23	1.98	0.42
5:E:46:ALA:O	5:E:50:VAL:HG23	2.20	0.42
7:G:565:ALA:N	7:G:568:GLU:OE1	2.53	0.42
8:H:178:ALA:HB1	8:H:181:MET:HE3	2.02	0.42
16:P:139:MET:SD	16:P:139:MET:N	2.93	0.42
18:S:22:LEU:O	18:S:56:GLU:HA	2.19	0.42
19:T:45:LEU:HD12	21:W:38:TYR:CD2	2.55	0.42
21:W:20:LYS:HD3	21:W:20:LYS:N	2.35	0.42
24:Z:97:MET:HG2	24:Z:103:TRP:CG	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:l:3:MET:HG3	36:l:7:MET:HE3	2.01	0.42
42:s:36:GLN:OE1	42:s:36:GLN:N	2.31	0.42
6:F:295:LEU:HD23	6:F:296:ALA:N	2.35	0.41
7:G:138:GLU:OE2	17:R:77:LYS:N	2.44	0.41
7:G:452:VAL:HA	7:G:491:ASN:OD1	2.20	0.41
7:G:465:ALA:HB1	7:G:656:VAL:HG12	2.01	0.41
12:L:173:LEU:HD13	13:M:405:MET:SD	2.59	0.41
12:L:237:MET:CE	12:L:303:ALA:HB2	2.50	0.41
12:L:433:THR:CG2	12:L:434:LYS:N	2.83	0.41
16:P:23:VAL:O	16:P:48:PRO:HD2	2.20	0.41
18:S:62:PRO:HG2	18:S:78:LEU:O	2.20	0.41
19:U:55:GLU:HG3	19:U:62:ILE:HG12	2.02	0.41
6:F:403:THR:HB	43:F:501:SF4:S4	2.60	0.41
7:G:306:MET:HB3	7:G:521:MET:SD	2.60	0.41
8:H:32:GLN:O	8:H:33:LEU:CB	2.68	0.41
11:K:20:LEU:HD12	12:L:592:LEU:HD12	2.00	0.41
12:L:119:LYS:NZ	48:L:702:CDL:OB4	2.44	0.41
12:L:595:ILE:CD1	14:N:100:MET:CE	2.98	0.41
13:M:33:LEU:HD21	31:g:62:PHE:CD1	2.55	0.41
14:N:95:LEU:CD2	14:N:148:LEU:HB3	2.49	0.41
14:N:255:PRO:HA	14:N:260:PHE:CG	2.55	0.41
15:O:188:ASP:OD2	15:O:191:GLU:N	2.47	0.41
16:P:193:LEU:O	16:P:257:PRO:HA	2.20	0.41
24:Z:29:LEU:HB3	24:Z:34:MET:HE2	2.02	0.41
34:j:26:GLU:HA	34:j:26:GLU:OE1	2.20	0.41
2:B:68:GLU:HG3	2:B:160:PRO:HB2	2.01	0.41
4:D:371:LYS:HE3	4:D:424:VAL:HG13	2.02	0.41
7:G:53:ARG:N	45:G:803:FES:S1	2.84	0.41
7:G:222:THR:CG2	7:G:241:SER:HB3	2.49	0.41
7:G:491:ASN:OD1	7:G:491:ASN:N	2.52	0.41
12:L:53:MET:HA	40:p:29:ILE:HD11	2.03	0.41
12:L:65:THR:CA	33:i:89:MET:HE2	2.42	0.41
12:L:127:THR:HG22	12:L:147:VAL:HG23	2.03	0.41
15:O:230:ASP:OD1	15:O:232:THR:HG22	2.20	0.41
17:R:80:LYS:HD2	17:R:81:THR:H	1.85	0.41
19:U:86:VAL:O	19:U:88:GLU:N	2.53	0.41
33:i:84:TYR:O	33:i:85:VAL:C	2.64	0.41
38:n:98:VAL:HG23	38:n:99:PRO:HD2	2.02	0.41
4:D:5:GLN:O	4:D:6:PRO:C	2.62	0.41
8:H:307:LEU:HA	8:H:310:PHE:CE2	2.56	0.41
10:J:55:VAL:HG21	11:K:41:PHE:CE1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:96:LEU:HB2	14:N:54:GLU:OE2	2.20	0.41
12:L:150:MET:HE3	12:L:150:MET:HA	2.03	0.41
12:L:216:LEU:HD22	12:L:259:LEU:HD21	2.02	0.41
47:L:701:3PE:O22	47:L:701:3PE:C1	2.68	0.41
32:h:110:VAL:HG22	32:h:114:MET:HE3	2.02	0.41
38:n:138:GLN:NE2	38:n:155:PRO:O	2.53	0.41
2:B:124:VAL:HG22	2:B:154:ILE:HB	2.02	0.41
6:F:93:LEU:CD2	6:F:95:VAL:HG23	2.51	0.41
14:N:73:ALA:HB1	14:N:98:THR:HG23	2.03	0.41
21:W:82:ASP:O	21:W:86:ILE:HG12	2.19	0.41
22:X:129:LYS:CG	26:b:65:VAL:HG13	2.51	0.41
2:B:42:LYS:HA	2:B:42:LYS:HD3	1.92	0.41
7:G:518:PRO:HA	7:G:519:PRO:HD3	1.92	0.41
8:H:11:VAL:HB	8:H:12:PRO:HD3	2.03	0.41
10:J:145:TYR:CD1	10:J:152:MET:HE3	2.55	0.41
11:K:20:LEU:O	12:L:588:PHE:HB3	2.19	0.41
12:L:326:PHE:CD1	12:L:326:PHE:C	2.98	0.41
13:M:157:SER:O	13:M:160:LEU:HB3	2.20	0.41
13:M:271:MET:O	13:M:274:SER:OG	2.33	0.41
31:g:80:ASP:CG	31:g:83:MET:HA	2.46	0.41
31:g:103:LEU:HD12	31:g:103:LEU:N	2.36	0.41
32:h:34:ILE:HA	32:h:37:ILE:HG22	2.01	0.41
40:p:128:LEU:HD23	40:p:128:LEU:C	2.45	0.41
7:G:156:CYS:N	43:G:802:SF4:S4	2.92	0.41
7:G:273:GLY:O	7:G:549:HIS:NE2	2.44	0.41
7:G:433:ALA:HB2	7:G:473:MET:HA	2.02	0.41
7:G:433:ALA:HB2	7:G:473:MET:HG3	2.03	0.41
8:H:20:LEU:HB2	25:a:12:MET:HE2	2.03	0.41
12:L:253:VAL:HB	12:L:310:LEU:HD21	2.02	0.41
12:L:332:HIS:HA	12:L:335:PHE:CZ	2.55	0.41
12:L:595:ILE:HD13	14:N:100:MET:CE	2.51	0.41
14:N:40:ILE:HG23	14:N:44:LEU:HD13	2.02	0.41
19:T:70:LEU:HD12	19:T:76:ILE:HA	2.02	0.41
29:e:44:HIS:CG	32:h:130:LYS:HE2	2.56	0.41
2:B:126:MET:SD	2:B:166:LEU:HD13	2.60	0.41
8:H:81:LEU:HD22	8:H:108:THR:HG23	2.03	0.41
12:L:310:LEU:HD12	12:L:313:MET:HE3	2.02	0.41
12:L:491:LEU:CD1	34:j:47:VAL:HG13	2.50	0.41
13:M:220:HIS:ND1	13:M:327:PHE:CZ	2.89	0.41
14:N:203:LEU:HD13	14:N:344:ILE:HG22	2.03	0.41
18:S:22:LEU:HD12	18:S:22:LEU:C	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:Z:104:LYS:HB2	24:Z:107:GLU:HB2	2.01	0.41
26:b:30:ILE:HG23	26:b:31:MET:N	2.36	0.41
39:o:51:THR:HA	40:p:118:GLU:HG2	2.03	0.41
1:A:72:LEU:HD11	1:A:94:LEU:HD22	2.03	0.41
4:D:182:GLU:OE1	9:I:60:SER:OG	2.33	0.41
4:D:424:VAL:HG23	4:D:424:VAL:O	2.21	0.41
5:E:132:THR:HG22	5:E:137:PHE:H	1.86	0.41
6:F:97:ALA:HB2	6:F:225:VAL:HG21	2.03	0.41
6:F:409:ASP:OD1	6:F:409:ASP:N	2.54	0.41
7:G:418:ARG:NE	7:G:418:ARG:HA	2.35	0.41
7:G:682:GLN:O	7:G:685:ALA:HB3	2.20	0.41
8:H:228:TYR:HA	8:H:231:ILE:HD12	2.03	0.41
8:H:273:ILE:HG23	8:H:277:TYR:CD2	2.55	0.41
9:I:66:GLU:OE2	9:I:151:TYR:OH	2.30	0.41
9:I:115:MET:HE3	9:I:115:MET:HB3	1.98	0.41
12:L:264:HIS:N	12:L:265:PRO:HD2	2.35	0.41
12:L:391:SER:OG	12:L:392:LYS:N	2.54	0.41
12:L:482:MET:SD	12:L:487:LYS:HA	2.61	0.41
12:L:489:THR:HA	12:L:492:ILE:HD11	2.03	0.41
12:L:513:MET:HE1	38:n:30:CYS:SG	2.60	0.41
13:M:390:ASN:OD1	13:M:390:ASN:N	2.54	0.41
48:M:502:CDL:H522	28:d:53:VAL:HG11	2.03	0.41
15:O:36:ASN:OD1	15:O:37:LYS:N	2.54	0.41
15:O:171:TYR:CE1	15:O:206:TYR:CD2	3.09	0.41
17:R:43:LEU:H	17:R:43:LEU:HD12	1.86	0.41
21:W:51:LEU:O	21:W:51:LEU:HD23	2.21	0.41
22:X:63:ASN:OD1	24:Z:80:ARG:NH2	2.54	0.41
31:g:119:PRO:O	40:p:166:ARG:NH1	2.47	0.41
32:h:81:ASP:OD1	32:h:81:ASP:N	2.53	0.41
1:A:53:MET:HE1	10:J:69:TYR:CD1	2.56	0.41
5:E:117:LEU:O	5:E:117:LEU:HD12	2.21	0.41
5:E:142:VAL:HG12	5:E:143:GLU:N	2.35	0.41
6:F:268:VAL:HG21	6:F:283:HIS:NE2	2.36	0.41
12:L:156:GLY:CA	12:L:164:ALA:HB1	2.50	0.41
12:L:419:THR:HA	12:L:422:TYR:CE2	2.56	0.41
13:M:282:LEU:HD21	13:M:359:TRP:HH2	1.85	0.41
17:R:45:ALA:C	17:R:47:GLN:OE1	2.64	0.41
22:X:52:PRO:HD2	24:Z:115:TRP:CG	2.56	0.41
27:c:30:TYR:CD1	27:c:30:TYR:C	2.99	0.41
35:k:83:PHE:CZ	35:k:87:LEU:HD22	2.56	0.41
1:A:78:ALA:HB2	10:J:145:TYR:HD2	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:165:ILE:HG23	14:N:42:PRO:HG3	2.04	0.40
13:M:264:LEU:HD23	13:M:264:LEU:HA	1.96	0.40
23:Y:9:SER:O	23:Y:12:GLU:HG2	2.22	0.40
23:Y:63:THR:O	23:Y:67:ILE:HG12	2.21	0.40
30:f:42:LEU:HD23	30:f:42:LEU:C	2.46	0.40
35:k:37:VAL:HG13	38:n:38:TYR:HB2	2.02	0.40
1:A:70:ALA:HB2	10:J:58:ILE:HD11	2.02	0.40
4:D:424:VAL:O	4:D:427:GLU:HG2	2.21	0.40
6:F:50:LEU:C	6:F:50:LEU:HD12	2.46	0.40
6:F:322:LEU:HD22	6:F:327:THR:OG1	2.21	0.40
6:F:428:ASP:OD1	6:F:428:ASP:C	2.65	0.40
8:H:40:VAL:HG21	8:H:47:GLN:OE1	2.21	0.40
12:L:78:MET:HE1	47:i:201:3PE:C36	2.51	0.40
12:L:124:PHE:CD2	12:L:247:LEU:HD11	2.56	0.40
48:M:502:CDL:H521	28:d:53:VAL:HG11	2.02	0.40
21:W:40:ALA:HB1	21:W:92:LEU:HD21	2.03	0.40
22:X:63:ASN:OD1	24:Z:80:ARG:NH1	2.54	0.40
24:Z:143:THR:HB	25:a:48:MET:HE1	2.03	0.40
4:D:165:THR:N	4:D:166:PRO:HD2	2.36	0.40
4:D:184:VAL:O	9:I:62:ARG:NH1	2.54	0.40
5:E:41:GLY:O	5:E:42:HIS:C	2.65	0.40
6:F:32:ARG:NH1	42:s:36:GLN:HE21	2.20	0.40
7:G:240:VAL:O	7:G:240:VAL:HG13	2.22	0.40
8:H:260:MET:HG3	25:a:17:VAL:HG23	2.02	0.40
9:I:5:VAL:HG23	41:r:112:LEU:HD22	2.03	0.40
11:K:15:SER:CB	11:K:36:MET:HG2	2.51	0.40
12:L:203:PHE:CE1	40:p:109:LEU:HD11	2.55	0.40
13:M:41:LEU:HB3	13:M:63:THR:HG22	2.03	0.40
13:M:81:GLN:NE2	31:g:45:TYR:O	2.53	0.40
29:e:49:THR:O	29:e:53:LYS:NZ	2.51	0.40
36:l:111:MET:HA	36:l:111:MET:HE3	2.02	0.40
2:B:44:ASP:OD1	2:B:44:ASP:N	2.48	0.40
6:F:87:ASP:OD1	6:F:87:ASP:C	2.64	0.40
7:G:451:VAL:HG11	7:G:473:MET:HE1	2.04	0.40
8:H:286:MET:HE2	8:H:286:MET:HB2	1.92	0.40
12:L:71:MET:O	12:L:72:ASN:C	2.63	0.40
12:L:208:ASP:O	12:L:270:ASN:ND2	2.54	0.40
12:L:316:THR:HA	12:L:319:MET:HE3	2.03	0.40
12:L:595:ILE:HD12	14:N:100:MET:HE1	2.03	0.40
13:M:134:THR:HG1	14:N:302:LEU:HD13	1.86	0.40
13:M:307:TRP:CD1	13:M:459:MET:HE1	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:41:GLU:HG3	15:O:231:PRO:HG2	2.04	0.40
15:O:52:PRO:O	15:O:107:GLN:NE2	2.47	0.40
15:O:135:LEU:HD22	15:O:152:TYR:CD1	2.57	0.40
24:Z:59:LEU:O	24:Z:63:ASP:OD2	2.39	0.40
39:o:16:PRO:HG3	39:o:104:GLU:OE1	2.21	0.40
39:o:62:LEU:O	39:o:62:LEU:HD12	2.21	0.40
3:C:155:TYR:HB3	21:W:94:GLU:OE1	2.21	0.40
4:D:44:VAL:O	4:D:44:VAL:CG1	2.69	0.40
4:D:338:MET:HA	4:D:344:SER:OG	2.20	0.40
6:F:12:PHE:HB3	6:F:271:GLU:OE1	2.22	0.40
13:M:126:LEU:HD21	13:M:153:THR:HG21	2.04	0.40
13:M:155:ILE:HD13	14:N:287:LEU:CD1	2.52	0.40
14:N:142:LEU:HB3	14:N:194:LEU:HD21	2.03	0.40
15:O:13:LYS:O	15:O:16:LYS:HB3	2.22	0.40
31:g:42:PHE:CE2	31:g:44:GLY:HA2	2.56	0.40
34:j:41:TRP:CD2	35:k:87:LEU:HD21	2.57	0.40
36:l:98:ASP:OD1	36:l:99:VAL:HG23	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	113/115 (98%)	108 (96%)	5 (4%)	0	100	100
2	B	154/224 (69%)	140 (91%)	14 (9%)	0	100	100
3	C	205/263 (78%)	195 (95%)	10 (5%)	0	100	100
4	D	427/463 (92%)	402 (94%)	25 (6%)	0	100	100
5	E	208/245 (85%)	192 (92%)	15 (7%)	1 (0%)	24	59
6	F	426/464 (92%)	387 (91%)	37 (9%)	2 (0%)	24	59
7	G	686/727 (94%)	621 (90%)	63 (9%)	2 (0%)	36	69

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
8	H	316/318 (99%)	297 (94%)	19 (6%)	0	100	100
9	I	176/212 (83%)	162 (92%)	14 (8%)	0	100	100
10	J	169/171 (99%)	159 (94%)	10 (6%)	0	100	100
11	K	96/98 (98%)	91 (95%)	5 (5%)	0	100	100
12	L	604/606 (100%)	561 (93%)	42 (7%)	1 (0%)	43	74
13	M	457/459 (100%)	439 (96%)	18 (4%)	0	100	100
14	N	342/345 (99%)	331 (97%)	11 (3%)	0	100	100
15	O	318/355 (90%)	307 (96%)	11 (4%)	0	100	100
16	P	323/377 (86%)	301 (93%)	21 (6%)	1 (0%)	36	69
17	R	53/116 (46%)	52 (98%)	1 (2%)	0	100	100
18	S	80/99 (81%)	67 (84%)	12 (15%)	1 (1%)	9	40
19	T	74/156 (47%)	68 (92%)	6 (8%)	0	100	100
19	U	84/156 (54%)	79 (94%)	5 (6%)	0	100	100
20	V	110/116 (95%)	103 (94%)	7 (6%)	0	100	100
21	W	107/131 (82%)	103 (96%)	3 (3%)	1 (1%)	14	47
22	X	169/172 (98%)	157 (93%)	12 (7%)	0	100	100
23	Y	138/143 (96%)	132 (96%)	6 (4%)	0	100	100
24	Z	139/144 (96%)	133 (96%)	6 (4%)	0	100	100
25	a	66/70 (94%)	64 (97%)	2 (3%)	0	100	100
26	b	78/84 (93%)	72 (92%)	6 (8%)	0	100	100
27	c	46/76 (60%)	45 (98%)	1 (2%)	0	100	100
28	d	118/120 (98%)	113 (96%)	5 (4%)	0	100	100
29	e	103/106 (97%)	96 (93%)	7 (7%)	0	100	100
30	f	51/57 (90%)	43 (84%)	7 (14%)	1 (2%)	6	32
31	g	99/151 (66%)	94 (95%)	5 (5%)	0	100	100
32	h	136/189 (72%)	132 (97%)	4 (3%)	0	100	100
33	i	90/127 (71%)	83 (92%)	7 (8%)	0	100	100
34	j	60/105 (57%)	56 (93%)	4 (7%)	0	100	100
35	k	73/104 (70%)	72 (99%)	1 (1%)	0	100	100
36	l	152/186 (82%)	144 (95%)	8 (5%)	0	100	100
37	m	124/129 (96%)	118 (95%)	6 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
38	n	175/179 (98%)	171 (98%)	4 (2%)	0	100	100
39	o	111/137 (81%)	102 (92%)	9 (8%)	0	100	100
40	p	166/176 (94%)	157 (95%)	9 (5%)	0	100	100
41	r	57/113 (50%)	53 (93%)	4 (7%)	0	100	100
42	s	29/104 (28%)	29 (100%)	0	0	100	100
All	All	7708/8888 (87%)	7231 (94%)	467 (6%)	10 (0%)	49	80

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	G	339	ASP
18	S	88	ARG
30	f	43	PHE
5	E	192	CYS
6	F	287	VAL
12	L	249	SER
7	G	187	ILE
21	W	98	VAL
6	F	115	PRO
16	P	292	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	103/103 (100%)	103 (100%)	0	100	100
2	B	132/185 (71%)	132 (100%)	0	100	100
3	C	189/227 (83%)	189 (100%)	0	100	100
4	D	370/394 (94%)	369 (100%)	1 (0%)	86	85
5	E	183/205 (89%)	183 (100%)	0	100	100
6	F	343/370 (93%)	342 (100%)	1 (0%)	86	85
7	G	567/610 (93%)	565 (100%)	2 (0%)	84	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	H	279/279 (100%)	278 (100%)	1 (0%)	84	83
9	I	152/178 (85%)	152 (100%)	0	100	100
10	J	136/136 (100%)	136 (100%)	0	100	100
11	K	87/87 (100%)	87 (100%)	0	100	100
12	L	548/548 (100%)	545 (100%)	3 (0%)	81	82
13	M	414/414 (100%)	414 (100%)	0	100	100
14	N	306/307 (100%)	306 (100%)	0	100	100
15	O	284/309 (92%)	284 (100%)	0	100	100
16	P	286/325 (88%)	285 (100%)	1 (0%)	86	85
17	R	43/96 (45%)	43 (100%)	0	100	100
18	S	56/80 (70%)	56 (100%)	0	100	100
19	T	70/135 (52%)	70 (100%)	0	100	100
19	U	79/135 (58%)	79 (100%)	0	100	100
20	V	100/102 (98%)	100 (100%)	0	100	100
21	W	103/114 (90%)	103 (100%)	0	100	100
22	X	153/154 (99%)	153 (100%)	0	100	100
23	Y	105/107 (98%)	105 (100%)	0	100	100
24	Z	122/123 (99%)	122 (100%)	0	100	100
25	a	58/60 (97%)	58 (100%)	0	100	100
26	b	71/73 (97%)	71 (100%)	0	100	100
27	c	42/67 (63%)	42 (100%)	0	100	100
28	d	107/107 (100%)	107 (100%)	0	100	100
29	e	93/94 (99%)	92 (99%)	1 (1%)	65	74
30	f	49/53 (92%)	49 (100%)	0	100	100
31	g	92/129 (71%)	92 (100%)	0	100	100
32	h	123/162 (76%)	123 (100%)	0	100	100
33	i	88/118 (75%)	88 (100%)	0	100	100
34	j	58/87 (67%)	58 (100%)	0	100	100
35	k	58/78 (74%)	58 (100%)	0	100	100
36	l	139/161 (86%)	139 (100%)	0	100	100
37	m	112/114 (98%)	112 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
38	n	162/164 (99%)	162 (100%)	0	100	100
39	o	106/121 (88%)	105 (99%)	1 (1%)	70	75
40	p	153/158 (97%)	153 (100%)	0	100	100
41	r	57/96 (59%)	57 (100%)	0	100	100
42	s	31/95 (33%)	31 (100%)	0	100	100
All	All	6809/7660 (89%)	6798 (100%)	11 (0%)	85	87

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	D	356	TYR
6	F	257	ASN
7	G	491	ASN
7	G	675	ASP
8	H	271	LEU
12	L	49	LEU
12	L	249	SER
12	L	383	MET
16	P	109	VAL
29	e	102	GLU
39	o	46	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (90) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	137	HIS
3	C	38	GLN
3	C	53	HIS
3	C	102	ASN
3	C	124	HIS
4	D	149	ASN
5	E	37	ASN
5	E	42	HIS
5	E	57	GLN
5	E	67	ASN
6	F	144	ASN
6	F	261	HIS
6	F	361	GLN
6	F	373	ASN

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Mol	Chain	Res	Type
6	F	402	HIS
6	F	416	GLN
7	G	7	ASN
7	G	117	GLN
7	G	472	ASN
7	G	491	ASN
7	G	582	GLN
7	G	618	GLN
7	G	643	GLN
7	G	644	GLN
7	G	682	GLN
8	H	97	ASN
8	H	124	ASN
9	I	51	ASN
10	J	85	ASN
11	K	25	HIS
12	L	113	ASN
12	L	207	ASN
12	L	543	ASN
12	L	594	ASN
13	M	48	ASN
13	M	51	ASN
13	M	88	ASN
13	M	92	GLN
13	M	103	GLN
14	N	2	ASN
14	N	63	GLN
14	N	186	HIS
14	N	273	ASN
14	N	289	ASN
14	N	309	ASN
15	O	114	HIS
15	O	141	GLN
15	O	184	GLN
15	O	287	HIS
15	O	309	ASN
16	P	58	HIS
16	P	131	HIS
16	P	181	HIS
18	S	30	GLN
18	S	61	GLN
20	V	36	HIS

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Mol	Chain	Res	Type
20	V	40	HIS
21	W	60	GLN
21	W	110	HIS
22	X	68	ASN
24	Z	84	GLN
26	b	61	ASN
27	c	35	HIS
28	d	3	ASN
29	e	33	HIS
29	e	81	GLN
29	e	96	HIS
31	g	109	ASN
31	g	117	GLN
32	h	95	GLN
32	h	124	GLN
32	h	143	ASN
33	i	25	GLN
35	k	65	ASN
36	l	54	GLN
38	n	77	GLN
38	n	168	HIS
39	o	3	HIS
39	o	81	HIS
39	o	83	GLN
39	o	109	GLN
40	p	106	GLN
40	p	121	GLN
40	p	130	GLN
40	p	160	GLN
40	p	171	GLN
41	r	46	HIS
41	r	50	ASN
42	s	34	ASN
42	s	37	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

9 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
12	FME	L	1	12	8,9,10	0.94	0	7,9,11	0.89	0
4	2MR	D	85	4	10,12,13	1.84	1 (10%)	5,13,15	2.77	1 (20%)
8	FME	H	1	8	8,9,10	0.93	0	7,9,11	1.10	1 (14%)
1	FME	A	1	1	8,9,10	0.99	0	7,9,11	0.66	0
13	FME	M	1	13	8,9,10	0.95	0	7,9,11	0.69	0
10	FME	J	1	10	8,9,10	0.93	0	7,9,11	0.89	0
33	SAC	i	1	33	7,8,9	1.00	0	8,9,11	1.74	2 (25%)
11	FME	K	1	11	8,9,10	0.95	0	7,9,11	1.61	1 (14%)
14	FME	N	1	14	8,9,10	0.94	0	7,9,11	0.95	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	FME	L	1	12	-	3/7/9/11	-
4	2MR	D	85	4	-	4/10/13/15	-
8	FME	H	1	8	-	5/7/9/11	-
1	FME	A	1	1	-	0/7/9/11	-
13	FME	M	1	13	-	2/7/9/11	-
10	FME	J	1	10	-	3/7/9/11	-
33	SAC	i	1	33	-	5/7/8/10	-
11	FME	K	1	11	-	2/7/9/11	-
14	FME	N	1	14	-	0/7/9/11	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	85	2MR	CZ-NE	5.12	1.45	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	85	2MR	NE-CZ-NH2	5.95	124.93	119.48
11	K	1	FME	C-CA-N	3.62	116.26	109.73
33	i	1	SAC	C2A-C1A-N	2.94	121.07	116.10
33	i	1	SAC	CA-N-C1A	2.39	127.56	123.15
8	H	1	FME	CA-N-CN	2.22	126.23	122.82

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	85	2MR	C-CA-CB-CG
8	H	1	FME	O1-CN-N-CA
8	H	1	FME	CB-CA-N-CN
8	H	1	FME	C-CA-CB-CG
10	J	1	FME	O1-CN-N-CA
10	J	1	FME	O-C-CA-CB
11	K	1	FME	N-CA-CB-CG
11	K	1	FME	C-CA-CB-CG
12	L	1	FME	O1-CN-N-CA
33	i	1	SAC	CB-CA-N-C1A
4	D	85	2MR	NE-CD-CG-CB
33	i	1	SAC	C2A-C1A-N-CA
33	i	1	SAC	OAC-C1A-N-CA
12	L	1	FME	CA-CB-CG-SD
8	H	1	FME	CB-CG-SD-CE
33	i	1	SAC	N-CA-CB-OG
13	M	1	FME	N-CA-CB-CG
8	H	1	FME	N-CA-CB-CG
12	L	1	FME	CB-CA-N-CN
4	D	85	2MR	N-CA-CB-CG
4	D	85	2MR	CA-CB-CG-CD
33	i	1	SAC	C-CA-CB-OG
10	J	1	FME	C-CA-CB-CG
13	M	1	FME	CB-CA-N-CN

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	M	1	FME	2	0
33	i	1	SAC	2	0
11	K	1	FME	2	0

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 31 ligands modelled in this entry, 1 is monoatomic - leaving 30 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
47	3PE	J	201	-	32,32,50	1.05	4 (12%)	35,37,55	1.22	2 (5%)
45	FES	G	803	7	0,4,4	-	-	-	-	-
51	NAP	P	501	-	49,52,52	2.40	14 (28%)	69,80,80	2.01	16 (23%)
46	FMN	F	502	-	33,33,33	4.63	19 (57%)	48,50,50	4.80	22 (45%)
49	LMT	L	704	-	36,36,36	1.22	6 (16%)	47,47,47	1.04	2 (4%)
48	CDL	d	201	-	66,66,99	1.09	7 (10%)	72,78,111	1.19	5 (6%)
43	SF4	B	201	2	0,12,12	-	-	-	-	-
48	CDL	N	401	-	64,64,99	1.10	8 (12%)	70,76,111	1.19	5 (7%)
47	3PE	i	201	-	41,41,50	0.93	4 (9%)	44,46,55	1.11	2 (4%)
48	CDL	L	702	-	73,73,99	1.02	7 (9%)	79,85,111	1.09	4 (5%)
50	GTP	O	401	-	29,33,34	2.72	12 (41%)	44,52,54	1.66	10 (22%)
53	EHZ	T	201	19	29,36,37	1.87	5 (17%)	35,44,47	1.60	5 (14%)
53	EHZ	U	201	19	29,36,37	1.85	5 (17%)	35,44,47	1.59	6 (17%)
45	FES	E	301	5	0,4,4	-	-	-	-	-
47	3PE	H	401	-	43,43,50	0.91	4 (9%)	46,48,55	1.14	2 (4%)
44	PC1	H	402	-	41,41,53	1.42	6 (14%)	47,49,61	1.26	3 (6%)
47	3PE	Y	401	-	40,40,50	0.97	4 (10%)	43,45,55	1.14	2 (4%)
44	PC1	B	202	-	40,40,53	1.40	5 (12%)	46,48,61	1.08	2 (4%)
43	SF4	F	501	6	0,12,12	-	-	-	-	-
49	LMT	L	703	-	36,36,36	1.17	5 (13%)	47,47,47	0.99	4 (8%)
48	CDL	d	202	-	62,62,99	1.12	6 (9%)	68,74,111	1.17	4 (5%)
48	CDL	g	201	-	69,69,99	1.06	8 (11%)	75,81,111	1.23	6 (8%)
43	SF4	G	802	7	0,12,12	-	-	-	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
43	SF4	G	801	7	0,12,12	-	-	-		
47	3PE	I	201	-	50,50,50	0.84	3 (6%)	53,55,55	1.13	2 (3%)
48	CDL	M	502	-	58,58,99	1.08	5 (8%)	63,69,111	1.03	3 (4%)
47	3PE	L	701	-	48,48,50	0.88	3 (6%)	51,53,55	1.05	3 (5%)
43	SF4	I	202	9	0,12,12	-	-	-		
43	SF4	I	203	9	0,12,12	-	-	-		
47	3PE	M	501	-	41,41,50	0.95	3 (7%)	44,46,55	1.08	3 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
47	3PE	J	201	-	-	17/36/36/54	-
45	FES	G	803	7	-	-	0/1/1/1
51	NAP	P	501	-	-	5/35/67/67	0/5/5/5
46	FMN	F	502	-	-	9/18/18/18	0/3/3/3
49	LMT	L	704	-	-	13/21/61/61	0/2/2/2
48	CDL	d	201	-	-	30/77/77/110	-
48	CDL	N	401	-	-	39/75/75/110	-
50	GTP	O	401	-	-	7/22/34/38	0/3/3/3
47	3PE	i	201	-	-	16/45/45/54	-
48	CDL	L	702	-	-	32/84/84/110	-
53	EHZ	T	201	19	-	16/42/44/45	-
53	EHZ	U	201	19	-	18/42/44/45	-
43	SF4	B	201	2	-	-	0/6/5/5
45	FES	E	301	5	-	-	0/1/1/1
47	3PE	H	401	-	-	17/47/47/54	-
44	PC1	H	402	-	-	15/45/45/57	-
47	3PE	Y	401	-	-	18/44/44/54	-
44	PC1	B	202	-	-	16/44/44/57	-
43	SF4	F	501	6	-	-	0/6/5/5
49	LMT	L	703	-	-	13/21/61/61	0/2/2/2
48	CDL	d	202	-	-	35/73/73/110	-
48	CDL	g	201	-	-	46/80/80/110	-
43	SF4	G	802	7	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
43	SF4	G	801	7	-	-	0/6/5/5
47	3PE	I	201	-	-	19/54/54/54	-
48	CDL	M	502	-	-	39/67/67/110	-
47	3PE	L	701	-	-	23/52/52/54	-
43	SF4	I	202	9	-	-	0/6/5/5
43	SF4	I	203	9	-	-	0/6/5/5
47	3PE	M	501	-	-	20/45/45/54	-

All (143) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
51	P	501	NAP	P2B-O2B	11.66	1.81	1.59
46	F	502	FMN	C6-C5A	11.45	1.57	1.40
46	F	502	FMN	C9A-C5A	10.26	1.58	1.41
46	F	502	FMN	C9-C8	-10.06	1.24	1.39
46	F	502	FMN	O2-C2	8.71	1.40	1.24
50	O	401	GTP	O6-C6	8.53	1.39	1.23
46	F	502	FMN	O4-C4	8.42	1.39	1.23
46	F	502	FMN	C8-C7	-6.99	1.23	1.40
53	U	201	EHZ	C15-N2	5.97	1.46	1.33
46	F	502	FMN	C4A-C4	5.96	1.66	1.44
53	T	201	EHZ	C15-N2	5.83	1.46	1.33
53	T	201	EHZ	C12-N1	5.80	1.46	1.33
46	F	502	FMN	C2-N3	-5.58	1.26	1.39
53	U	201	EHZ	C12-N1	5.51	1.45	1.33
46	F	502	FMN	C4A-N5	4.67	1.39	1.30
50	O	401	GTP	C2-N2	4.62	1.45	1.34
50	O	401	GTP	C2-N1	4.37	1.48	1.37
51	P	501	NAP	C4N-C3N	4.30	1.46	1.39
50	O	401	GTP	C5-N7	4.24	1.47	1.39
51	P	501	NAP	C5N-C4N	4.13	1.47	1.38
50	O	401	GTP	C2-N3	3.95	1.42	1.33
51	P	501	NAP	PN-O5D	3.70	1.74	1.59
44	B	202	PC1	O31-C31	3.68	1.44	1.33
51	P	501	NAP	C3N-C7N	-3.67	1.45	1.50
44	H	402	PC1	O21-C21	3.62	1.44	1.34
44	H	402	PC1	O31-C31	3.54	1.43	1.33
44	B	202	PC1	O21-C21	3.44	1.44	1.34
46	F	502	FMN	C8M-C8	3.31	1.57	1.51
46	F	502	FMN	C4A-C10	3.28	1.53	1.44
51	P	501	NAP	O2B-C2B	-3.17	1.32	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
46	F	502	FMN	C9A-N10	3.01	1.46	1.41
49	L	703	LMT	O2'-C2'	-2.84	1.36	1.43
49	L	703	LMT	O3'-C3'	-2.84	1.36	1.43
49	L	704	LMT	O3'-C3'	-2.83	1.36	1.43
51	P	501	NAP	C7N-N7N	2.83	1.38	1.33
48	M	502	CDL	OB8-CB7	2.75	1.41	1.33
53	T	201	EHZ	C9-S1	2.74	1.82	1.76
47	M	501	3PE	O21-C2	-2.74	1.39	1.46
50	O	401	GTP	C8-N9	-2.72	1.31	1.37
48	L	702	CDL	OB6-CB4	-2.67	1.39	1.46
48	L	702	CDL	OB8-CB7	2.64	1.41	1.33
47	I	201	3PE	O21-C2	-2.63	1.40	1.46
48	g	201	CDL	OA6-CA4	-2.63	1.40	1.46
48	d	201	CDL	OB8-CB7	2.63	1.41	1.33
49	L	704	LMT	O2'-C2'	-2.63	1.36	1.43
47	i	201	3PE	O21-C2	-2.61	1.40	1.46
49	L	704	LMT	O2B-C2B	-2.60	1.36	1.43
48	N	401	CDL	OB6-CB4	-2.60	1.40	1.46
48	M	502	CDL	OB6-CB5	2.59	1.41	1.34
48	d	202	CDL	OA8-CA7	2.58	1.40	1.33
48	d	202	CDL	OB6-CB4	-2.58	1.40	1.46
44	B	202	PC1	O21-C2	-2.58	1.40	1.46
48	d	202	CDL	OB8-CB7	2.57	1.40	1.33
53	U	201	EHZ	C9-S1	2.56	1.82	1.76
48	d	202	CDL	OA6-CA4	-2.55	1.40	1.46
46	F	502	FMN	P-O2P	-2.55	1.45	1.54
48	N	401	CDL	OB8-CB7	2.53	1.40	1.33
47	L	701	3PE	O21-C2	-2.53	1.40	1.46
53	U	201	EHZ	O4-C15	-2.52	1.18	1.23
53	T	201	EHZ	O4-C15	-2.52	1.18	1.23
48	N	401	CDL	OA6-CA4	-2.51	1.40	1.46
48	d	201	CDL	OA6-CA4	-2.50	1.40	1.46
47	Y	401	3PE	O21-C2	-2.50	1.40	1.46
50	O	401	GTP	C2'-C3'	-2.50	1.46	1.52
49	L	703	LMT	O2B-C2B	-2.49	1.37	1.43
48	g	201	CDL	OB6-CB5	2.47	1.41	1.34
50	O	401	GTP	O4'-C1'	2.47	1.47	1.42
46	F	502	FMN	C5A-N5	2.46	1.44	1.39
48	d	201	CDL	OB6-CB5	2.46	1.41	1.34
47	J	201	3PE	O21-C2	-2.46	1.40	1.46
49	L	704	LMT	O3B-C3B	-2.44	1.37	1.43
48	g	201	CDL	OA8-CA6	-2.44	1.39	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
47	L	701	3PE	O31-C31	2.43	1.40	1.33
51	P	501	NAP	C6N-C5N	-2.43	1.33	1.38
46	F	502	FMN	P-O3P	-2.42	1.45	1.54
46	F	502	FMN	C10-N1	2.42	1.38	1.33
48	g	201	CDL	OB8-CB7	2.42	1.40	1.33
48	d	201	CDL	OA8-CA6	-2.40	1.39	1.45
49	L	703	LMT	O3B-C3B	-2.40	1.37	1.43
50	O	401	GTP	C1'-N9	-2.40	1.41	1.48
48	N	401	CDL	OB6-CB5	2.39	1.41	1.34
48	M	502	CDL	OA8-CA7	2.38	1.40	1.33
48	L	702	CDL	OA8-CA7	2.37	1.40	1.33
47	Y	401	3PE	O31-C3	-2.37	1.39	1.45
48	d	202	CDL	OB6-CB5	2.37	1.41	1.34
51	P	501	NAP	C2N-C3N	-2.36	1.35	1.39
48	N	401	CDL	OA8-CA7	2.36	1.40	1.33
47	I	201	3PE	O31-C3	-2.36	1.39	1.45
48	L	702	CDL	OA8-CA6	-2.36	1.39	1.45
48	d	201	CDL	OB6-CB4	-2.35	1.40	1.46
47	H	401	3PE	O21-C2	-2.33	1.40	1.46
47	M	501	3PE	O31-C31	2.32	1.40	1.33
51	P	501	NAP	O4B-C4B	-2.32	1.39	1.45
44	H	402	PC1	C22-C21	2.32	1.57	1.50
48	g	201	CDL	OB6-CB4	-2.32	1.40	1.46
47	H	401	3PE	O31-C3	-2.30	1.39	1.45
53	T	201	EHZ	O3-C12	-2.30	1.18	1.23
47	J	201	3PE	O31-C31	2.29	1.40	1.33
50	O	401	GTP	C4-N9	-2.28	1.32	1.38
44	H	402	PC1	O21-C2	-2.28	1.40	1.46
49	L	703	LMT	O4'-C4B	-2.27	1.37	1.43
48	d	201	CDL	OA8-CA7	2.27	1.40	1.33
48	d	201	CDL	OA6-CA5	2.27	1.40	1.34
53	U	201	EHZ	O3-C12	-2.25	1.18	1.23
47	i	201	3PE	O31-C3	-2.25	1.40	1.45
47	Y	401	3PE	O31-C31	2.25	1.39	1.33
48	N	401	CDL	OA6-CA5	2.24	1.40	1.34
48	M	502	CDL	OA8-CA6	-2.24	1.40	1.45
48	L	702	CDL	OA6-CA4	-2.24	1.41	1.46
50	O	401	GTP	PG-O2G	-2.23	1.46	1.54
50	O	401	GTP	PG-O3G	-2.23	1.46	1.54
46	F	502	FMN	O4'-C4'	-2.23	1.38	1.43
44	B	202	PC1	C22-C21	2.23	1.57	1.50
49	L	704	LMT	O1'-C1'	-2.23	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
47	L	701	3PE	O21-C21	2.21	1.40	1.34
47	Y	401	3PE	O21-C21	2.21	1.40	1.34
47	M	501	3PE	O31-C3	-2.20	1.40	1.45
47	H	401	3PE	O31-C31	2.19	1.39	1.33
47	J	201	3PE	O31-C3	-2.19	1.40	1.45
48	N	401	CDL	OA8-CA6	-2.19	1.40	1.45
51	P	501	NAP	O3D-C3D	-2.17	1.37	1.43
47	J	201	3PE	O21-C21	2.17	1.40	1.34
44	H	402	PC1	P-O11	2.16	1.68	1.59
44	B	202	PC1	P-O11	2.16	1.68	1.59
47	i	201	3PE	O31-C31	2.14	1.39	1.33
48	L	702	CDL	OA6-CA5	2.14	1.40	1.34
44	H	402	PC1	P-O13	2.14	1.67	1.59
49	L	704	LMT	O4'-C4B	-2.12	1.38	1.43
48	N	401	CDL	OB8-CB6	-2.12	1.40	1.45
48	g	201	CDL	OA6-CA5	2.11	1.40	1.34
51	P	501	NAP	O3B-C3B	-2.11	1.38	1.43
46	F	502	FMN	C7M-C7	2.10	1.55	1.51
48	L	702	CDL	OB6-CB5	2.10	1.40	1.34
47	I	201	3PE	O31-C31	2.09	1.39	1.33
48	g	201	CDL	OA8-CA7	2.09	1.39	1.33
47	H	401	3PE	O21-C21	2.08	1.40	1.34
51	P	501	NAP	O2D-C2D	-2.07	1.38	1.43
48	M	502	CDL	OB6-CB4	-2.07	1.41	1.46
51	P	501	NAP	O5D-C5D	-2.06	1.36	1.44
47	i	201	3PE	O21-C21	2.06	1.40	1.34
48	d	202	CDL	OA8-CA6	-2.04	1.40	1.45
46	F	502	FMN	C1'-C2'	2.04	1.55	1.52
48	g	201	CDL	OB8-CB6	-2.02	1.40	1.45

All (113) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
46	F	502	FMN	C6-C5A-N5	15.21	145.07	118.51
46	F	502	FMN	C4-C4A-N5	15.12	139.75	118.23
46	F	502	FMN	C6-C5A-C9A	-12.19	101.69	118.94
46	F	502	FMN	C5A-N5-C4A	9.65	134.11	118.07
46	F	502	FMN	C9A-C5A-N5	-8.46	113.23	122.43
46	F	502	FMN	C9A-N10-C10	7.95	133.16	120.77
51	P	501	NAP	PN-O3-PA	-7.54	106.96	132.83
51	P	501	NAP	C5N-C4N-C3N	-6.90	112.18	120.34
46	F	502	FMN	C5A-C9A-N10	-6.82	110.92	117.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
46	F	502	FMN	C10-N1-C2	6.70	130.30	116.90
46	F	502	FMN	C10-C4A-N5	-6.04	112.03	124.86
53	T	201	EHZ	C8-C9-S1	5.58	120.53	113.63
44	H	402	PC1	O21-C21-C22	5.14	122.58	111.50
46	F	502	FMN	C4-C4A-C10	-5.10	108.22	116.79
53	U	201	EHZ	C8-C9-S1	5.00	119.81	113.63
48	g	201	CDL	OA6-CA5-C11	4.75	121.73	111.50
48	d	201	CDL	OA6-CA5-C11	4.68	121.59	111.50
47	i	201	3PE	O21-C21-C22	4.59	121.39	111.50
47	H	401	3PE	O21-C21-C22	4.52	121.24	111.50
47	Y	401	3PE	O21-C21-C22	4.49	121.19	111.50
48	L	702	CDL	OA6-CA5-C11	4.46	121.11	111.50
50	O	401	GTP	C8-N9-C4	4.29	114.21	106.05
51	P	501	NAP	O7N-C7N-C3N	4.28	124.76	119.63
48	M	502	CDL	OB6-CB5-C51	4.27	120.71	111.50
47	J	201	3PE	O21-C21-C22	4.19	120.54	111.50
48	N	401	CDL	OA6-CA5-C11	4.18	120.52	111.50
48	d	202	CDL	OA6-CA5-C11	4.10	120.34	111.50
48	N	401	CDL	OB6-CB5-C51	4.10	120.33	111.50
46	F	502	FMN	C9-C9A-N10	4.04	127.29	121.84
47	I	201	3PE	O21-C21-C22	4.01	120.14	111.50
44	B	202	PC1	O21-C21-C22	3.93	119.97	111.50
48	d	201	CDL	OB6-CB5-C51	3.85	119.80	111.50
51	P	501	NAP	C2N-C3N-C4N	3.82	122.59	118.26
48	d	202	CDL	OB6-CB5-C51	3.81	119.71	111.50
48	g	201	CDL	OB6-CB5-C51	3.67	119.41	111.50
47	M	501	3PE	O21-C21-C22	3.53	119.12	111.50
51	P	501	NAP	O2B-P2B-O1X	-3.46	96.05	109.39
50	O	401	GTP	C2-N1-C6	-3.39	118.92	125.10
47	J	201	3PE	O31-C31-C32	3.38	122.53	111.91
50	O	401	GTP	PA-O3A-PB	-3.36	121.30	132.83
47	L	701	3PE	O31-C31-C32	3.04	121.44	111.91
50	O	401	GTP	PB-O3B-PG	-2.99	122.58	132.83
51	P	501	NAP	PA-O5B-C5B	-2.98	104.20	121.68
51	P	501	NAP	PN-O5D-C5D	-2.95	104.39	121.68
46	F	502	FMN	C9A-C9-C8	2.93	125.21	119.30
50	O	401	GTP	O2G-PG-O3B	2.93	114.48	104.64
50	O	401	GTP	O3G-PG-O3B	2.91	114.40	104.64
48	L	702	CDL	OB6-CB5-C51	2.91	117.76	111.50
46	F	502	FMN	C4A-C10-N1	-2.89	118.02	124.73
46	F	502	FMN	O3P-P-O5'	2.86	114.34	106.73
47	H	401	3PE	O31-C31-C32	2.85	120.86	111.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
47	L	701	3PE	O21-C21-C22	2.84	117.61	111.50
53	U	201	EHZ	C16-C15-N2	2.83	122.22	116.58
48	M	502	CDL	OB8-CB7-C71	2.80	120.70	111.91
44	H	402	PC1	O31-C31-C32	2.80	120.68	111.91
48	d	202	CDL	OA8-CA7-C31	2.78	120.62	111.91
48	d	202	CDL	OB8-CB7-C71	2.76	120.57	111.91
51	P	501	NAP	O7N-C7N-N7N	-2.75	118.68	122.58
47	I	201	3PE	O31-C31-C32	2.70	120.37	111.91
46	F	502	FMN	O2P-P-O5'	2.66	113.82	106.73
51	P	501	NAP	O4B-C4B-C3B	2.65	110.36	105.11
48	g	201	CDL	OA8-CA7-C31	2.64	120.20	111.91
46	F	502	FMN	C1'-N10-C9A	-2.63	116.12	120.51
48	L	702	CDL	OB8-CB7-C71	2.63	120.16	111.91
48	d	201	CDL	OA8-CA7-C31	2.62	120.14	111.91
48	g	201	CDL	OB8-CB7-C71	2.61	120.11	111.91
53	U	201	EHZ	C19-C17-C16	2.60	113.33	108.82
49	L	704	LMT	C3'-C4'-C5'	-2.60	104.97	110.93
46	F	502	FMN	O5'-P-O1P	2.59	113.74	106.47
48	g	201	CDL	CA4-OA6-CA5	-2.59	111.42	117.79
48	L	702	CDL	C33-C32-C31	-2.58	103.90	113.19
46	F	502	FMN	C5A-C6-C7	2.58	125.45	120.71
53	T	201	EHZ	C10-S1-C9	2.55	109.80	101.87
51	P	501	NAP	C5B-C4B-C3B	-2.54	105.66	115.18
51	P	501	NAP	C2A-N1A-C6A	-2.54	114.41	118.77
51	P	501	NAP	O3X-P2B-O2X	2.53	117.31	107.64
47	M	501	3PE	O31-C31-C32	2.52	119.82	111.91
47	Y	401	3PE	O31-C31-C32	2.52	119.81	111.91
48	d	201	CDL	OB8-CB7-C71	2.52	119.81	111.91
49	L	703	LMT	C1'-O5'-C5'	-2.47	108.84	113.69
50	O	401	GTP	C5-C6-N1	2.46	119.45	113.19
46	F	502	FMN	N10-C10-N1	2.46	125.43	118.35
44	B	202	PC1	O31-C31-C32	2.41	119.47	111.91
47	M	501	3PE	C2-O21-C21	-2.35	112.01	117.79
51	P	501	NAP	O2N-PN-O1N	2.35	123.84	112.24
46	F	502	FMN	C9-C8-C7	2.34	123.03	119.67
53	T	201	EHZ	C19-C17-C16	2.34	112.87	108.82
50	O	401	GTP	O6-C6-C5	-2.34	120.40	126.60
48	N	401	CDL	OB8-CB7-C71	2.33	119.22	111.91
46	F	502	FMN	C4-N3-C2	2.33	129.95	125.64
51	P	501	NAP	O5D-PN-O1N	-2.29	100.11	109.07
53	U	201	EHZ	C10-C11-N1	-2.28	107.62	112.42
53	U	201	EHZ	O2-C9-S1	-2.27	119.67	122.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
49	L	703	LMT	C3'-C4'-C5'	-2.24	105.79	110.93
53	T	201	EHZ	C16-C15-N2	2.24	121.03	116.58
48	N	401	CDL	OA8-CA7-C31	2.22	118.87	111.91
53	U	201	EHZ	C14-C13-C12	-2.22	108.66	112.36
47	i	201	3PE	O31-C31-C32	2.21	118.84	111.91
46	F	502	FMN	C6-C7-C8	2.20	122.83	119.67
51	P	501	NAP	N3A-C4A-N9A	2.19	130.69	127.08
48	N	401	CDL	CB4-OB6-CB5	-2.18	112.44	117.79
44	H	402	PC1	C2-O21-C21	-2.17	112.44	117.79
50	O	401	GTP	O2B-PB-O1B	-2.17	101.53	112.24
53	T	201	EHZ	C7-C8-C9	-2.13	109.02	113.89
49	L	703	LMT	O5B-C5B-C4B	2.12	113.55	109.69
51	P	501	NAP	C6N-C5N-C4N	2.10	122.49	119.44
50	O	401	GTP	O2A-PA-O1A	-2.10	101.85	112.24
48	g	201	CDL	OA6-CA5-OA7	-2.07	118.71	123.70
49	L	704	LMT	O5'-C1'-O1'	-2.06	105.09	109.97
48	d	201	CDL	CA6-CA4-CA3	-2.03	106.98	111.79
49	L	703	LMT	O5B-C5B-C6B	2.01	111.44	106.44
47	L	701	3PE	C23-C22-C21	-2.00	106.33	113.62
48	M	502	CDL	OA8-CA7-C31	2.00	118.19	111.91

There are no chirality outliers.

All (463) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
44	B	202	PC1	C1-O11-P-O14
44	B	202	PC1	O13-C11-C12-N
44	B	202	PC1	O22-C21-O21-C2
44	B	202	PC1	C22-C21-O21-C2
44	H	402	PC1	O21-C2-C3-O31
46	F	502	FMN	N10-C1'-C2'-O2'
46	F	502	FMN	C1'-C2'-C3'-C4'
46	F	502	FMN	C5'-O5'-P-O2P
46	F	502	FMN	C5'-O5'-P-O3P
47	H	401	3PE	C1-O11-P-O12
47	H	401	3PE	C1-O11-P-O14
47	H	401	3PE	C2-C1-O11-P
47	H	401	3PE	C22-C21-O21-C2
47	I	201	3PE	C1-O11-P-O14
47	J	201	3PE	C1-O11-P-O12
47	L	701	3PE	C1-O11-P-O12
47	L	701	3PE	C1-O11-P-O14

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Mol	Chain	Res	Type	Atoms
47	M	501	3PE	C11-O13-P-O11
47	M	501	3PE	C11-O13-P-O14
47	M	501	3PE	O22-C21-O21-C2
47	M	501	3PE	C22-C21-O21-C2
47	i	201	3PE	C1-O11-P-O12
47	i	201	3PE	C11-O13-P-O12
47	i	201	3PE	C12-C11-O13-P
47	i	201	3PE	C22-C21-O21-C2
48	L	702	CDL	CA2-OA2-PA1-OA3
48	L	702	CDL	CA2-OA2-PA1-OA4
48	L	702	CDL	CB2-OB2-PB2-OB3
48	L	702	CDL	C51-CB5-OB6-CB4
48	M	502	CDL	O1-C1-CB2-OB2
48	M	502	CDL	CA2-OA2-PA1-OA3
48	M	502	CDL	CA2-OA2-PA1-OA4
48	M	502	CDL	CA2-OA2-PA1-OA5
48	M	502	CDL	CA3-OA5-PA1-OA3
48	M	502	CDL	CB2-OB2-PB2-OB3
48	N	401	CDL	CA2-OA2-PA1-OA3
48	N	401	CDL	CA2-OA2-PA1-OA4
48	N	401	CDL	CA2-OA2-PA1-OA5
48	N	401	CDL	CA3-OA5-PA1-OA3
48	N	401	CDL	CA3-OA5-PA1-OA4
48	N	401	CDL	CA4-CA3-OA5-PA1
48	N	401	CDL	C11-CA5-OA6-CA4
48	N	401	CDL	CB2-OB2-PB2-OB4
48	d	201	CDL	O1-C1-CB2-OB2
48	d	201	CDL	CA2-C1-CB2-OB2
48	d	201	CDL	CA2-OA2-PA1-OA3
48	d	201	CDL	C51-CB5-OB6-CB4
48	d	202	CDL	O1-C1-CB2-OB2
48	d	202	CDL	CA3-OA5-PA1-OA2
48	d	202	CDL	CA3-OA5-PA1-OA3
48	d	202	CDL	CA3-OA5-PA1-OA4
48	g	201	CDL	CA2-OA2-PA1-OA3
48	g	201	CDL	CA2-OA2-PA1-OA4
48	g	201	CDL	CB2-OB2-PB2-OB3
48	g	201	CDL	CB2-OB2-PB2-OB4
48	g	201	CDL	CB3-OB5-PB2-OB3
48	g	201	CDL	CB3-OB5-PB2-OB4
49	L	703	LMT	C2'-C1'-O1'-C1
49	L	703	LMT	O5'-C1'-O1'-C1

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Mol	Chain	Res	Type	Atoms
49	L	704	LMT	O5'-C1'-O1'-C1
50	O	401	GTP	C5'-O5'-PA-O3A
50	O	401	GTP	C5'-O5'-PA-O1A
50	O	401	GTP	C5'-O5'-PA-O2A
53	T	201	EHZ	O1-C7-C8-C9
53	T	201	EHZ	C6-C7-C8-C9
53	T	201	EHZ	S1-C10-C11-N1
53	T	201	EHZ	C15-C16-C17-C20
53	T	201	EHZ	O5-C16-C17-C18
53	T	201	EHZ	O5-C16-C17-C20
53	T	201	EHZ	C16-C17-C20-O6
53	T	201	EHZ	C19-C17-C20-O6
53	T	201	EHZ	O2-C9-S1-C10
53	T	201	EHZ	C8-C9-S1-C10
53	U	201	EHZ	C5-C6-C7-C8
53	U	201	EHZ	S1-C10-C11-N1
53	U	201	EHZ	C12-C13-C14-N2
53	U	201	EHZ	C16-C15-N2-C14
53	U	201	EHZ	C15-C16-C17-C18
53	U	201	EHZ	C15-C16-C17-C19
53	U	201	EHZ	C15-C16-C17-C20
53	U	201	EHZ	O5-C16-C17-C18
53	U	201	EHZ	O5-C16-C17-C19
53	U	201	EHZ	O5-C16-C17-C20
53	U	201	EHZ	O2-C9-S1-C10
53	U	201	EHZ	C8-C9-S1-C10
47	I	201	3PE	O32-C31-O31-C3
48	d	202	CDL	OA9-CA7-OA8-CA6
48	d	202	CDL	C31-CA7-OA8-CA6
47	H	401	3PE	O32-C31-O31-C3
48	L	702	CDL	OB9-CB7-OB8-CB6
48	d	201	CDL	OB9-CB7-OB8-CB6
48	g	201	CDL	OB9-CB7-OB8-CB6
47	H	401	3PE	O22-C21-O21-C2
47	I	201	3PE	O22-C21-O21-C2
48	L	702	CDL	OB7-CB5-OB6-CB4
48	N	401	CDL	OA7-CA5-OA6-CA4
48	d	201	CDL	OB7-CB5-OB6-CB4
47	H	401	3PE	C32-C31-O31-C3
47	I	201	3PE	C32-C31-O31-C3
48	L	702	CDL	C71-CB7-OB8-CB6
48	g	201	CDL	C71-CB7-OB8-CB6

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Mol	Chain	Res	Type	Atoms
47	I	201	3PE	C22-C21-O21-C2
48	d	201	CDL	C71-CB7-OB8-CB6
47	i	201	3PE	O22-C21-O21-C2
47	Y	401	3PE	O32-C31-O31-C3
48	M	502	CDL	OB9-CB7-OB8-CB6
48	N	401	CDL	OB9-CB7-OB8-CB6
49	L	703	LMT	O5B-C1B-O1B-C4'
53	U	201	EHZ	O4-C15-N2-C14
48	M	502	CDL	C51-CB5-OB6-CB4
48	g	201	CDL	C11-CA5-OA6-CA4
48	g	201	CDL	C51-CB5-OB6-CB4
51	P	501	NAP	O4D-C4D-C5D-O5D
44	H	402	PC1	C32-C31-O31-C3
47	Y	401	3PE	C32-C31-O31-C3
48	M	502	CDL	C71-CB7-OB8-CB6
48	N	401	CDL	C71-CB7-OB8-CB6
48	g	201	CDL	OA7-CA5-OA6-CA4
48	N	401	CDL	C1-CA2-OA2-PA1
48	d	201	CDL	C1-CA2-OA2-PA1
49	L	703	LMT	O5'-C5'-C6'-O6'
48	M	502	CDL	CA2-C1-CB2-OB2
48	d	202	CDL	CA2-C1-CB2-OB2
48	M	502	CDL	OB7-CB5-OB6-CB4
48	g	201	CDL	OB7-CB5-OB6-CB4
44	H	402	PC1	O32-C31-O31-C3
47	J	201	3PE	C32-C31-O31-C3
48	d	201	CDL	C31-CA7-OA8-CA6
48	g	201	CDL	C31-CA7-OA8-CA6
47	L	701	3PE	C32-C33-C34-C35
48	L	702	CDL	OA5-CA3-CA4-OA6
49	L	703	LMT	C4'-C5'-C6'-O6'
49	L	704	LMT	C2'-C1'-O1'-C1
47	J	201	3PE	O32-C31-O31-C3
48	d	201	CDL	OA9-CA7-OA8-CA6
47	Y	401	3PE	C31-C32-C33-C34
48	M	502	CDL	CB5-C51-C52-C53
48	d	202	CDL	CB5-C51-C52-C53
49	L	704	LMT	O5'-C5'-C6'-O6'
48	L	702	CDL	CB5-C51-C52-C53
48	d	201	CDL	CB5-C51-C52-C53
48	g	201	CDL	CA5-C11-C12-C13
53	U	201	EHZ	C5-C6-C7-O1

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Mol	Chain	Res	Type	Atoms
48	g	201	CDL	OA9-CA7-OA8-CA6
47	L	701	3PE	C32-C31-O31-C3
48	d	202	CDL	CA5-C11-C12-C13
48	N	401	CDL	O1-C1-CB2-OB2
44	B	202	PC1	C1-O11-P-O13
47	H	401	3PE	C1-O11-P-O13
47	I	201	3PE	C1-O11-P-O13
47	I	201	3PE	C11-O13-P-O11
47	J	201	3PE	C1-O11-P-O13
47	L	701	3PE	C1-O11-P-O13
47	Y	401	3PE	C1-O11-P-O13
47	i	201	3PE	C1-O11-P-O13
47	i	201	3PE	C11-O13-P-O11
48	L	702	CDL	CA2-OA2-PA1-OA5
48	L	702	CDL	CA3-OA5-PA1-OA2
48	L	702	CDL	CB2-OB2-PB2-OB5
48	M	502	CDL	CA3-OA5-PA1-OA2
48	M	502	CDL	CB2-OB2-PB2-OB5
48	N	401	CDL	CA3-OA5-PA1-OA2
48	N	401	CDL	CB2-OB2-PB2-OB5
48	d	202	CDL	CB2-OB2-PB2-OB5
48	g	201	CDL	CA2-OA2-PA1-OA5
48	g	201	CDL	CA3-OA5-PA1-OA2
48	g	201	CDL	CB2-OB2-PB2-OB5
48	g	201	CDL	CB3-OB5-PB2-OB2
48	M	502	CDL	C31-CA7-OA8-CA6
48	d	202	CDL	C78-C79-C80-C81
48	N	401	CDL	CA2-C1-CB2-OB2
47	J	201	3PE	C34-C35-C36-C37
53	T	201	EHZ	C18-C17-C20-O6
47	I	201	3PE	C32-C33-C34-C35
53	U	201	EHZ	C2-C3-C4-C5
48	M	502	CDL	C53-C54-C55-C56
48	d	202	CDL	C75-C76-C77-C78
48	N	401	CDL	O1-C1-CA2-OA2
48	g	201	CDL	O1-C1-CB2-OB2
47	i	201	3PE	C25-C26-C27-C28
47	I	201	3PE	C31-C32-C33-C34
48	N	401	CDL	OA6-CA4-CA6-OA8
48	N	401	CDL	C51-C52-C53-C54
47	L	701	3PE	O32-C31-O31-C3
47	Y	401	3PE	C33-C34-C35-C36

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Mol	Chain	Res	Type	Atoms
48	N	401	CDL	C75-C76-C77-C78
48	N	401	CDL	OB7-CB5-OB6-CB4
48	d	202	CDL	OA7-CA5-OA6-CA4
48	N	401	CDL	C51-CB5-OB6-CB4
48	d	202	CDL	C11-CA5-OA6-CA4
53	U	201	EHZ	C21-C1-C2-C3
47	I	201	3PE	C21-C22-C23-C24
48	L	702	CDL	C77-C78-C79-C80
49	L	704	LMT	C7-C8-C9-C10
47	L	701	3PE	C2-C3-O31-C31
47	L	701	3PE	O13-C11-C12-N
48	d	201	CDL	C55-C56-C57-C58
44	H	402	PC1	C27-C28-C29-C2A
48	g	201	CDL	C51-C52-C53-C54
47	M	501	3PE	C3B-C3C-C3D-C3E
48	d	202	CDL	C32-C33-C34-C35
48	M	502	CDL	OA9-CA7-OA8-CA6
49	L	703	LMT	C11-C10-C9-C8
47	I	201	3PE	C25-C26-C27-C28
46	F	502	FMN	O2'-C2'-C3'-O3'
48	N	401	CDL	CA5-C11-C12-C13
47	M	501	3PE	C35-C36-C37-C38
48	L	702	CDL	C14-C15-C16-C17
46	F	502	FMN	O2'-C2'-C3'-C4'
44	H	402	PC1	C22-C23-C24-C25
48	M	502	CDL	OA6-CA4-CA6-OA8
44	B	202	PC1	C37-C38-C39-C3A
48	N	401	CDL	C78-C79-C80-C81
48	d	201	CDL	C57-C58-C59-C60
48	g	201	CDL	C19-C20-C21-C22
47	M	501	3PE	C3A-C3B-C3C-C3D
48	M	502	CDL	CA3-CA4-CA6-OA8
47	i	201	3PE	C33-C34-C35-C36
48	g	201	CDL	C23-C24-C25-C26
48	L	702	CDL	C11-CA5-OA6-CA4
48	L	702	CDL	OA7-CA5-OA6-CA4
47	M	501	3PE	O21-C2-C3-O31
48	g	201	CDL	C21-C22-C23-C24
48	d	201	CDL	C31-C32-C33-C34
47	M	501	3PE	C37-C38-C39-C3A
48	N	401	CDL	C76-C77-C78-C79
48	N	401	CDL	C53-C54-C55-C56

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Mol	Chain	Res	Type	Atoms
47	M	501	3PE	C1-O11-P-O13
47	Y	401	3PE	C11-O13-P-O11
48	d	201	CDL	CA2-OA2-PA1-OA5
48	d	202	CDL	C51-C52-C53-C54
47	I	201	3PE	O11-C1-C2-C3
48	L	702	CDL	OA5-CA3-CA4-CA6
47	I	201	3PE	C2D-C2E-C2F-C2G
47	i	201	3PE	C24-C25-C26-C27
49	L	703	LMT	C5'-C4'-O1B-C1B
48	L	702	CDL	C52-C53-C54-C55
53	T	201	EHZ	C3-C4-C5-C6
44	B	202	PC1	C1-C2-C3-O31
48	N	401	CDL	CB3-CB4-CB6-OB8
48	g	201	CDL	CB3-CB4-CB6-OB8
47	L	701	3PE	O21-C21-C22-C23
48	d	202	CDL	C80-C81-C82-C83
47	J	201	3PE	C22-C21-O21-C2
47	I	201	3PE	C37-C38-C39-C3A
48	L	702	CDL	C73-C74-C75-C76
47	H	401	3PE	C3-C2-O21-C21
48	L	702	CDL	CA6-CA4-OA6-CA5
46	F	502	FMN	C5'-O5'-P-O1P
47	L	701	3PE	C27-C28-C29-C2A
48	L	702	CDL	OB5-CB3-CB4-OB6
48	g	201	CDL	OB5-CB3-CB4-OB6
47	i	201	3PE	C22-C23-C24-C25
47	H	401	3PE	C36-C37-C38-C39
48	d	202	CDL	C76-C77-C78-C79
47	M	501	3PE	C31-C32-C33-C34
49	L	703	LMT	C4B-C5B-C6B-O6B
48	g	201	CDL	CB5-C51-C52-C53
50	O	401	GTP	O4'-C4'-C5'-O5'
47	L	701	3PE	O11-C1-C2-C3
48	L	702	CDL	OB5-CB3-CB4-CB6
48	M	502	CDL	OB5-CB3-CB4-CB6
48	g	201	CDL	C72-C73-C74-C75
47	L	701	3PE	C22-C23-C24-C25
48	d	201	CDL	CA5-C11-C12-C13
47	M	501	3PE	C33-C34-C35-C36
48	M	502	CDL	C55-C56-C57-C58
48	d	202	CDL	C71-CB7-OB8-CB6
47	J	201	3PE	O31-C31-C32-C33

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Mol	Chain	Res	Type	Atoms
47	L	701	3PE	C36-C37-C38-C39
48	N	401	CDL	CA7-C31-C32-C33
48	M	502	CDL	C75-C76-C77-C78
47	J	201	3PE	C1-C2-C3-O31
48	M	502	CDL	CB3-CB4-CB6-OB8
48	N	401	CDL	CA3-CA4-CA6-OA8
48	g	201	CDL	CA3-CA4-CA6-OA8
47	M	501	3PE	C22-C23-C24-C25
53	U	201	EHZ	C1-C21-C22-C23
44	B	202	PC1	C3B-C3C-C3D-C3E
47	M	501	3PE	O11-C1-C2-O21
44	B	202	PC1	O21-C2-C3-O31
47	L	701	3PE	O21-C2-C3-O31
48	L	702	CDL	OB6-CB4-CB6-OB8
48	M	502	CDL	OB6-CB4-CB6-OB8
48	N	401	CDL	OB6-CB4-CB6-OB8
48	g	201	CDL	OB6-CB4-CB6-OB8
48	d	202	CDL	C74-C75-C76-C77
47	L	701	3PE	C24-C25-C26-C27
47	J	201	3PE	O22-C21-O21-C2
48	g	201	CDL	C20-C21-C22-C23
49	L	703	LMT	C5-C6-C7-C8
49	L	704	LMT	C2B-C1B-O1B-C4'
48	M	502	CDL	C76-C77-C78-C79
48	g	201	CDL	C17-C18-C19-C20
44	H	402	PC1	C34-C35-C36-C37
51	P	501	NAP	C3D-C4D-C5D-O5D
50	O	401	GTP	PB-O3B-PG-O1G
48	M	502	CDL	C52-C51-CB5-OB6
48	d	201	CDL	C54-C55-C56-C57
49	L	703	LMT	C9-C10-C11-C12
47	i	201	3PE	C3-C2-O21-C21
48	M	502	CDL	CB6-CB4-OB6-CB5
48	d	201	CDL	CB3-CB4-OB6-CB5
48	d	202	CDL	CA6-CA4-OA6-CA5
47	H	401	3PE	C34-C35-C36-C37
49	L	704	LMT	O5B-C1B-O1B-C4'
44	H	402	PC1	C1-C2-C3-O31
48	d	202	CDL	CB3-CB4-CB6-OB8
44	H	402	PC1	C37-C38-C39-C3A
53	U	201	EHZ	C3-C4-C5-C6
47	H	401	3PE	O11-C1-C2-O21

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Mol	Chain	Res	Type	Atoms
47	I	201	3PE	O11-C1-C2-O21
48	M	502	CDL	OB5-CB3-CB4-OB6
47	H	401	3PE	C22-C23-C24-C25
48	L	702	CDL	C71-C72-C73-C74
48	M	502	CDL	C31-C32-C33-C34
48	d	202	CDL	OB9-CB7-OB8-CB6
48	N	401	CDL	C71-C72-C73-C74
48	d	202	CDL	OB7-CB5-OB6-CB4
47	J	201	3PE	C2-C3-O31-C31
48	L	702	CDL	C75-C76-C77-C78
48	L	702	CDL	C55-C56-C57-C58
44	B	202	PC1	C3C-C3D-C3E-C3F
44	H	402	PC1	C1-O11-P-O13
48	d	201	CDL	CB2-OB2-PB2-OB5
48	g	201	CDL	O1-C1-CA2-OA2
48	M	502	CDL	C58-C59-C60-C61
48	L	702	CDL	C1-CA2-OA2-PA1
47	I	201	3PE	C11-O13-P-O14
47	M	501	3PE	C1-O11-P-O12
47	M	501	3PE	C1-O11-P-O14
47	Y	401	3PE	C1-O11-P-O12
47	Y	401	3PE	C1-O11-P-O14
47	Y	401	3PE	C11-O13-P-O12
47	Y	401	3PE	C11-O13-P-O14
47	i	201	3PE	C11-O13-P-O14
48	L	702	CDL	CA3-OA5-PA1-OA4
48	M	502	CDL	CA3-OA5-PA1-OA4
48	M	502	CDL	CB2-OB2-PB2-OB4
48	N	401	CDL	CB2-OB2-PB2-OB3
48	d	201	CDL	CA2-OA2-PA1-OA4
48	d	202	CDL	CB2-OB2-PB2-OB3
48	d	202	CDL	CB2-OB2-PB2-OB4
48	g	201	CDL	CA3-OA5-PA1-OA4
47	H	401	3PE	O11-C1-C2-C3
47	Y	401	3PE	O11-C1-C2-C3
48	g	201	CDL	OB5-CB3-CB4-CB6
48	d	202	CDL	C51-CB5-OB6-CB4
44	B	202	PC1	C12-C11-O13-P
47	I	201	3PE	C12-C11-O13-P
47	M	501	3PE	C12-C11-O13-P
49	L	703	LMT	C3'-C4'-O1B-C1B
47	L	701	3PE	O11-C1-C2-O21

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Mol	Chain	Res	Type	Atoms
47	Y	401	3PE	O11-C1-C2-O21
48	d	201	CDL	OB5-CB3-CB4-OB6
49	L	704	LMT	C5-C6-C7-C8
44	H	402	PC1	O13-C11-C12-N
47	J	201	3PE	O21-C2-C3-O31
48	d	202	CDL	OB6-CB4-CB6-OB8
48	g	201	CDL	C24-C25-C26-C27
48	d	202	CDL	C71-C72-C73-C74
48	L	702	CDL	C56-C57-C58-C59
48	d	202	CDL	C72-C73-C74-C75
48	d	201	CDL	C71-C72-C73-C74
47	L	701	3PE	C1-C2-O21-C21
48	M	502	CDL	CA4-CA6-OA8-CA7
47	I	201	3PE	C35-C36-C37-C38
48	M	502	CDL	C1-CB2-OB2-PB2
44	B	202	PC1	C38-C39-C3A-C3B
47	i	201	3PE	O11-C1-C2-O21
48	N	401	CDL	C84-C85-C86-C87
49	L	704	LMT	O1'-C1-C2-C3
48	g	201	CDL	OA6-CA4-CA6-OA8
44	B	202	PC1	C11-O13-P-O11
47	L	701	3PE	C11-O13-P-O11
48	d	201	CDL	CA3-OA5-PA1-OA2
48	d	202	CDL	CB3-OB5-PB2-OB2
47	Y	401	3PE	C35-C36-C37-C38
49	L	703	LMT	O1'-C1-C2-C3
47	M	501	3PE	C1-C2-C3-O31
51	P	501	NAP	PN-O3-PA-O1A
48	d	201	CDL	C32-C33-C34-C35
47	L	701	3PE	O22-C21-C22-C23
47	Y	401	3PE	C2-C1-O11-P
48	M	502	CDL	OA5-CA3-CA4-CA6
49	L	704	LMT	C4'-C5'-C6'-O6'
48	M	502	CDL	C56-C57-C58-C59
44	H	402	PC1	C24-C25-C26-C27
48	M	502	CDL	OA5-CA3-CA4-OA6
47	i	201	3PE	O21-C21-C22-C23
47	Y	401	3PE	O21-C2-C3-O31
47	L	701	3PE	C2-C1-O11-P
48	N	401	CDL	C1-CB2-OB2-PB2
48	g	201	CDL	C53-C54-C55-C56
47	L	701	3PE	C1-C2-C3-O31

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Mol	Chain	Res	Type	Atoms
48	g	201	CDL	CB3-CB4-OB6-CB5
48	d	202	CDL	CA7-C31-C32-C33
48	M	502	CDL	C51-C52-C53-C54
47	i	201	3PE	O11-C1-C2-C3
48	N	401	CDL	OB5-CB3-CB4-CB6
44	B	202	PC1	C35-C36-C37-C38
48	d	201	CDL	C51-C52-C53-C54
48	d	201	CDL	C52-C53-C54-C55
47	I	201	3PE	C2-C1-O11-P
49	L	704	LMT	C6-C7-C8-C9
48	d	201	CDL	OB5-CB3-CB4-CB6
44	H	402	PC1	C39-C3A-C3B-C3C
47	M	501	3PE	C24-C25-C26-C27
49	L	703	LMT	C4-C5-C6-C7
47	J	201	3PE	O32-C31-C32-C33
53	T	201	EHZ	C21-C22-C23-C24
48	L	702	CDL	C12-C13-C14-C15
47	J	201	3PE	O21-C21-C22-C23
48	g	201	CDL	C12-C11-CA5-OA6
47	H	401	3PE	C38-C39-C3A-C3B
47	Y	401	3PE	O31-C31-C32-C33
48	L	702	CDL	CB3-CB4-CB6-OB8
44	B	202	PC1	O11-C1-C2-O21
48	g	201	CDL	C52-C51-CB5-OB6
50	O	401	GTP	PB-O3B-PG-O2G
50	O	401	GTP	PB-O3B-PG-O3G
53	T	201	EHZ	C15-C16-C17-C18
53	T	201	EHZ	C15-C16-C17-C19
47	J	201	3PE	O11-C1-C2-C3
47	M	501	3PE	O11-C1-C2-C3
47	H	401	3PE	C24-C25-C26-C27
44	H	402	PC1	C33-C34-C35-C36
47	L	701	3PE	C39-C3A-C3B-C3C
48	N	401	CDL	C12-C11-CA5-OA6
44	H	402	PC1	C3A-C3B-C3C-C3D
48	g	201	CDL	C18-C19-C20-C21
47	H	401	3PE	C35-C36-C37-C38
51	P	501	NAP	O4B-C4B-C5B-O5B
47	L	701	3PE	C26-C27-C28-C29
51	P	501	NAP	PN-O3-PA-O2A
49	L	704	LMT	O5B-C5B-C6B-O6B
48	d	201	CDL	C52-C51-CB5-OB6

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Mol	Chain	Res	Type	Atoms
48	g	201	CDL	C12-C11-CA5-OA7
47	J	201	3PE	C25-C26-C27-C28
53	T	201	EHZ	C2-C1-C21-C22
48	L	702	CDL	C17-C18-C19-C20
48	M	502	CDL	CB3-OB5-PB2-OB4
48	d	201	CDL	CB2-OB2-PB2-OB3
48	d	202	CDL	CB3-OB5-PB2-OB3
49	L	704	LMT	C3-C4-C5-C6
48	N	401	CDL	OA5-CA3-CA4-CA6
48	N	401	CDL	C12-C11-CA5-OA7
47	J	201	3PE	O22-C21-C22-C23
47	Y	401	3PE	O32-C31-C32-C33
48	g	201	CDL	C52-C51-CB5-OB7
46	F	502	FMN	C1'-C2'-C3'-O3'
48	g	201	CDL	CB6-CB4-OB6-CB5
48	g	201	CDL	C32-C31-CA7-OA8
44	B	202	PC1	C3F-C3G-C3H-C3I
44	H	402	PC1	O21-C21-C22-C23
48	d	202	CDL	C52-C51-CB5-OB6
48	N	401	CDL	CB2-C1-CA2-OA2
48	d	201	CDL	C52-C51-CB5-OB7
46	F	502	FMN	N10-C1'-C2'-C3'
47	J	201	3PE	O11-C1-C2-O21
47	Y	401	3PE	C34-C35-C36-C37
49	L	704	LMT	C2-C1-O1'-C1'
48	d	202	CDL	C52-C51-CB5-OB7

There are no ring outliers.

19 monomers are involved in 29 short contacts:

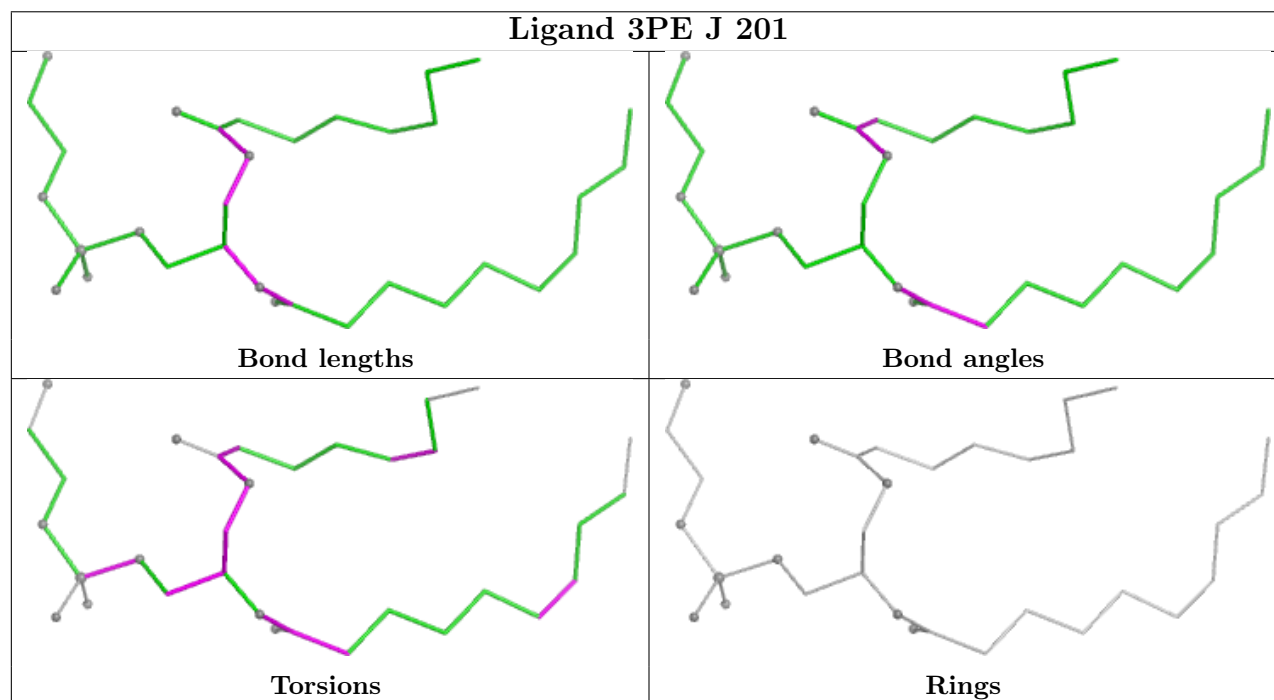
Mol	Chain	Res	Type	Clashes	Symm-Clashes
45	G	803	FES	1	0
51	P	501	NAP	1	0
46	F	502	FMN	2	0
49	L	704	LMT	1	0
47	i	201	3PE	3	0
48	L	702	CDL	1	0
50	O	401	GTP	2	0
53	T	201	EHZ	1	0
53	U	201	EHZ	1	0
45	E	301	FES	1	0
47	H	401	3PE	1	0

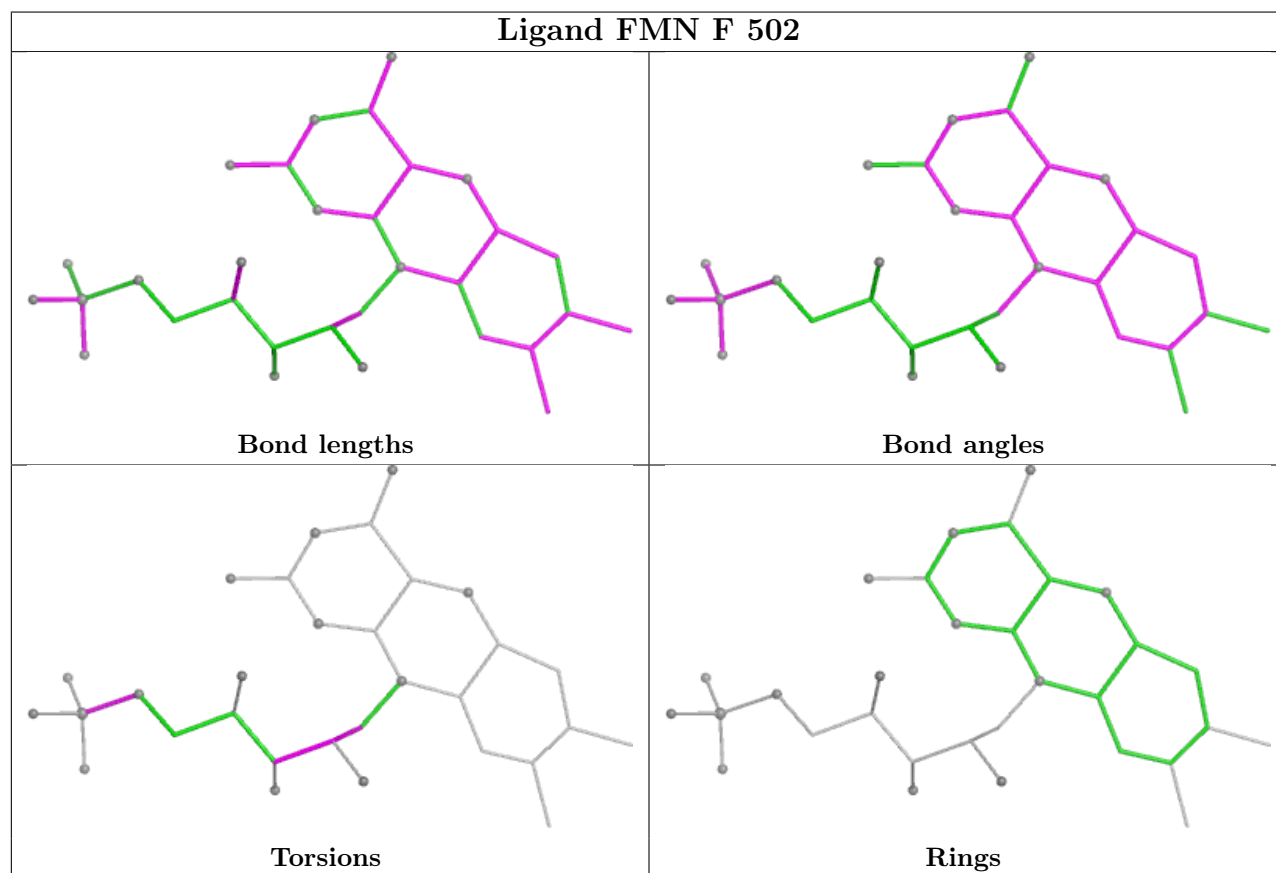
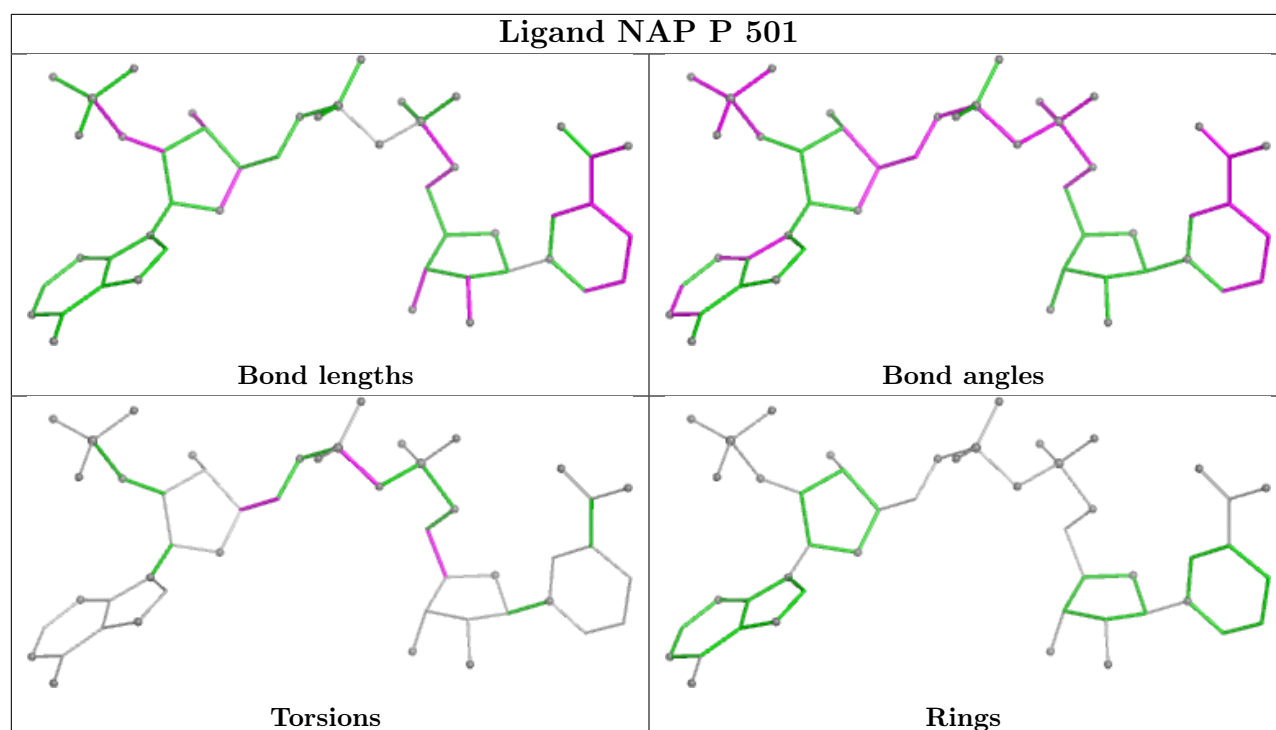
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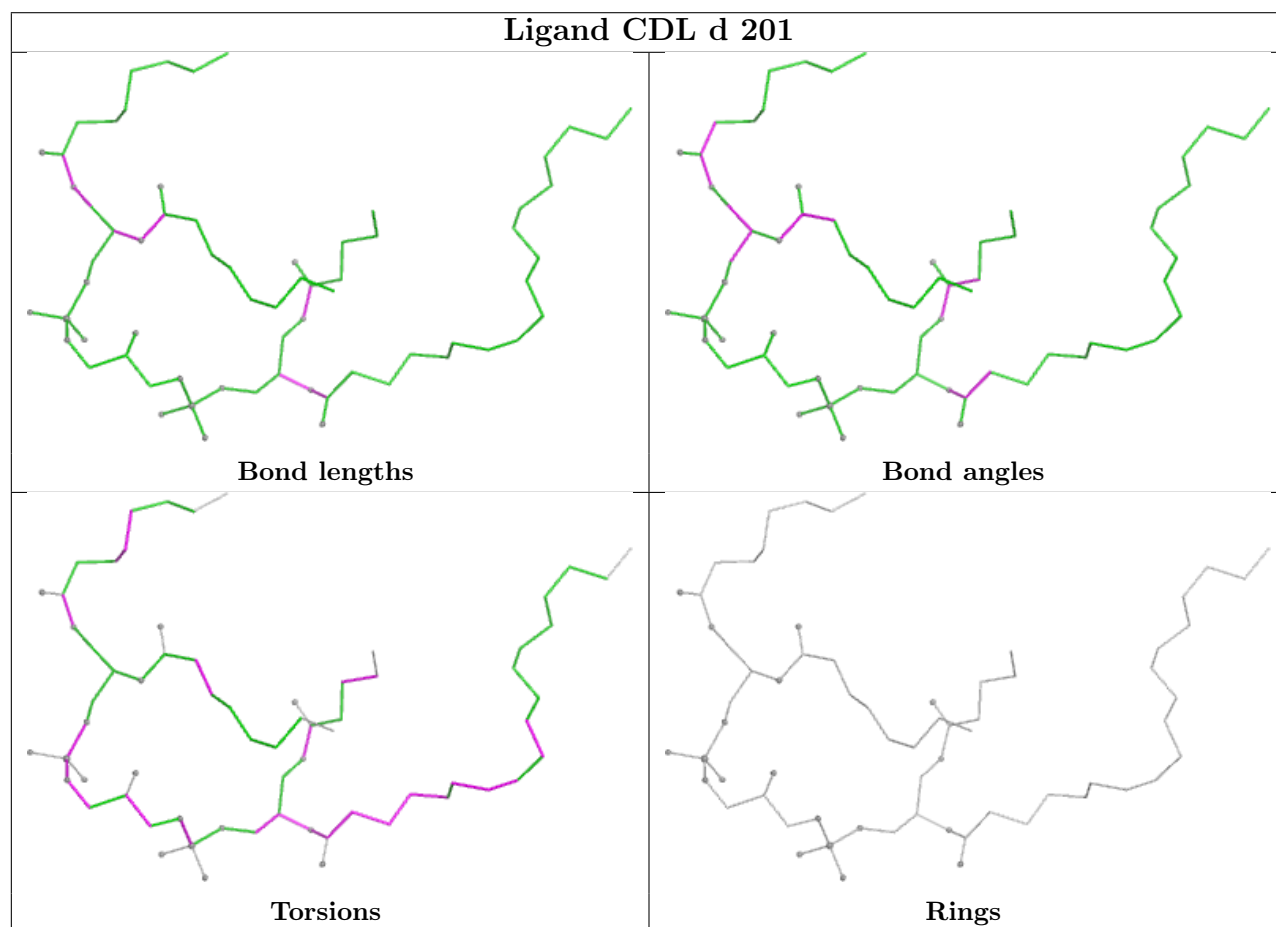
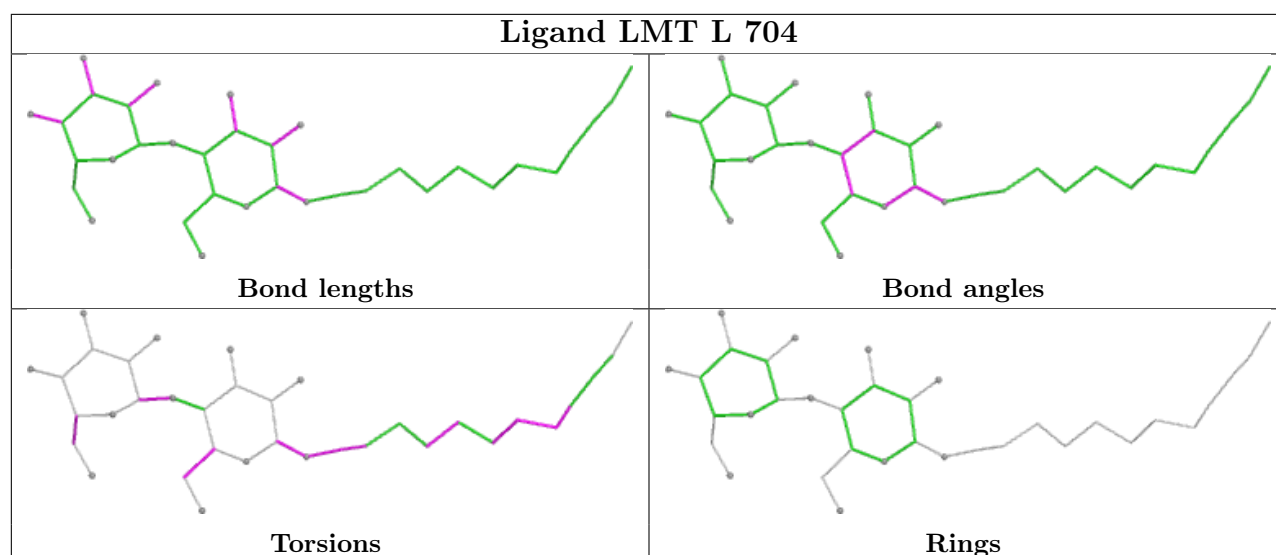
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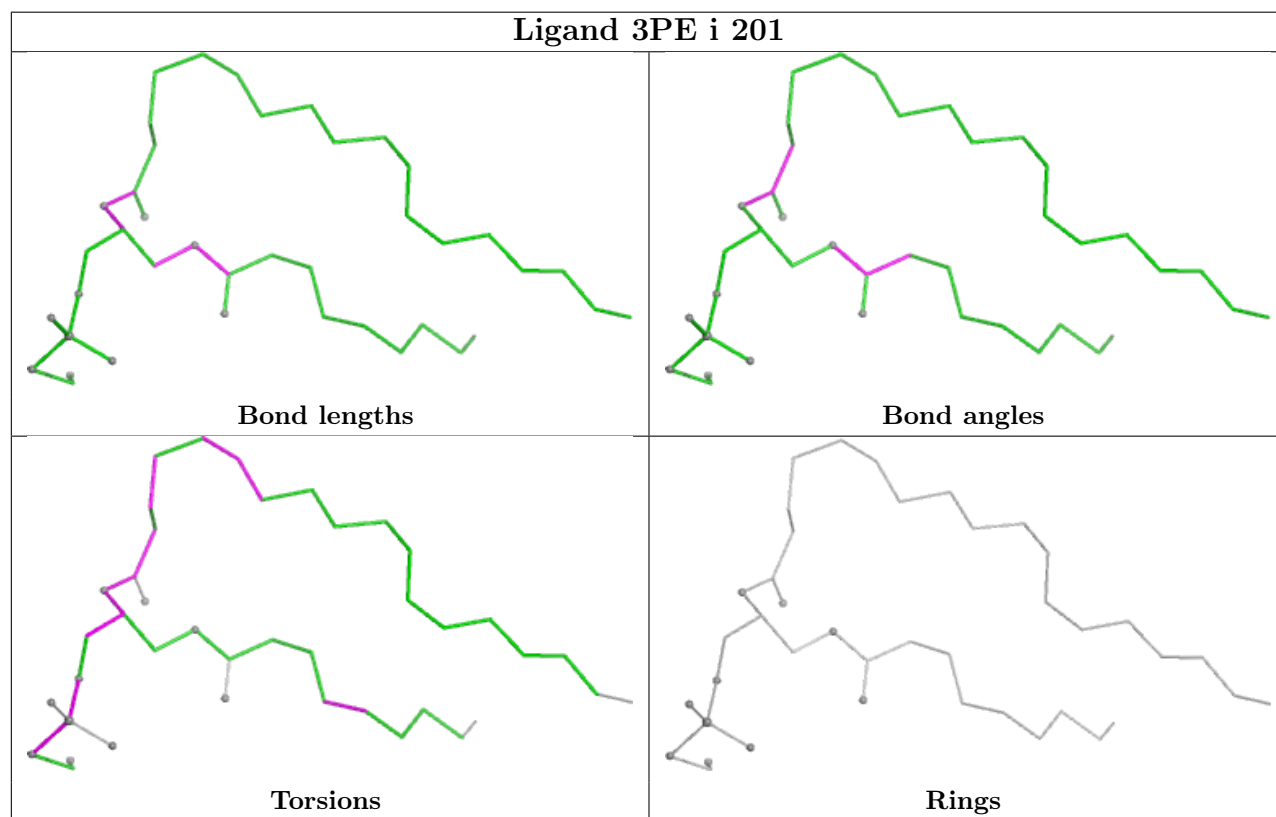
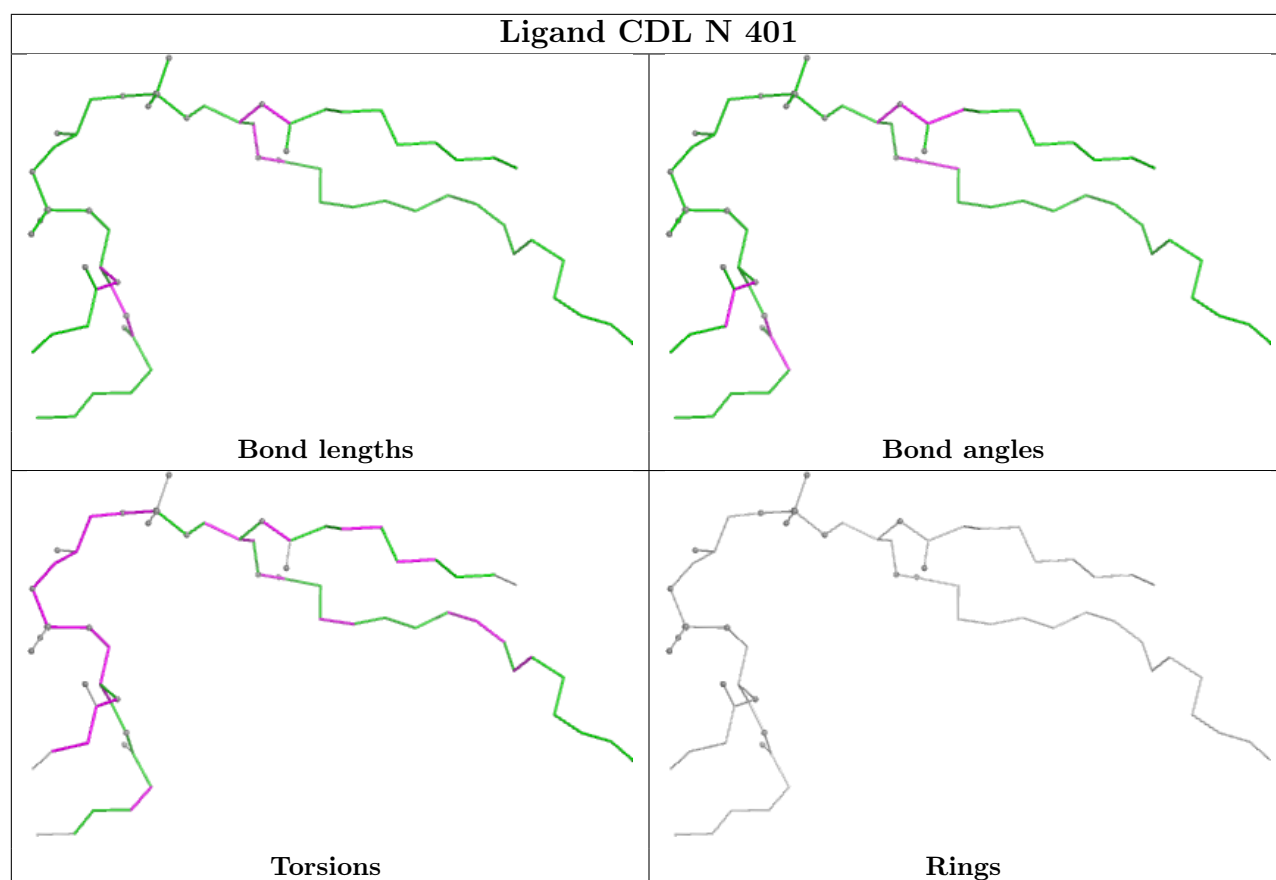
Mol	Chain	Res	Type	Clashes	Symm-Clashes
44	H	402	PC1	3	0
43	F	501	SF4	1	0
43	G	802	SF4	2	0
47	I	201	3PE	1	0
48	M	502	CDL	3	0
47	L	701	3PE	1	0
43	I	202	SF4	2	0
47	M	501	3PE	1	0

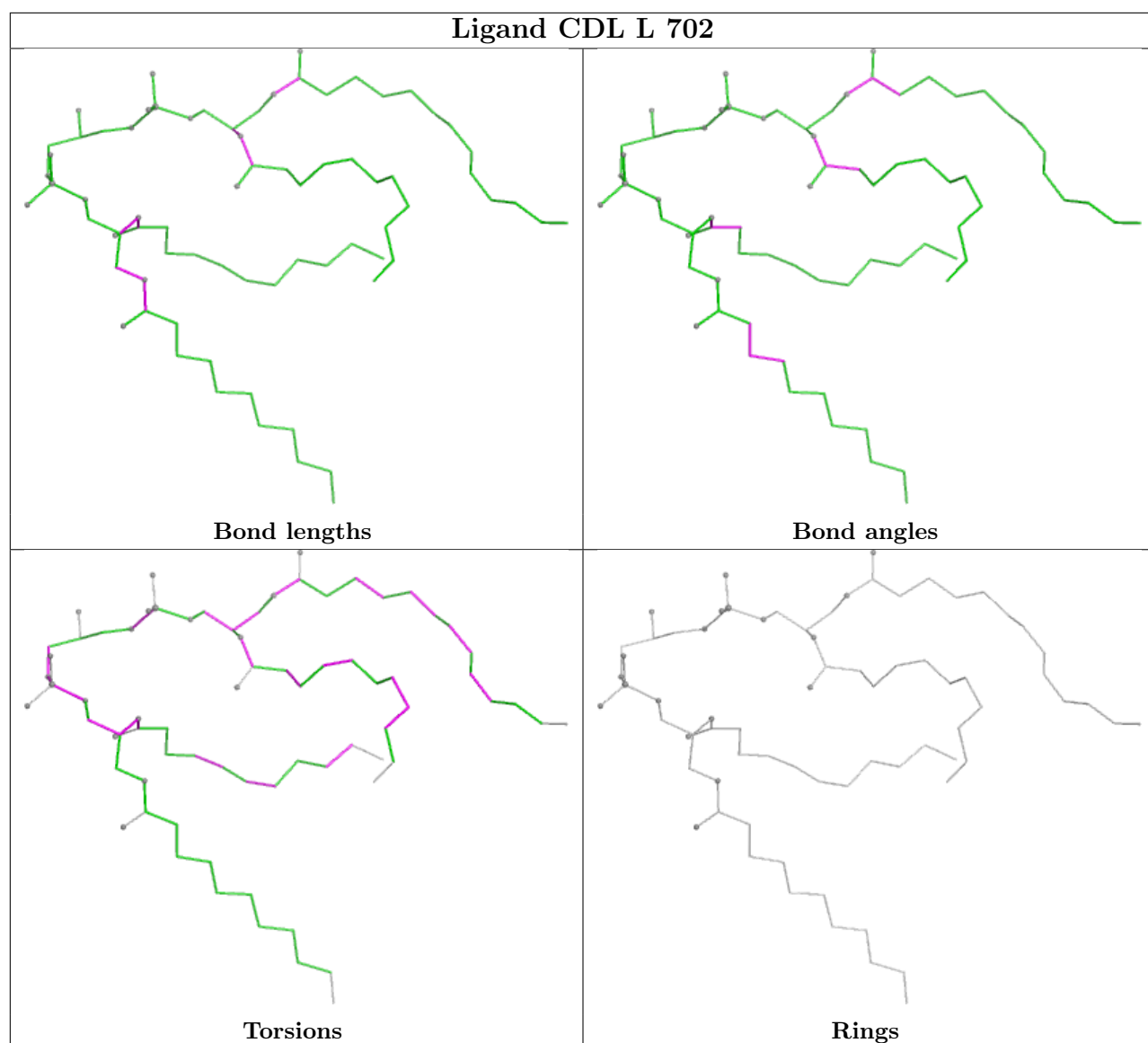
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

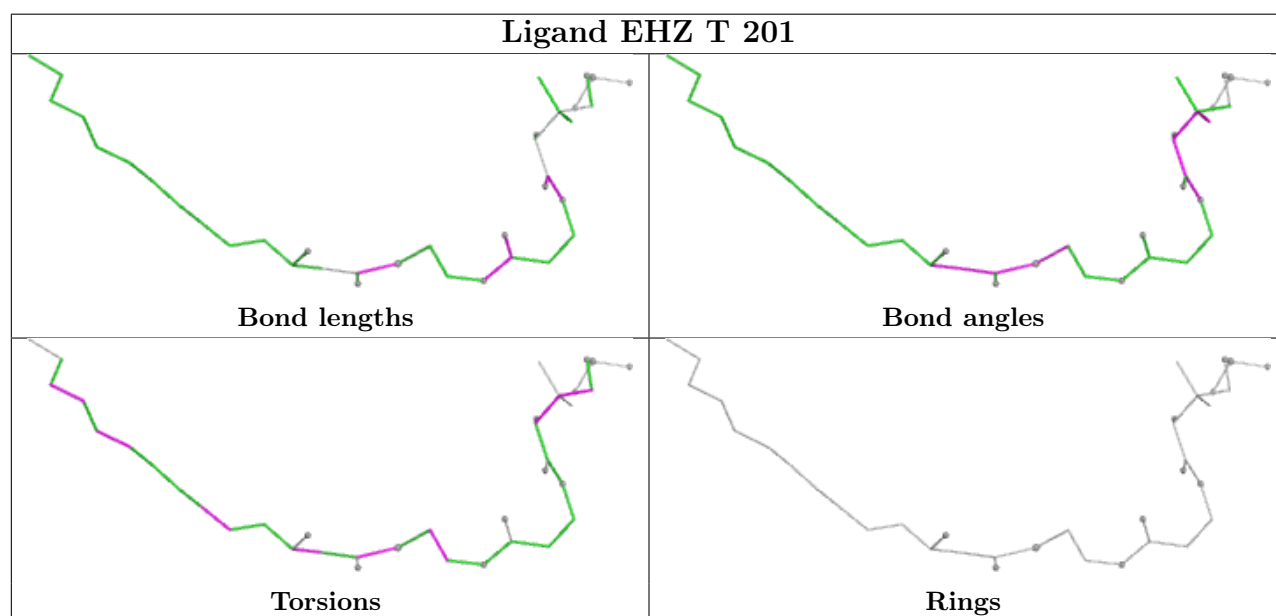
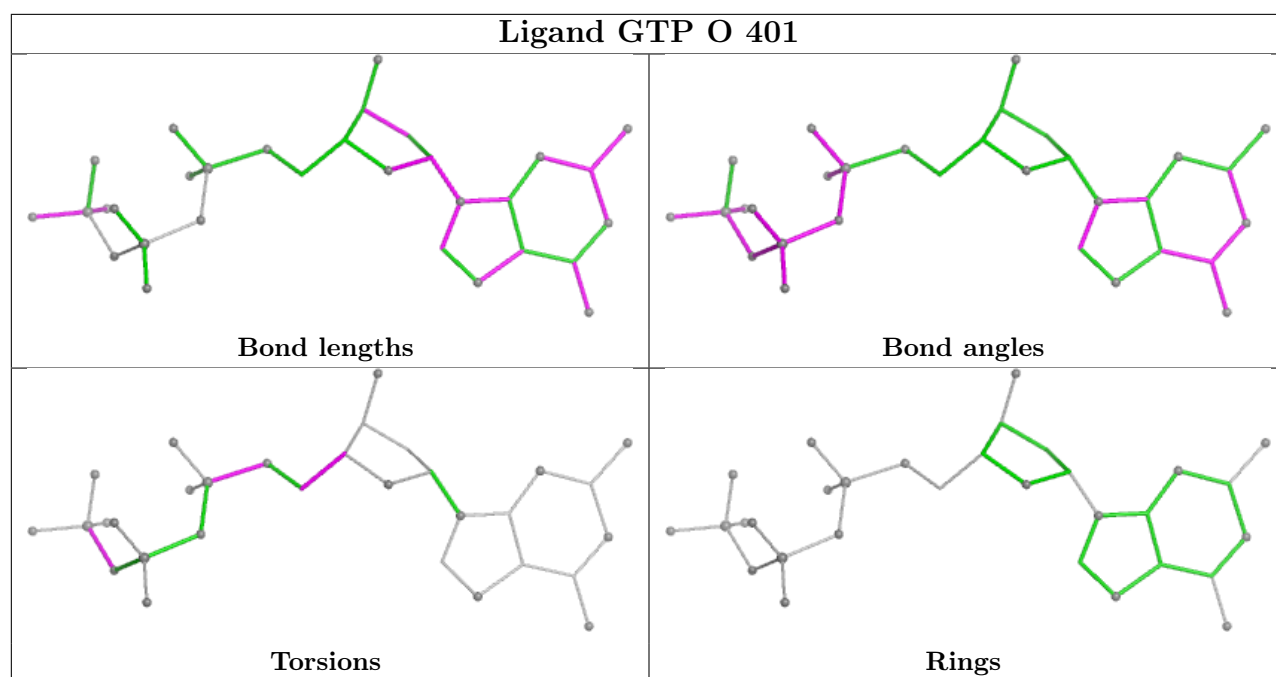


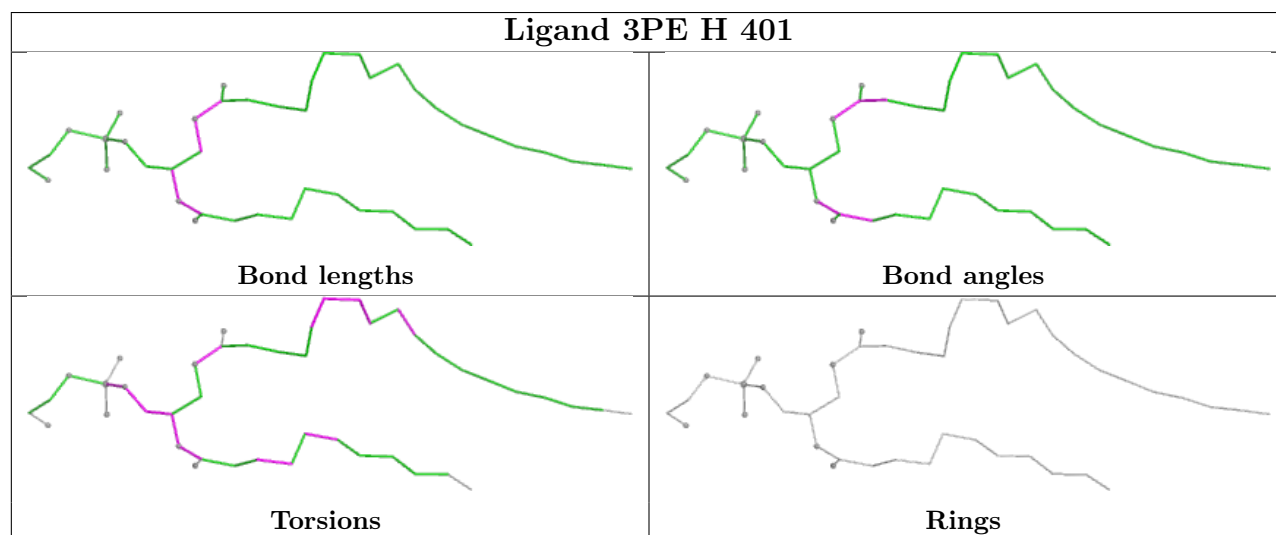
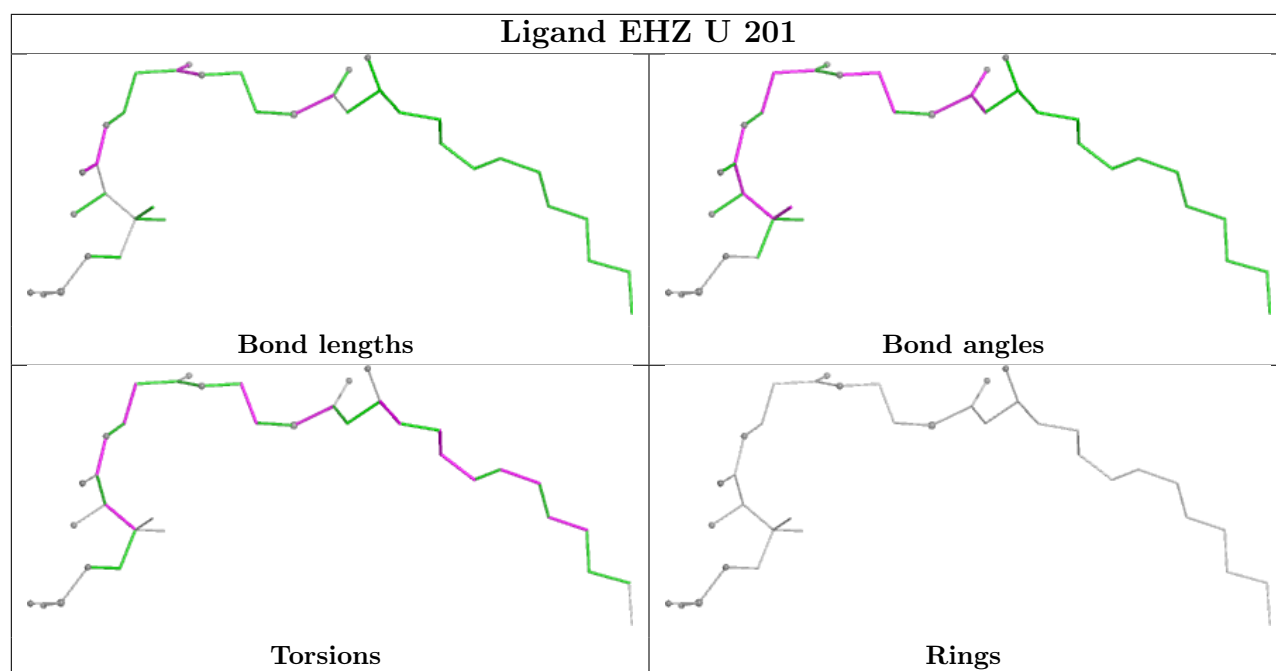


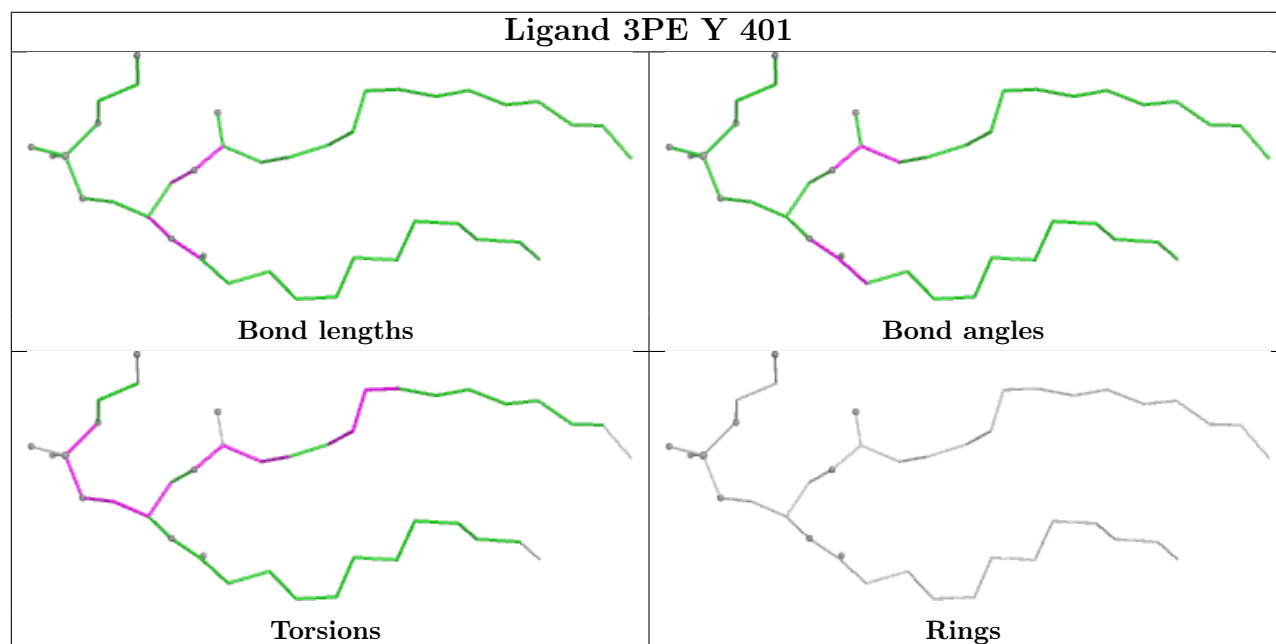
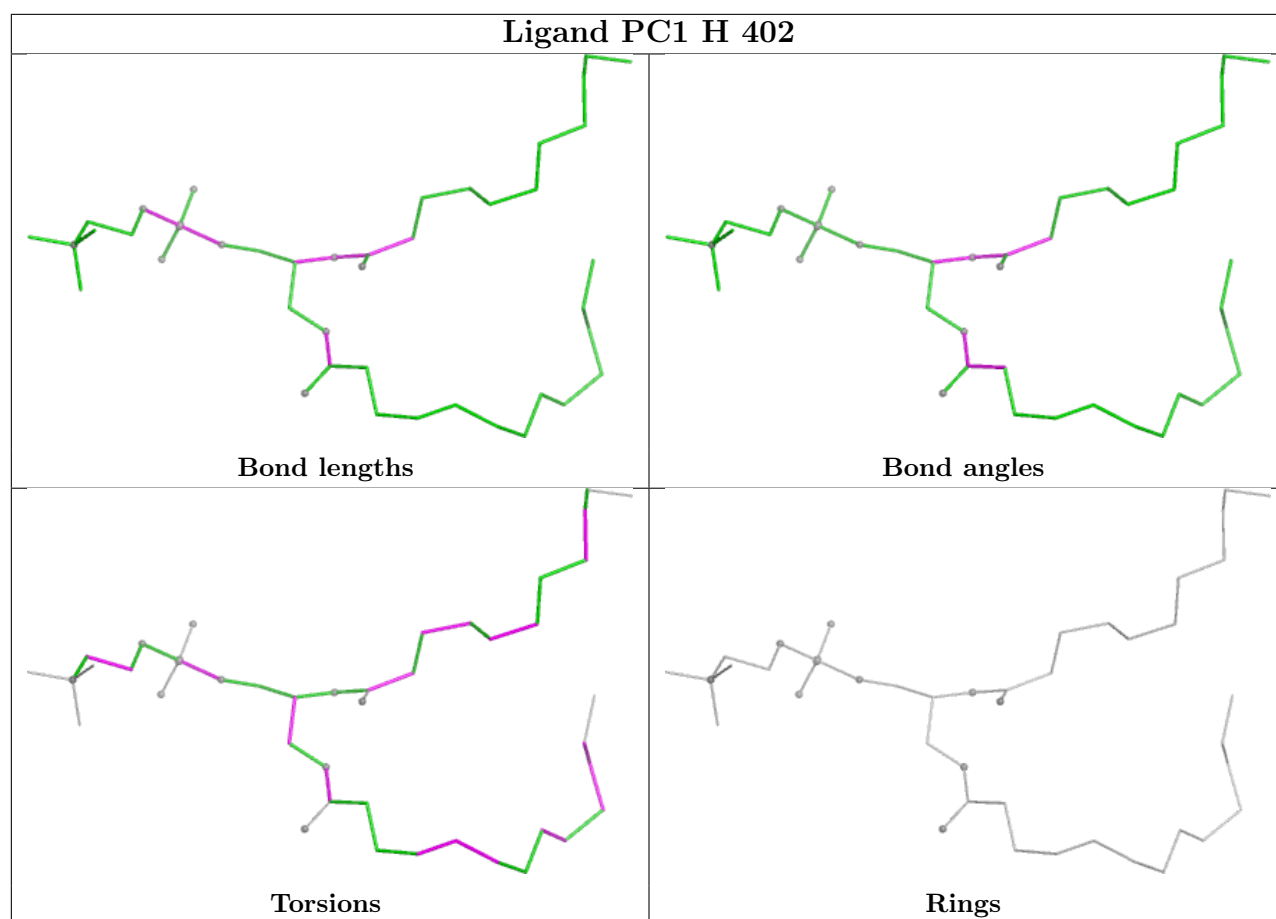


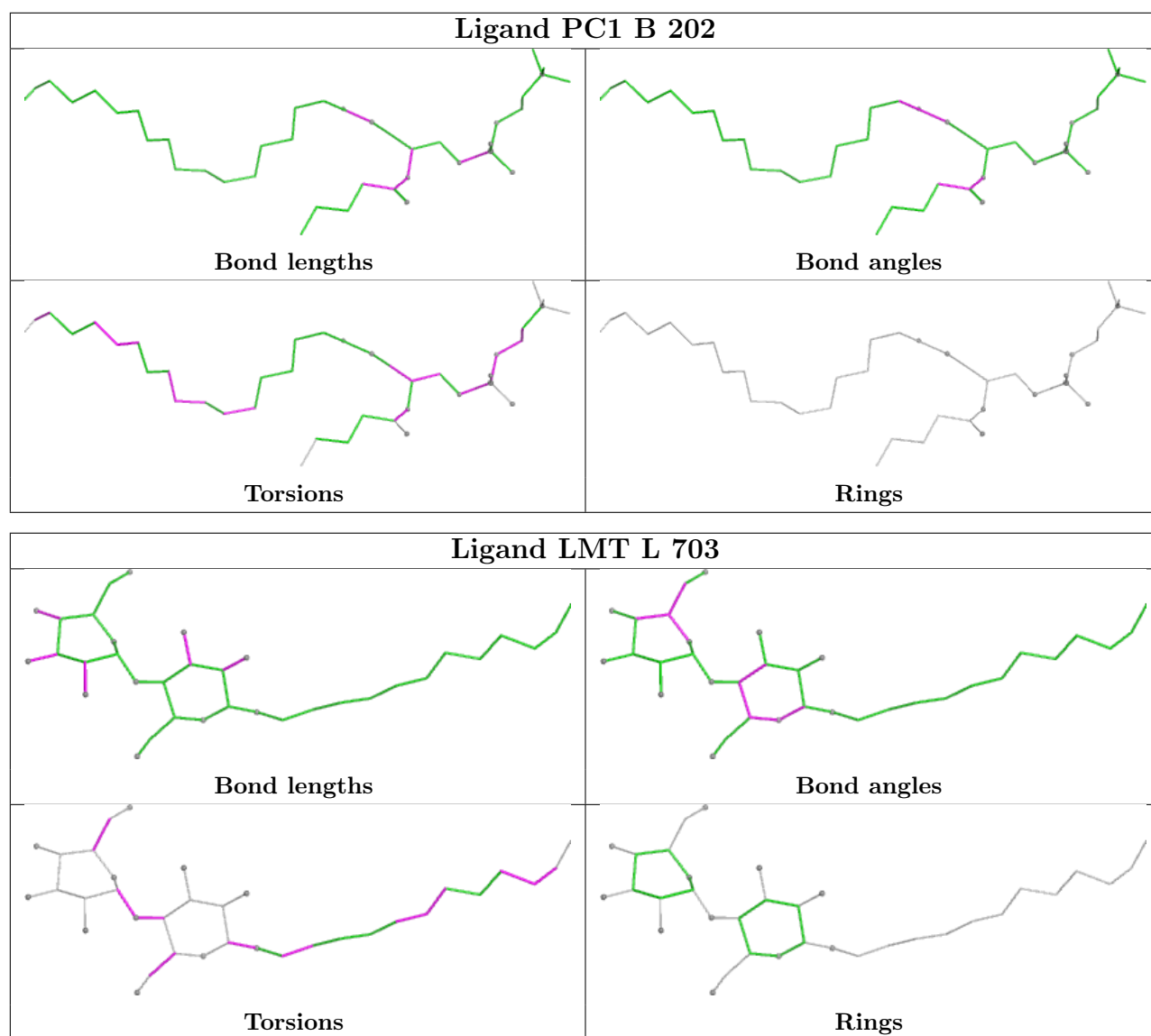


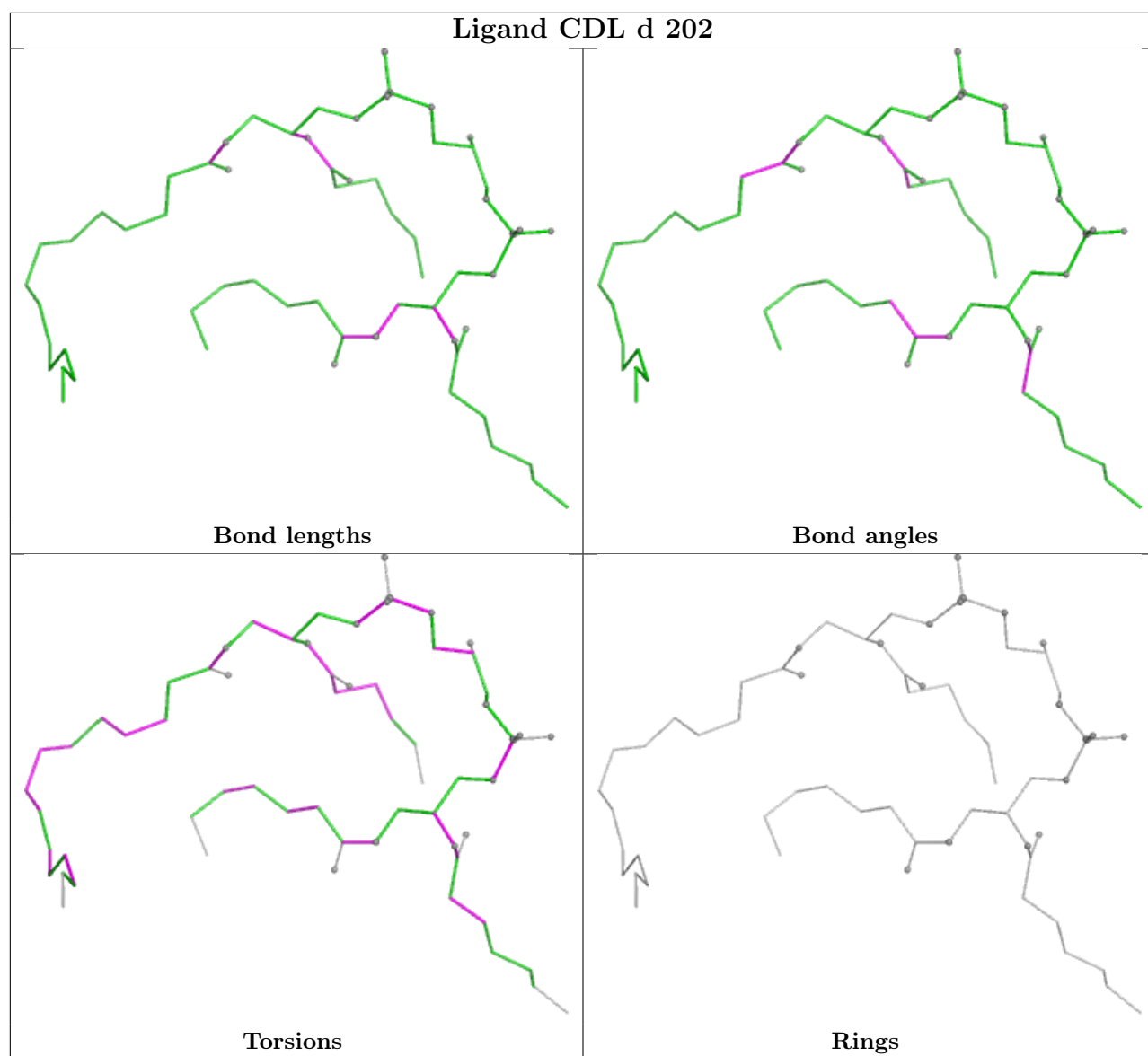


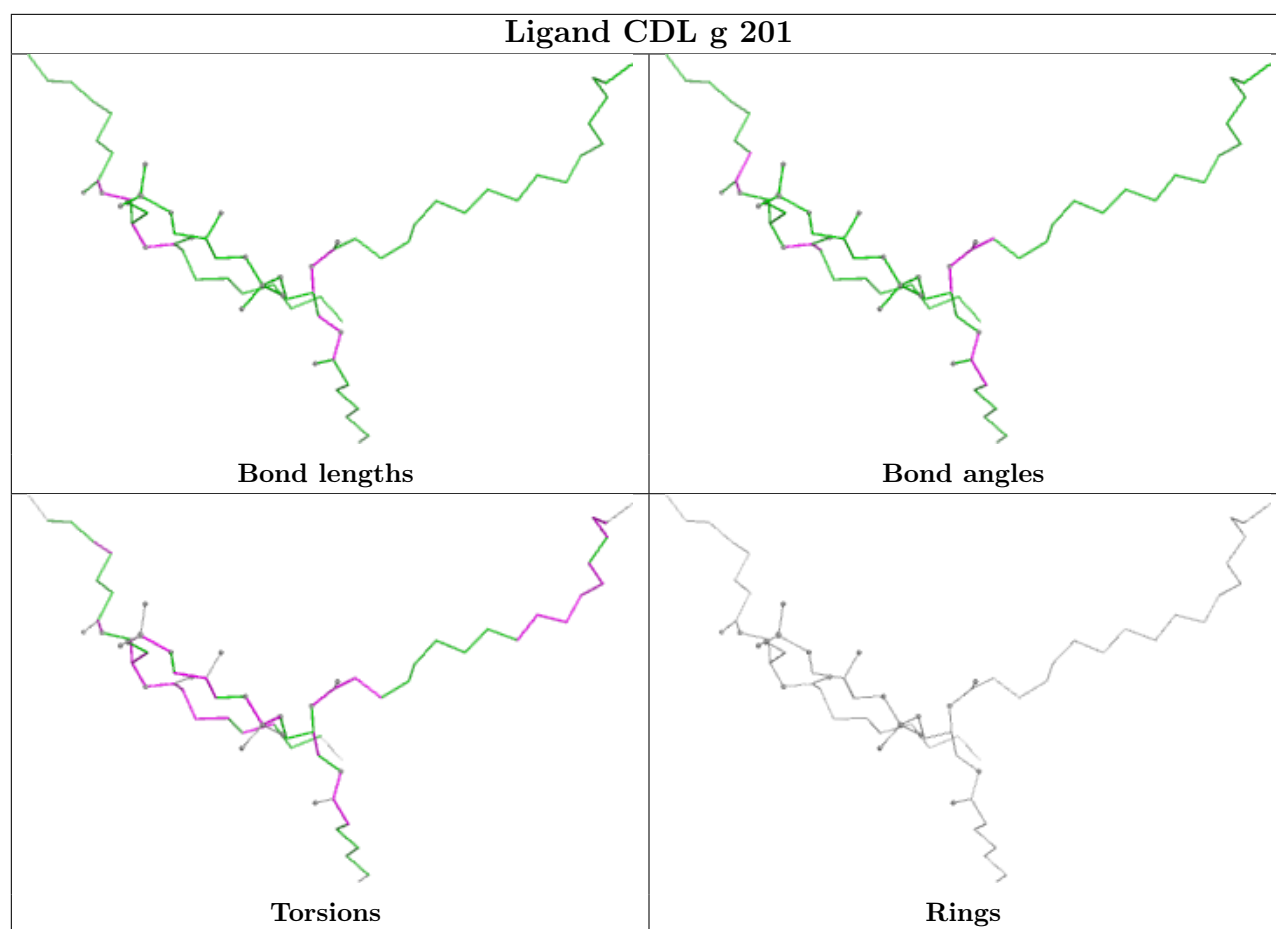


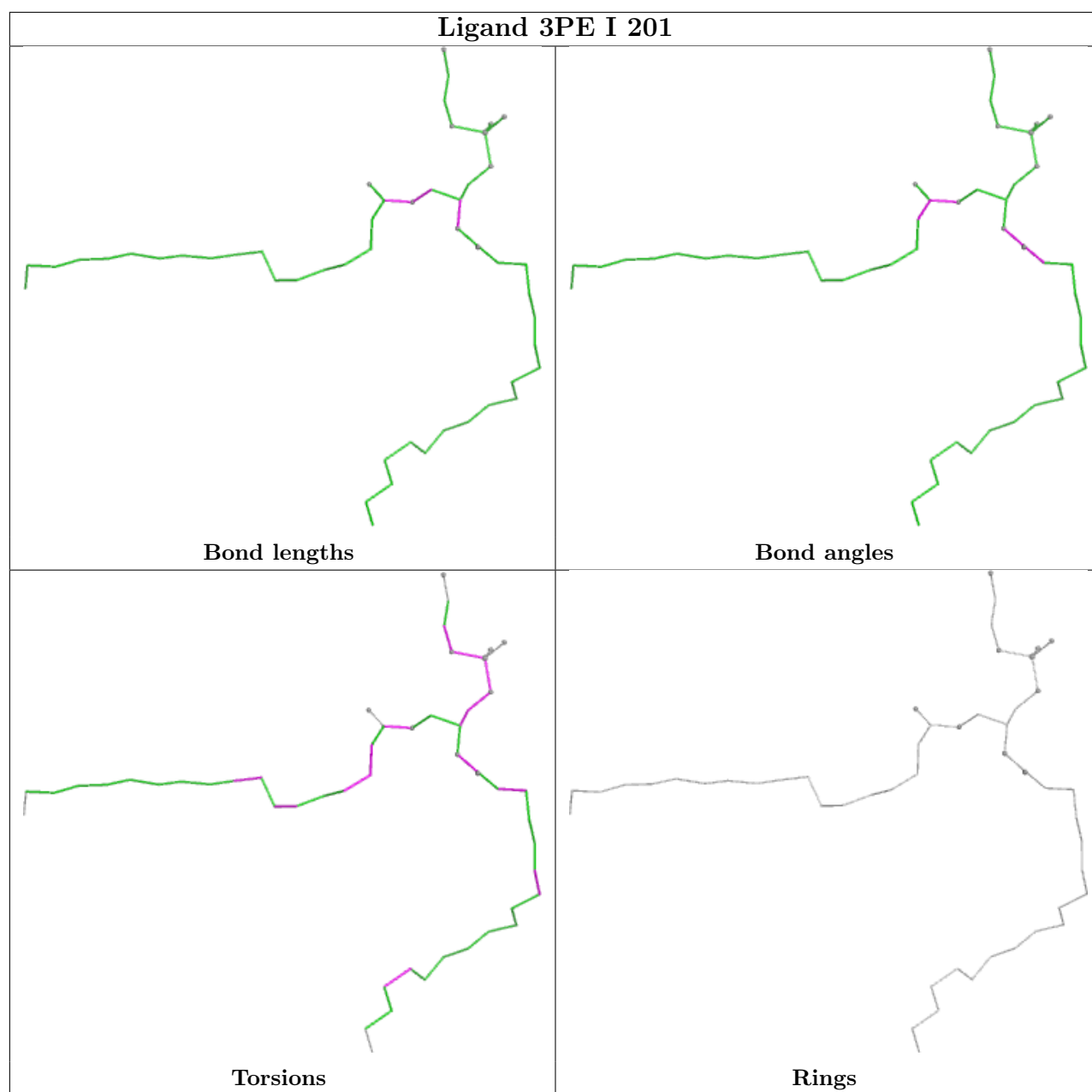


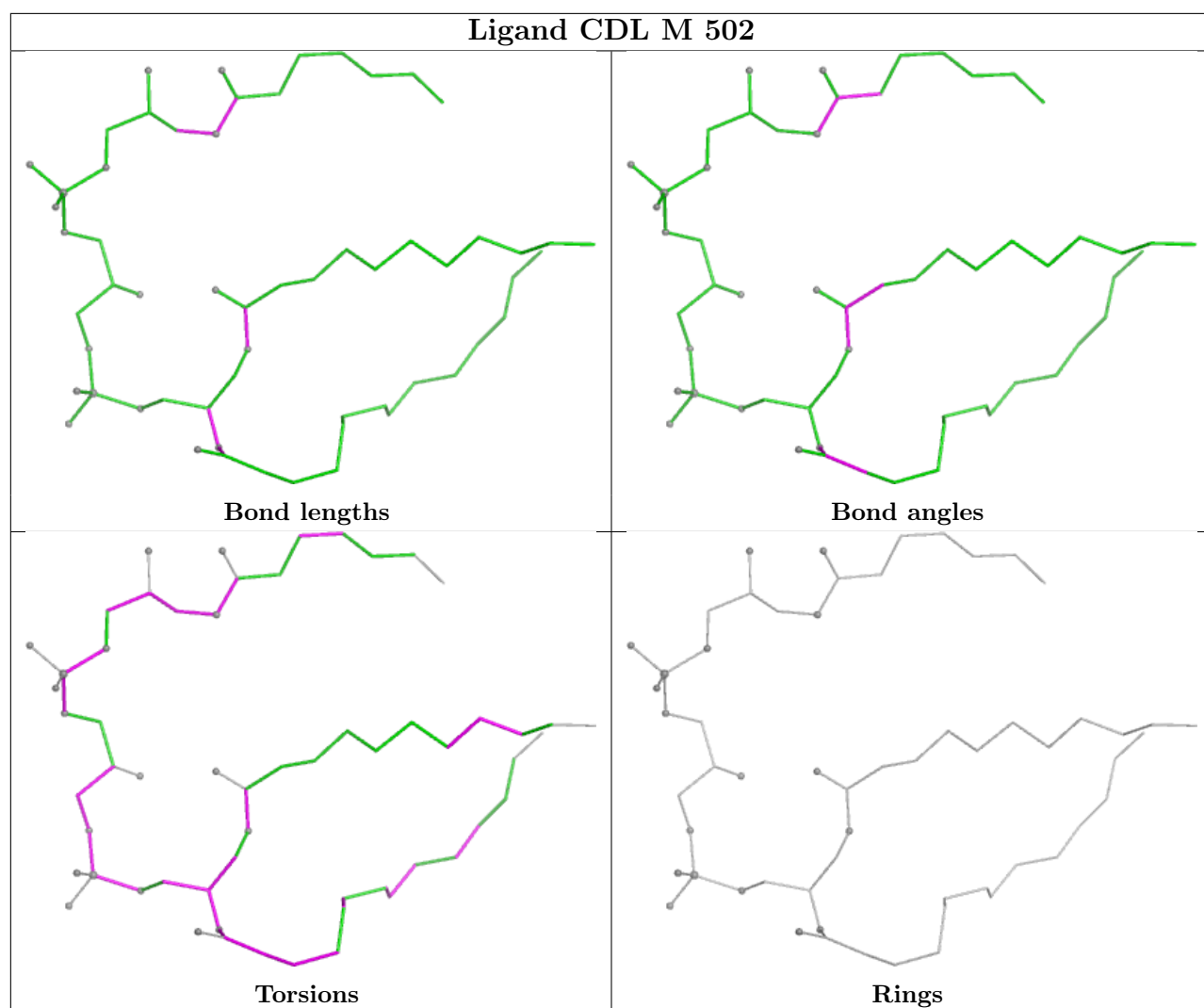


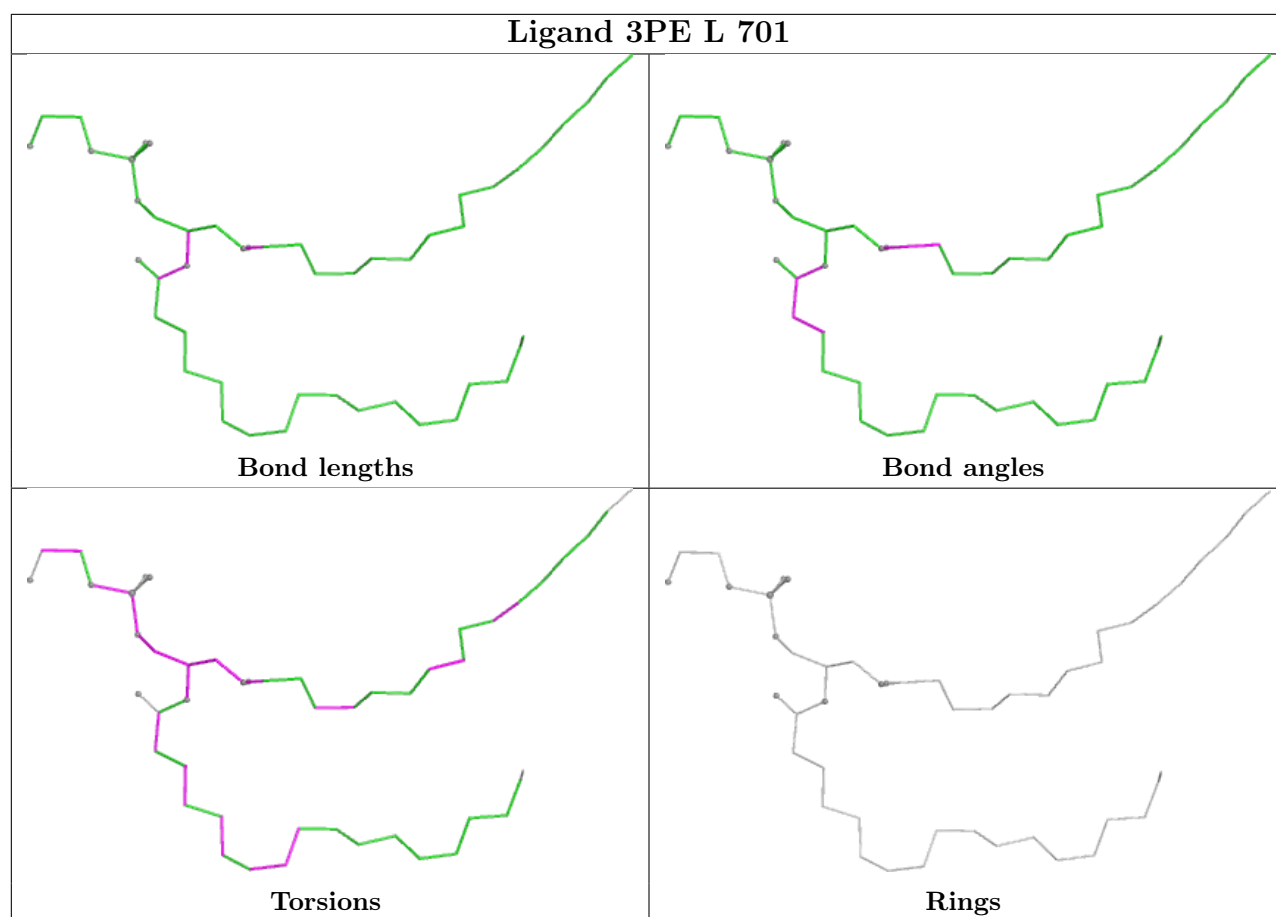


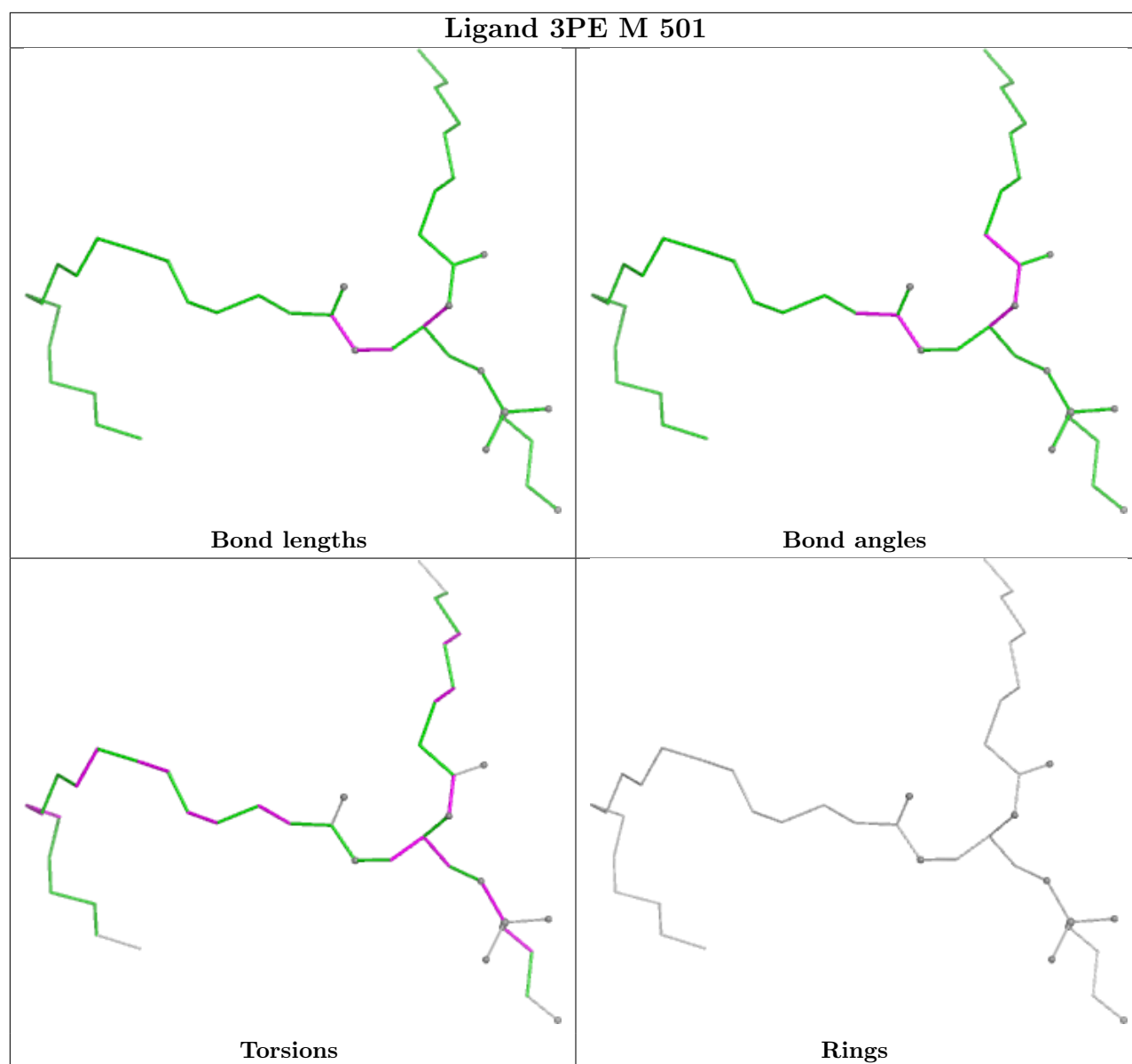












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

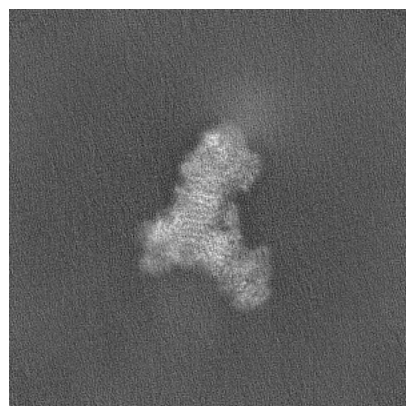
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-16398. These allow visual inspection of the internal detail of the map and identification of artifacts.

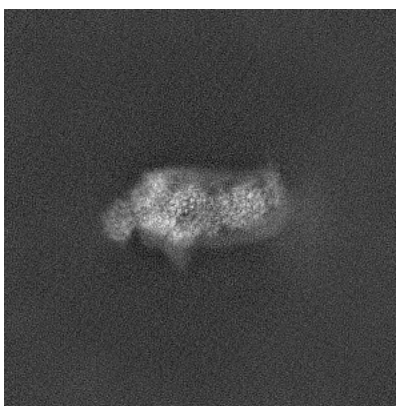
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

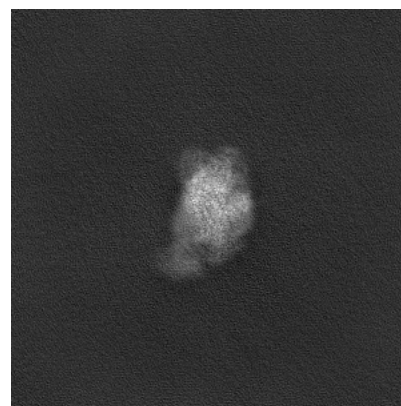
6.1.1 Primary map



X

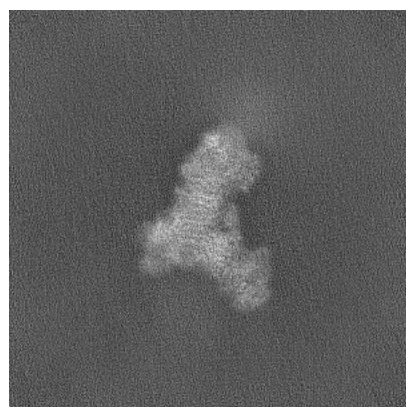


Y

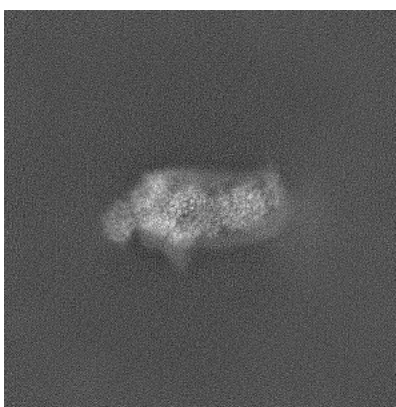


Z

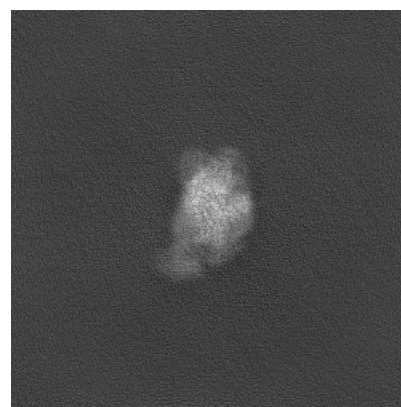
6.1.2 Raw map



X



Y

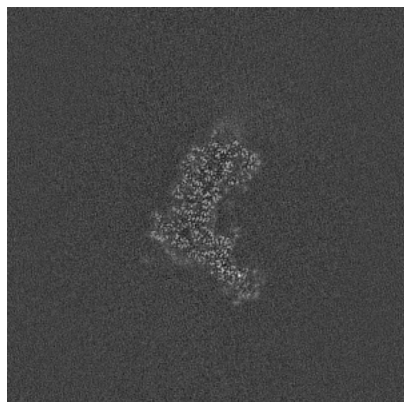


Z

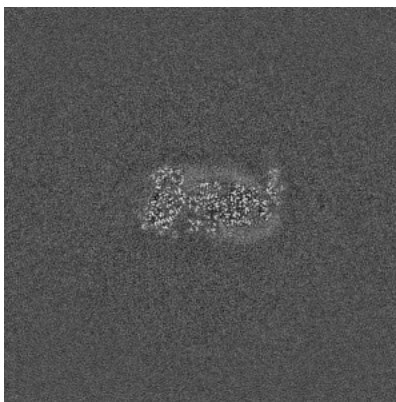
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 225

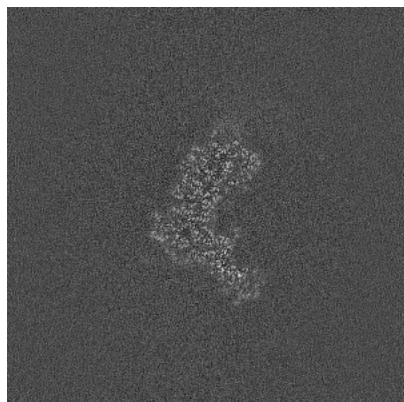


Y Index: 225

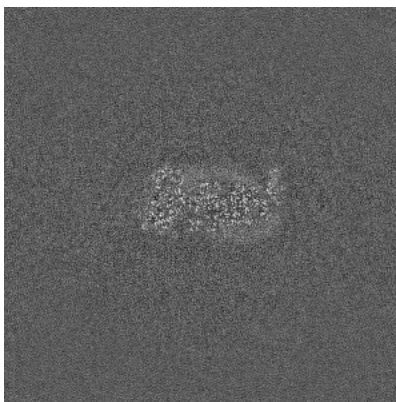


Z Index: 225

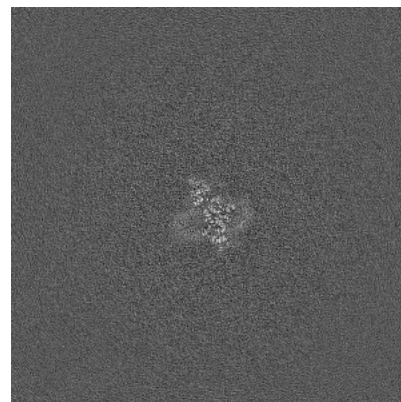
6.2.2 Raw map



X Index: 225



Y Index: 225

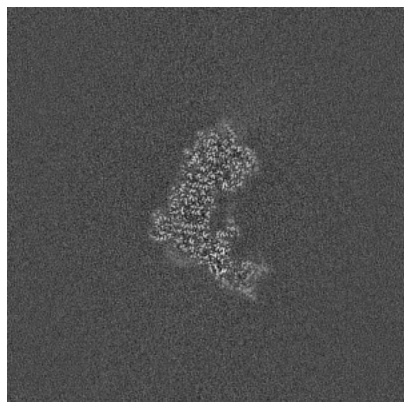


Z Index: 225

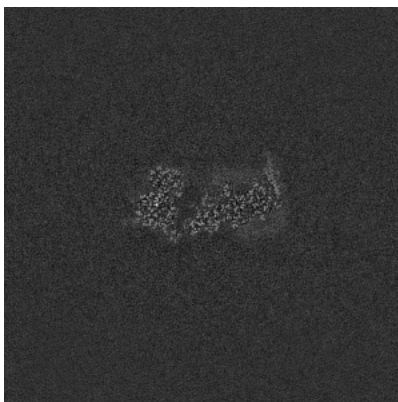
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

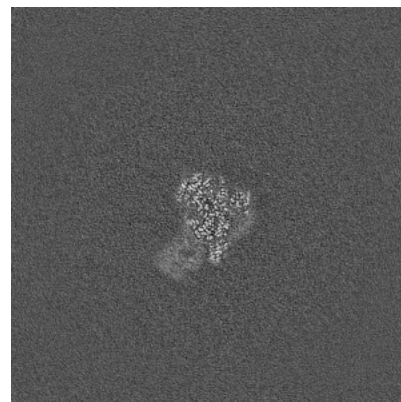
6.3.1 Primary map



X Index: 231

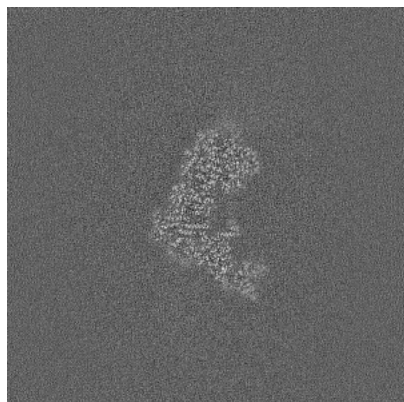


Y Index: 234

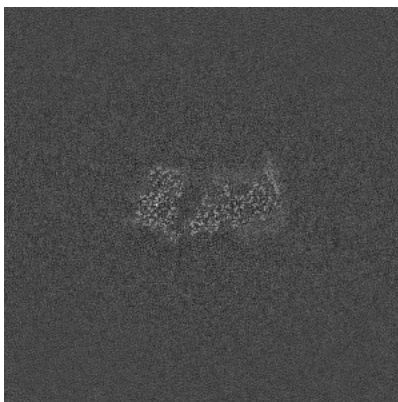


Z Index: 190

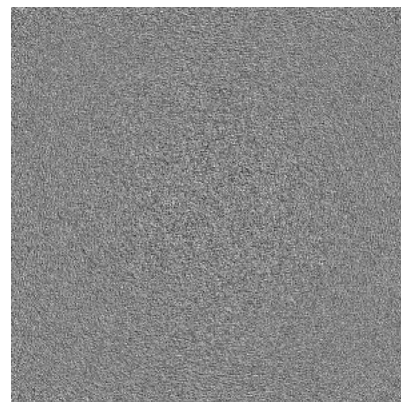
6.3.2 Raw map



X Index: 230



Y Index: 234

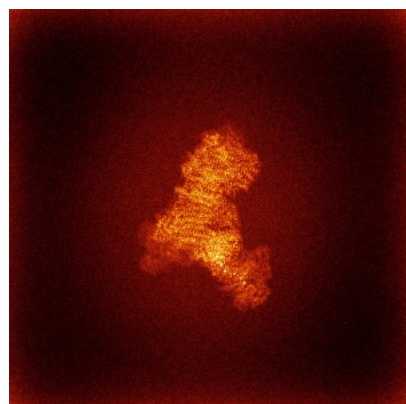


Z Index: 0

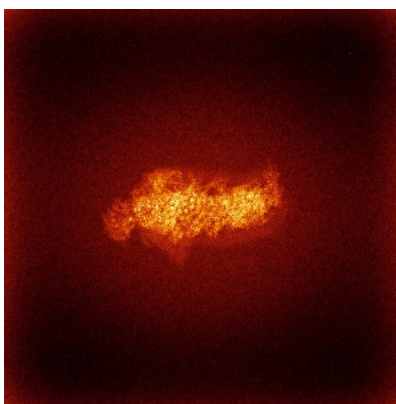
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

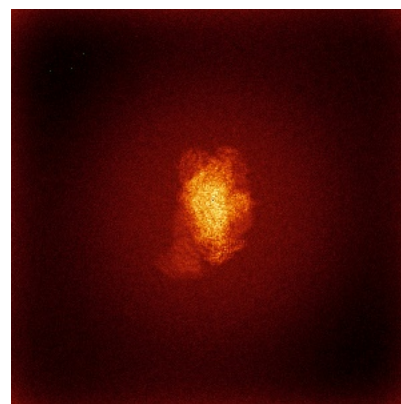
6.4.1 Primary map



X

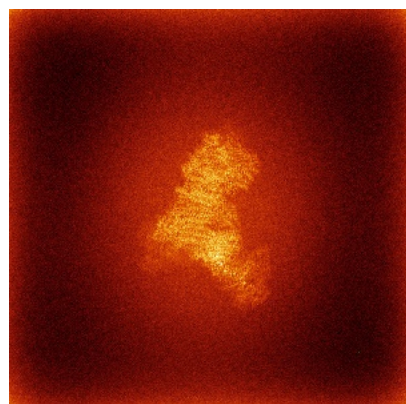


Y

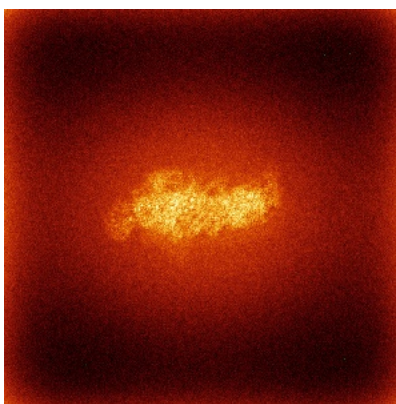


Z

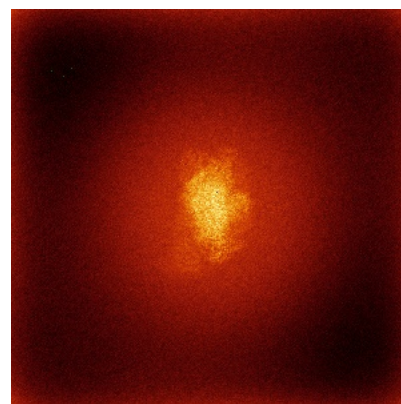
6.4.2 Raw map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



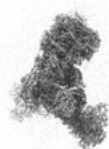
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 1.9. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

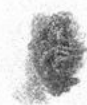
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

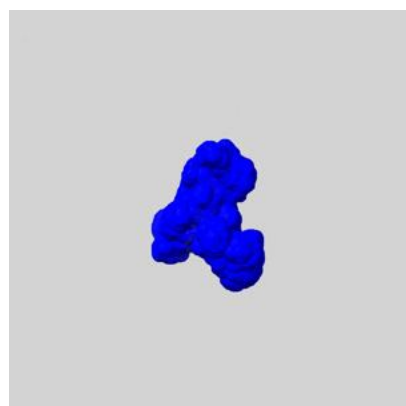
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

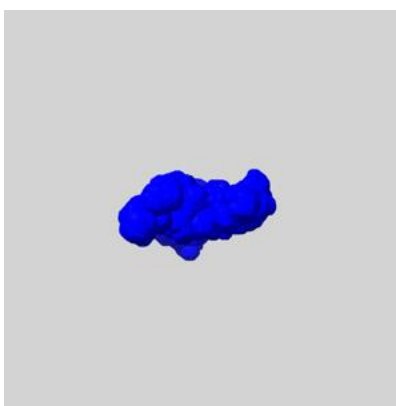
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

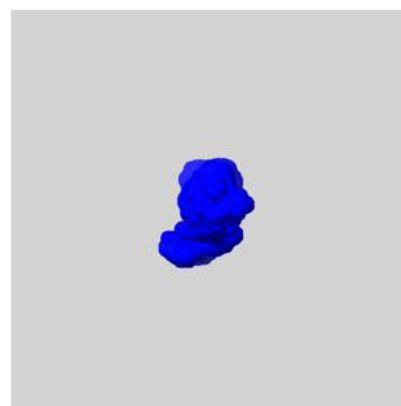
6.6.1 emd_16398_msk_1.map [i](#)



X



Y

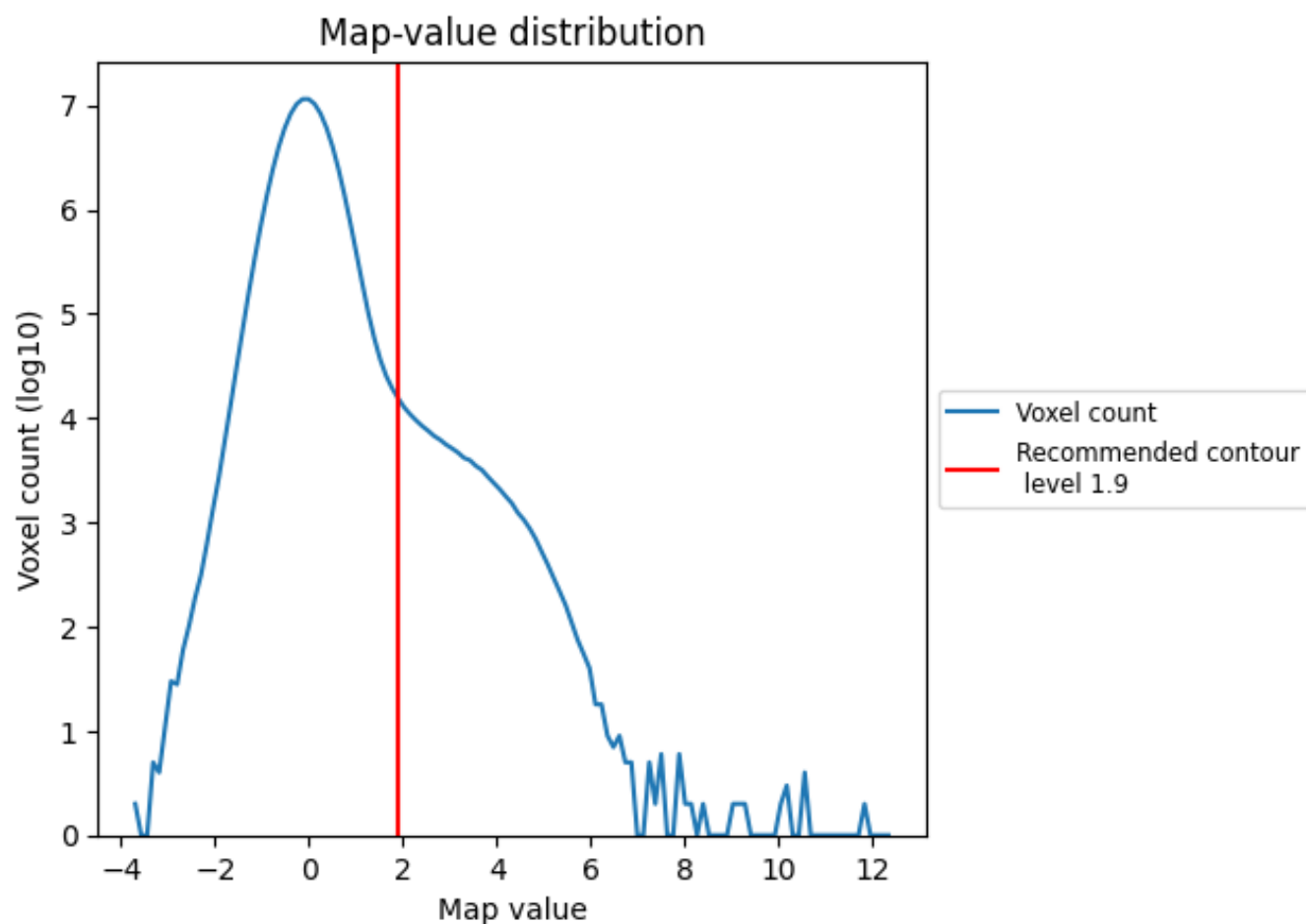


Z

7 Map analysis [i](#)

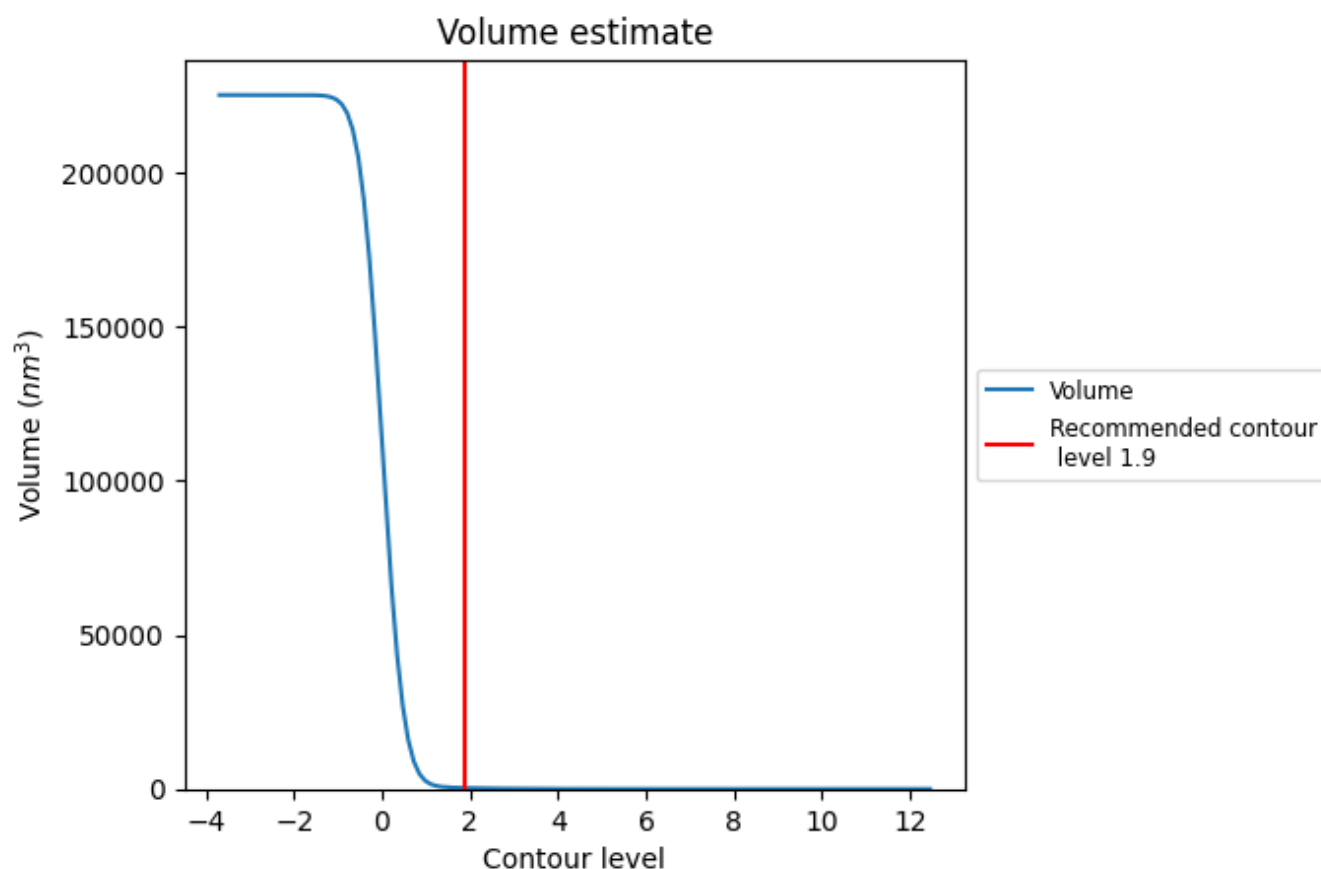
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

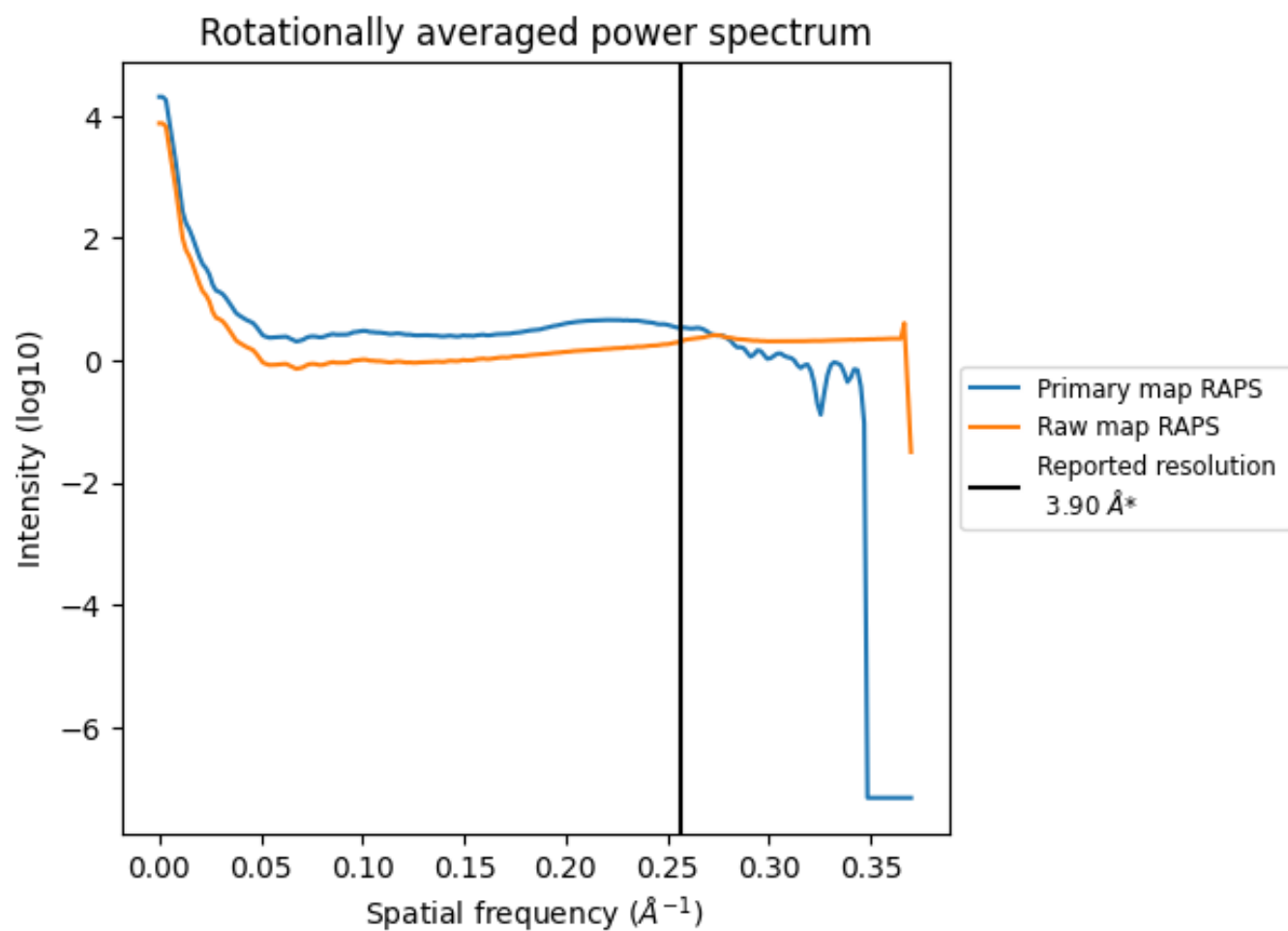
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 309 nm^3 ; this corresponds to an approximate mass of 279 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

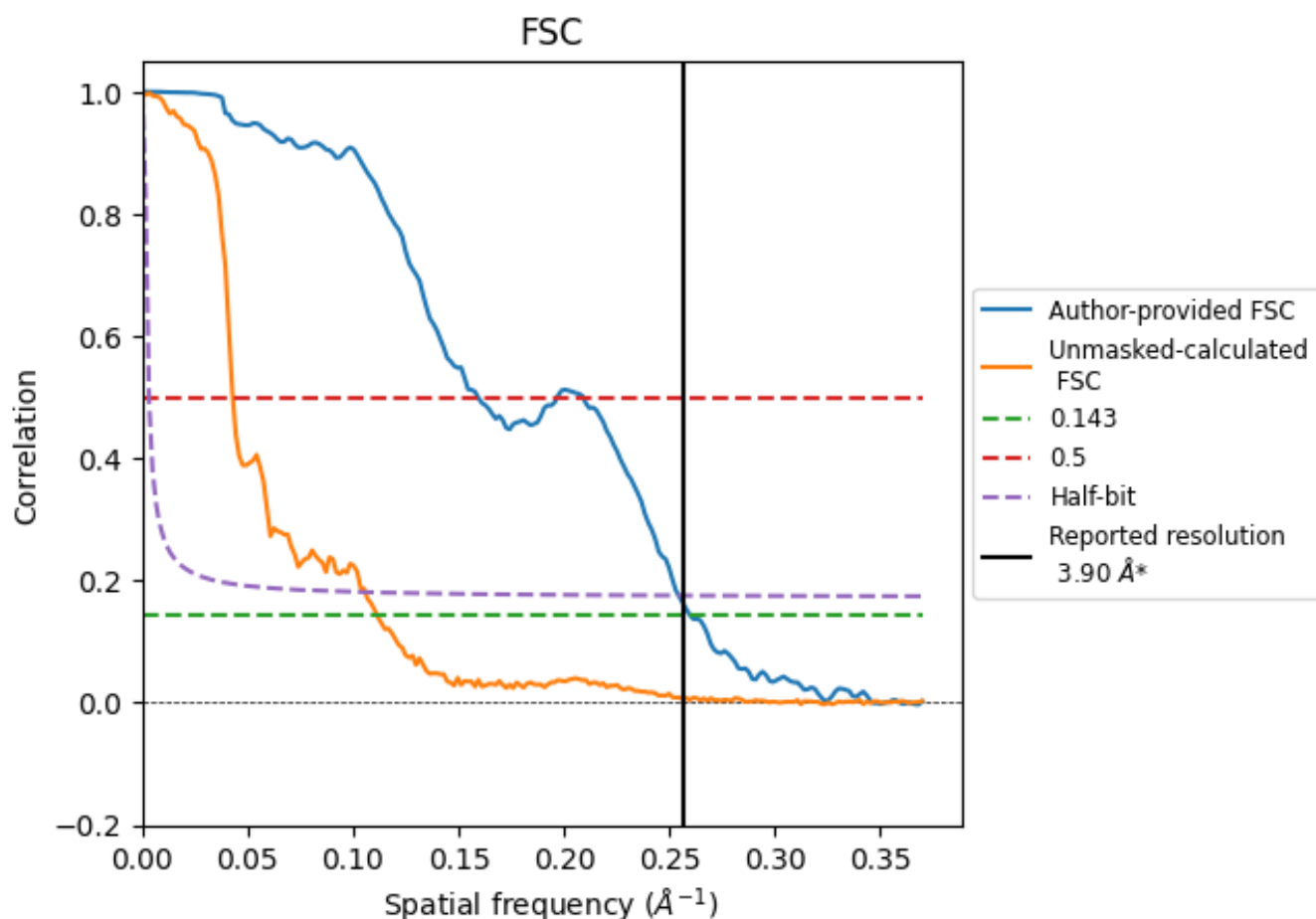


*Reported resolution corresponds to spatial frequency of $0.256 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.256 Å⁻¹

8.2 Resolution estimates [i](#)

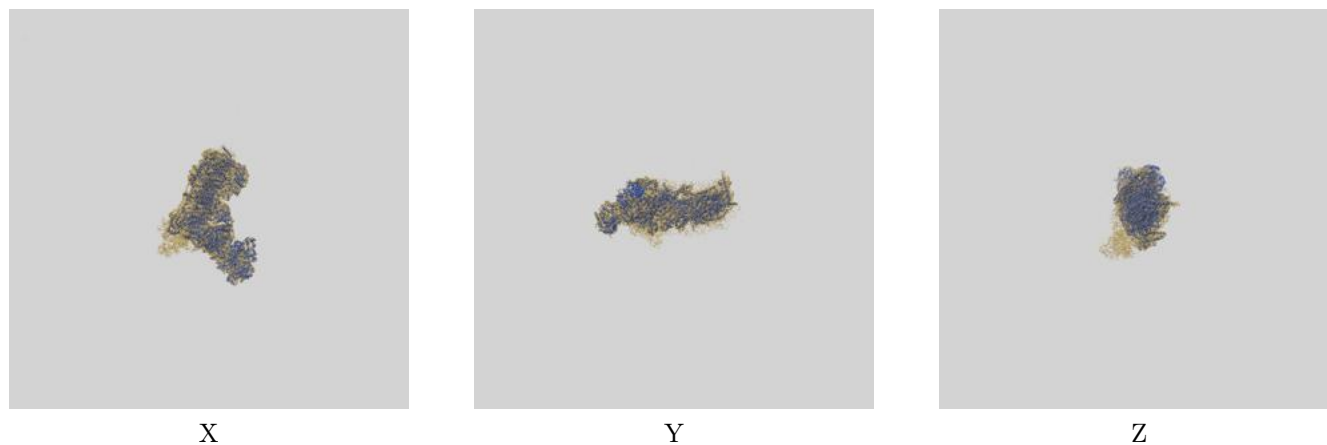
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	3.85	6.28	3.93
Unmasked-calculated*	8.98	23.26	9.45

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 8.98 differs from the reported value 3.9 by more than 10 %

9 Map-model fit [i](#)

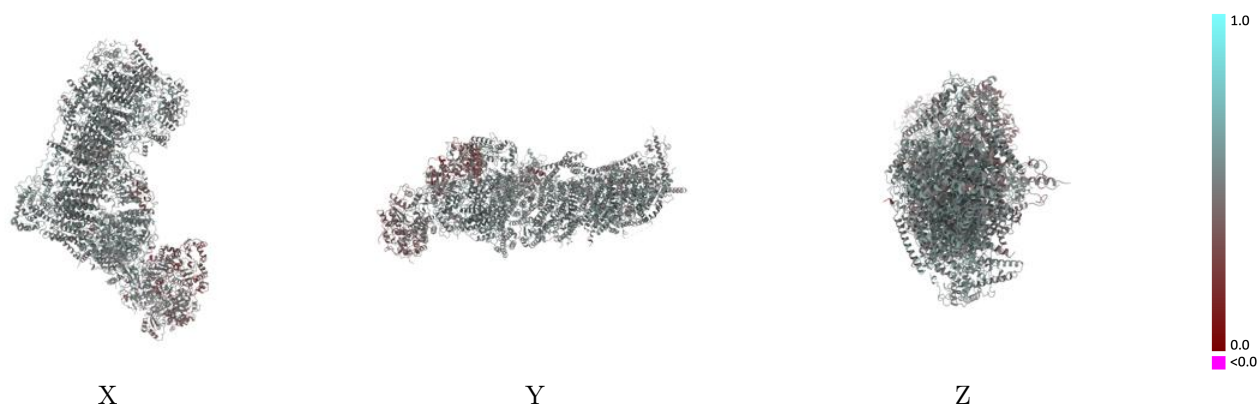
This section contains information regarding the fit between EMDB map EMD-16398 and PDB model 8C2S. Per-residue inclusion information can be found in [section 3](#) on [page 19](#).

9.1 Map-model overlay [i](#)



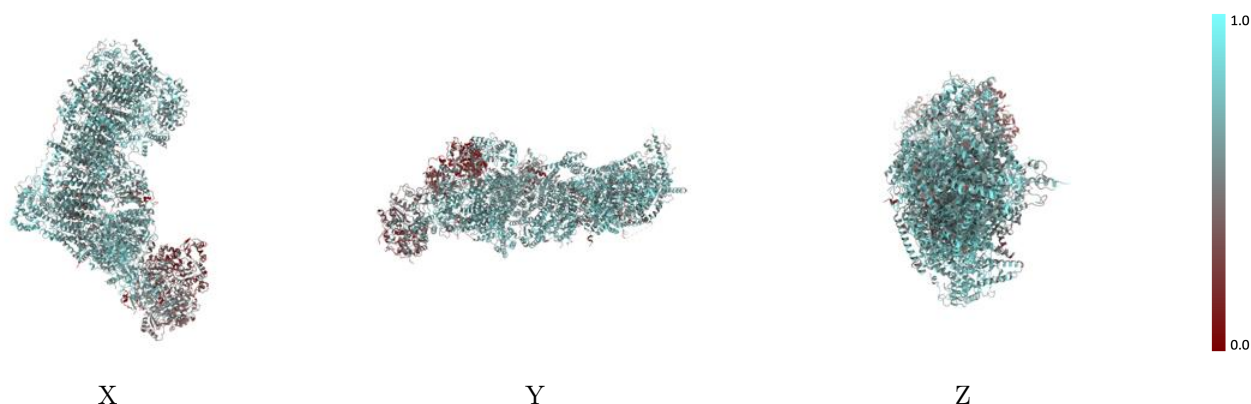
The images above show the 3D surface view of the map at the recommended contour level 1.9 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



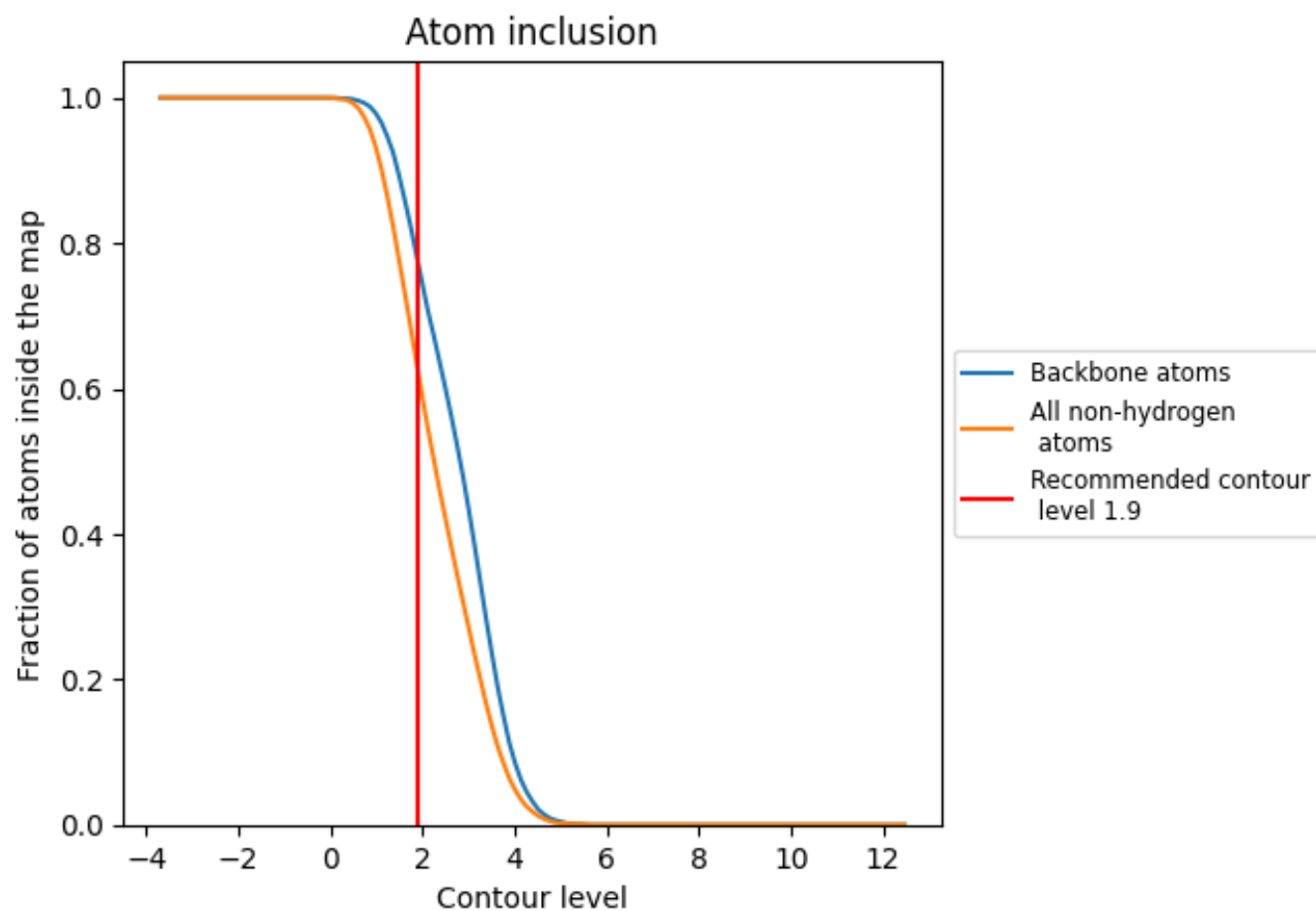
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (1.9).

9.4 Atom inclusion [i](#)



At the recommended contour level, 77% of all backbone atoms, 62% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (1.9) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6230	 0.5000
A	 0.6350	 0.5270
B	 0.7280	 0.5340
C	 0.6770	 0.5370
D	 0.7060	 0.5340
E	 0.4210	 0.4290
F	 0.4510	 0.4320
G	 0.4260	 0.4360
H	 0.6810	 0.5250
I	 0.6890	 0.5130
J	 0.6100	 0.5090
K	 0.6750	 0.5300
L	 0.6520	 0.5170
M	 0.6920	 0.5330
N	 0.6860	 0.5280
O	 0.7070	 0.5200
P	 0.6580	 0.5110
R	 0.4240	 0.4760
S	 0.2750	 0.3480
T	 0.4150	 0.4040
U	 0.6000	 0.4840
V	 0.6780	 0.5160
W	 0.5830	 0.4940
X	 0.6900	 0.5140
Y	 0.6270	 0.4890
Z	 0.6720	 0.5090
a	 0.7110	 0.5220
b	 0.6690	 0.5010
c	 0.6250	 0.4990
d	 0.6360	 0.5200
e	 0.6980	 0.5300
f	 0.5890	 0.4780
g	 0.6280	 0.5170
h	 0.7010	 0.5270
i	 0.6350	 0.5110



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Chain	Atom inclusion	Q-score
j	 0.6380	 0.4900
k	 0.6350	 0.4870
l	 0.7040	 0.5210
m	 0.6950	 0.5120
n	 0.6910	 0.5110
o	 0.6220	 0.4730
p	 0.6820	 0.5060
r	 0.5660	 0.5250
s	 0.2770	 0.3610