



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 12:31 PM EDT

PDB ID : 2IKC / pdb_00002ikc
Title : Crystal structure of sheep lactoperoxidase at 3.25 Å resolution reveals the binding sites for formate
Authors : Sheikh, I.A.; Singh, N.; Singh, A.K.; Sharma, S.; Singh, T.P.
Deposited on : 2006-10-02
Resolution : 3.25 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

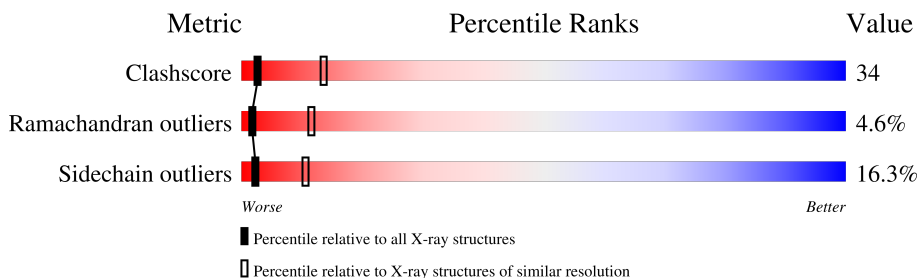
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1660 (3.30-3.22)
Ramachandran outliers	187476	1630 (3.30-3.22)
Sidechain outliers	187428	1629 (3.30-3.22)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	595	39% 43% 14% .
1	B	595	41% 42% 13% .
2	C	2	100%
2	D	2	50% 50%
2	F	2	50% 50%
2	G	2	100%
3	E	3	67% 33%
3	H	3	67% 33%

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
8	FMT	A	3002	-	-	X	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 10008 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 milk lactoperoxidase.

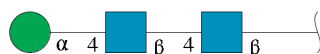
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	S	0	0	0
			4764	3024	853	861	26			
1	B	595	Total	C	N	O	S	0	0	0
			4764	3024	853	861	26			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	3	Total	C	N	O	0	0	0
			39	22	2	15			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	H	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

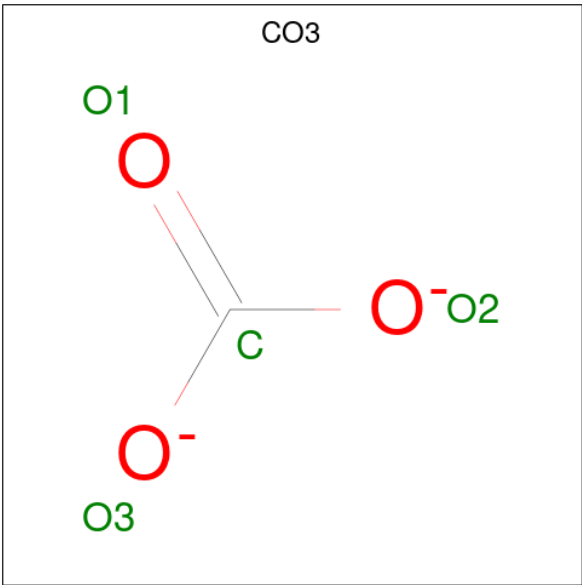


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

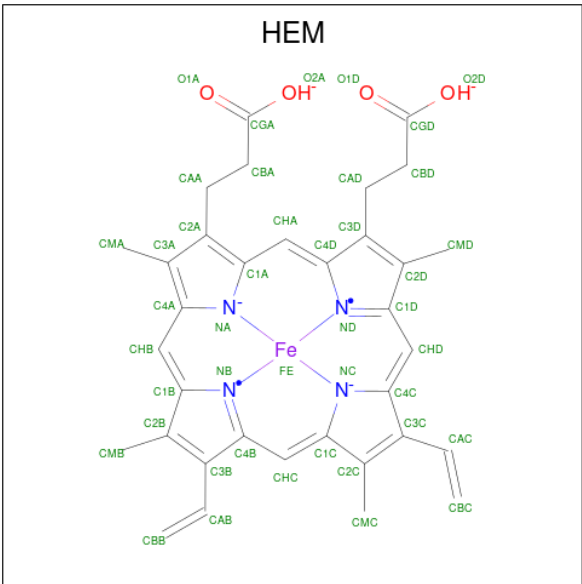
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		
5	B	1	Total	Ca	0	0
			1	1		

- Molecule 6 is CARBONATE ION (CCD ID: CO3) (formula: CO_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	1	3		
6	B	1	Total	C	O	0	0
			4	1	3		

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



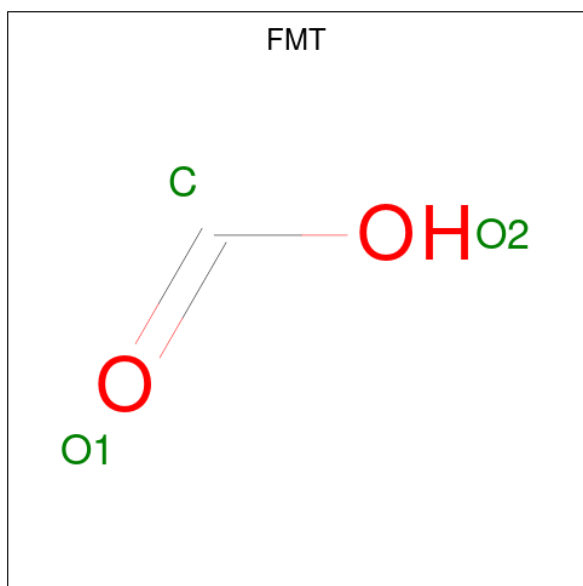
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	Fe	N	O	
			43	34	1	4	4	
							0	0

- Molecule 8 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O		
			3	1	2		
						0	0
8	A	1	Total	C	O		
			3	1	2		
						0	0
8	A	1	Total	C	O		
			3	1	2		
						0	0
8	B	1	Total	C	O		
			3	1	2		
						0	0
8	B	1	Total	C	O		
			3	1	2		
						0	0
8	B	1	Total	C	O		
			3	1	2		
						0	0

- Molecule 9 is water.

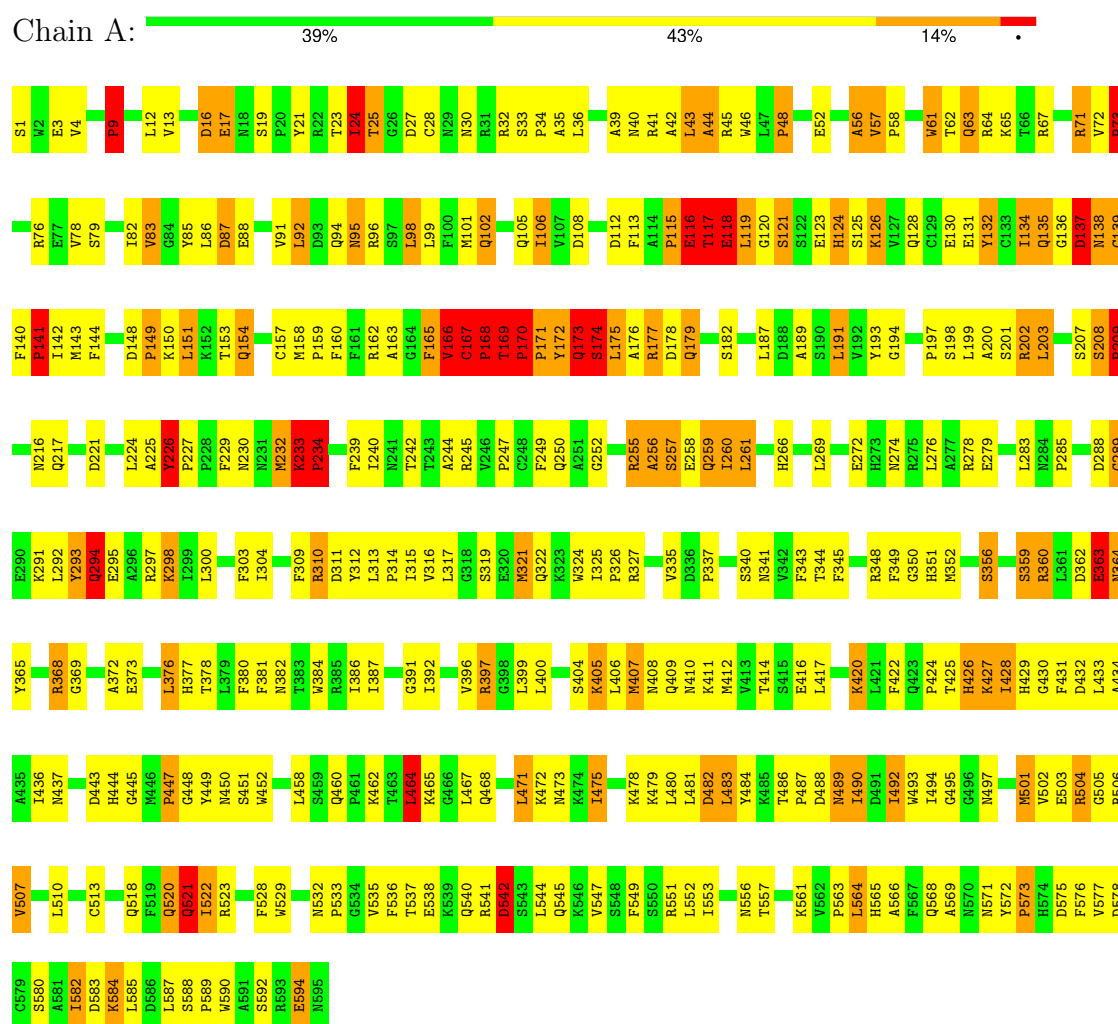
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	86	Total	O		
			86	86		
					0	0
9	B	62	Total	O		
			62	62		
					0	0

3 Residue-property plots

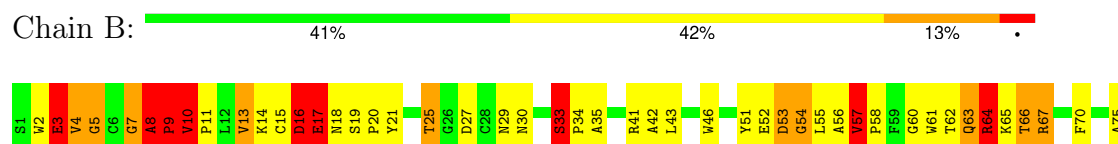
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

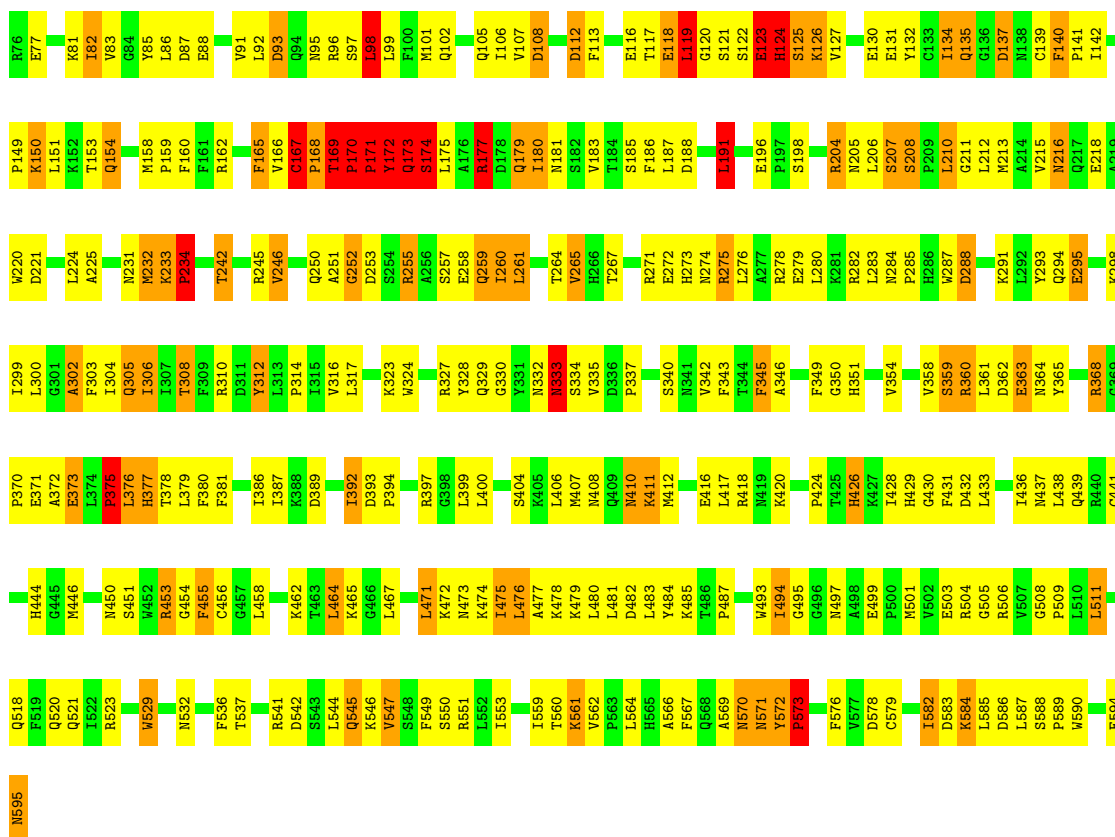
Note EDS was not executed.

• Molecule 1: milk lactoperoxidase



• Molecule 1: milk lactoperoxidase





- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  100%

MAG1
MAG2

- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50%

MAG1
MAG2

- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50%

MAG1
MAG2

- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%

MAG1
MAG2

- Molecule 3: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  67% 33%

MAG1
MAG2
MAN3

- Molecule 3: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  67% 33%

MAG1
MAG2
MAN3

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.08 Å 72.59 Å 84.47 Å 85.20° 84.07° 75.41°	Depositor
Resolution (Å)	19.97 – 3.25	Depositor
% Data completeness (in resolution range)	96.4 (19.97-3.25)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.187 , 0.230	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10008	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, FMT, HEM, MAN, CO3, CA

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	0.91	19/4889 (0.4%)	1.61	132/6629 (2.0%)
1	B	0.91	20/4889 (0.4%)	1.65	129/6629 (1.9%)
All	All	0.91	39/9778 (0.4%)	1.63	261/13258 (2.0%)

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
1	B	0	1

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	67	ARG	N-CA	30.59	1.83	1.46
1	A	24	ILE	N-CA	10.70	1.59	1.46
1	A	172	TYR	N-CA	9.71	1.58	1.46
1	B	122	SER	CA-C	9.40	1.65	1.52
1	A	118	GLU	N-CA	-8.85	1.35	1.45
1	A	9	PRO	N-CA	8.75	1.58	1.47
1	B	123	GLU	N-CA	8.68	1.57	1.46
1	A	173	GLN	N-CA	8.66	1.57	1.46
1	A	173	GLN	CA-C	8.26	1.63	1.52
1	B	330	GLY	N-CA	7.54	1.53	1.45
1	A	171	PRO	CA-C	7.49	1.63	1.52
1	A	172	TYR	C-N	6.84	1.43	1.33
1	A	174	SER	CA-C	6.84	1.61	1.52
1	A	44	ALA	N-CA	6.65	1.54	1.46
1	A	172	TYR	CA-C	6.64	1.61	1.52
1	B	172	TYR	N-CA	6.50	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	573	PRO	N-CA	6.43	1.56	1.47
1	B	572	TYR	C-N	6.30	1.43	1.33
1	B	330	GLY	C-N	6.11	1.42	1.33
1	B	330	GLY	CA-C	6.09	1.58	1.52
1	A	209	PRO	CA-C	-6.08	1.42	1.52
1	B	108	ASP	N-CA	5.85	1.53	1.46
1	B	124	HIS	CB-CG	5.78	1.58	1.50
1	B	572	TYR	N-CA	5.77	1.54	1.46
1	A	171	PRO	N-CA	5.75	1.54	1.47
1	B	9	PRO	CA-C	5.74	1.60	1.52
1	B	121	SER	CA-C	5.70	1.60	1.53
1	B	118	GLU	N-CA	-5.68	1.39	1.46
1	B	107	VAL	C-N	5.67	1.41	1.33
1	A	117	THR	C-N	-5.66	1.25	1.33
1	B	171	PRO	N-CA	5.46	1.54	1.47
1	B	122	SER	C-N	5.46	1.41	1.33
1	A	141	PRO	CA-C	-5.38	1.46	1.52
1	B	118	GLU	CA-C	-5.25	1.47	1.53
1	A	169	THR	CA-C	5.17	1.59	1.52
1	B	117	THR	C-N	-5.12	1.26	1.33
1	A	112	ASP	C-N	5.12	1.40	1.33
1	A	112	ASP	CA-CB	5.08	1.63	1.53
1	A	173	GLN	C-N	5.01	1.40	1.33

All (261) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	233	LYS	CA-C-N	36.67	165.67	119.84
1	B	233	LYS	C-N-CA	36.67	165.67	119.84
1	B	67	ARG	N-CA-CB	-19.57	79.85	110.57
1	B	66	THR	CA-C-N	-17.38	99.01	123.00
1	B	66	THR	C-N-CA	-17.38	99.01	123.00
1	A	24	ILE	N-CA-C	15.59	130.54	111.05
1	B	573	PRO	CA-N-CD	-14.21	92.10	112.00
1	A	233	LYS	CA-C-N	13.76	137.04	119.84
1	A	233	LYS	C-N-CA	13.76	137.04	119.84
1	A	482	ASP	N-CA-C	-12.89	97.54	113.01
1	B	119	LEU	N-CA-C	-12.54	91.91	110.46
1	B	124	HIS	CA-CB-CG	12.47	126.27	113.80
1	A	112	ASP	CA-CB-CG	12.24	124.84	112.60
1	A	171	PRO	N-CA-C	12.19	137.59	112.47
1	B	482	ASP	N-CA-C	-11.90	98.73	113.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	233	LYS	C-N-CD	-11.43	78.12	125.00
1	B	573	PRO	N-CA-C	11.18	130.41	113.75
1	A	105	GLN	OE1-CD-NE2	-10.76	111.84	122.60
1	B	333	ASN	CA-CB-CG	10.68	123.28	112.60
1	A	172	TYR	CA-CB-CG	10.68	133.12	113.90
1	A	41	ARG	CA-C-N	-10.61	107.58	123.07
1	A	41	ARG	C-N-CA	-10.61	107.58	123.07
1	A	166	VAL	N-CA-C	10.31	130.79	109.34
1	A	460	GLN	OE1-CD-NE2	-10.31	112.29	122.60
1	A	179	GLN	OE1-CD-NE2	-10.00	112.60	122.60
1	A	173	GLN	OE1-CD-NE2	-9.92	112.68	122.60
1	B	173	GLN	OE1-CD-NE2	-9.91	112.69	122.60
1	B	154	GLN	OE1-CD-NE2	-9.90	112.70	122.60
1	B	250	GLN	OE1-CD-NE2	-9.90	112.70	122.60
1	A	154	GLN	OE1-CD-NE2	-9.88	112.72	122.60
1	A	294	GLN	OE1-CD-NE2	-9.85	112.75	122.60
1	A	173	GLN	N-CA-C	9.84	131.75	110.80
1	A	135	GLN	OE1-CD-NE2	-9.78	112.82	122.60
1	B	135	GLN	OE1-CD-NE2	-9.78	112.82	122.60
1	B	305	GLN	OE1-CD-NE2	-9.61	112.99	122.60
1	A	94	GLN	OE1-CD-NE2	-9.54	113.06	122.60
1	A	63	GLN	OE1-CD-NE2	-9.44	113.16	122.60
1	A	16	ASP	N-CA-C	-9.42	100.70	110.97
1	A	259	GLN	OE1-CD-NE2	-9.31	113.29	122.60
1	A	102	GLN	OE1-CD-NE2	-9.26	113.34	122.60
1	A	171	PRO	O-C-N	-9.13	110.31	122.64
1	A	295	GLU	N-CA-C	-9.11	100.96	112.90
1	B	174	SER	N-CA-C	9.07	130.12	110.80
1	B	259	GLN	OE1-CD-NE2	-9.07	113.53	122.60
1	B	171	PRO	CA-C-N	9.04	138.81	121.54
1	B	171	PRO	C-N-CA	9.04	138.81	121.54
1	A	233	LYS	C-N-CD	-9.00	88.09	125.00
1	A	128	GLN	OE1-CD-NE2	-8.87	113.73	122.60
1	A	56	ALA	CA-C-N	-8.82	116.24	122.59
1	A	56	ALA	C-N-CA	-8.82	116.24	122.59
1	B	482	ASP	CA-CB-CG	8.78	121.38	112.60
1	A	209	PRO	CA-N-CD	-8.77	99.73	112.00
1	B	571	ASN	CB-CA-C	8.71	133.01	113.33
1	A	115	PRO	CA-C-N	-8.60	105.33	120.97
1	A	115	PRO	C-N-CA	-8.60	105.33	120.97
1	B	234	PRO	CA-N-CD	-8.57	100.00	112.00
1	A	172	TYR	CB-CA-C	-8.38	93.74	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	204	ARG	N-CA-C	8.27	122.39	109.07
1	B	122	SER	N-CA-C	-8.19	93.36	110.80
1	B	112	ASP	N-CA-C	8.14	121.44	109.07
1	A	501	MET	N-CA-C	7.97	120.60	110.24
1	B	122	SER	CA-C-N	7.91	136.65	121.54
1	B	122	SER	C-N-CA	7.91	136.65	121.54
1	A	369	GLY	CA-C-N	7.89	127.53	119.56
1	A	369	GLY	C-N-CA	7.89	127.53	119.56
1	B	316	VAL	N-CA-C	-7.84	104.11	111.48
1	A	521	GLN	N-CA-C	7.78	119.39	111.07
1	A	397	ARG	N-CA-C	-7.73	103.30	112.89
1	B	584	LYS	N-CA-C	7.71	124.83	114.12
1	A	191	LEU	N-CA-C	-7.70	103.34	112.89
1	B	453	ARG	N-CA-C	-7.66	102.90	111.71
1	B	57	VAL	CA-C-N	7.61	127.61	119.78
1	B	57	VAL	C-N-CA	7.61	127.61	119.78
1	B	7	GLY	N-CA-C	7.59	131.17	113.18
1	A	172	TYR	N-CA-C	7.58	126.94	110.80
1	A	316	VAL	N-CA-C	-7.56	104.42	111.45
1	A	584	LYS	N-CA-C	7.56	124.94	113.61
1	B	116	GLU	N-CA-C	-7.51	100.41	110.55
1	B	83	VAL	N-CA-C	7.49	119.16	111.00
1	B	330	GLY	N-CA-C	7.40	122.00	112.68
1	A	149	PRO	CA-N-CD	-7.29	101.79	112.00
1	A	460	GLN	CB-CA-C	7.28	118.25	110.17
1	A	364	ASN	N-CA-C	-7.27	103.72	112.59
1	B	306	ILE	CB-CA-C	-7.23	102.56	112.24
1	A	24	ILE	N-CA-CB	-7.21	98.31	110.65
1	A	234	PRO	CA-N-CD	-7.20	101.92	112.00
1	B	123	GLU	N-CA-C	7.17	126.08	110.80
1	A	170	PRO	CB-CA-C	7.16	119.66	110.92
1	A	153	THR	N-CA-C	7.16	122.14	113.41
1	B	154	GLN	CG-CD-NE2	7.09	127.04	116.40
1	A	179	GLN	CG-CD-NE2	7.09	127.03	116.40
1	A	83	VAL	N-CA-C	7.04	119.86	111.05
1	B	3	GLU	N-CA-C	-7.04	103.63	113.20
1	A	226	TYR	CA-C-N	-7.03	113.14	120.38
1	A	226	TYR	C-N-CA	-7.03	113.14	120.38
1	B	122	SER	CB-CA-C	7.03	124.40	110.42
1	B	368	ARG	N-CA-C	7.00	119.96	109.25
1	A	460	GLN	CG-CD-NE2	6.91	126.76	116.40
1	B	312	TYR	N-CA-C	6.89	122.02	112.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	173	GLN	CG-CD-NE2	6.88	126.72	116.40
1	A	105	GLN	CG-CD-NE2	6.87	126.70	116.40
1	A	175	LEU	N-CA-C	6.86	120.86	109.95
1	B	125	SER	N-CA-C	-6.83	104.21	112.54
1	B	225	ALA	N-CA-C	6.82	120.99	110.42
1	B	177	ARG	N-CA-C	6.79	120.22	110.10
1	A	171	PRO	CA-C-O	-6.79	108.24	120.60
1	B	126	LYS	N-CA-C	-6.78	103.91	111.71
1	B	180	ILE	N-CA-C	6.78	118.91	109.55
1	A	166	VAL	CB-CA-C	-6.78	100.18	111.29
1	B	67	ARG	N-CA-C	6.72	119.89	109.07
1	B	375	PRO	CA-N-CD	-6.70	102.62	112.00
1	A	165	PHE	CA-C-N	6.69	134.01	121.97
1	A	165	PHE	C-N-CA	6.69	134.01	121.97
1	B	508	GLY	CA-C-N	6.69	126.19	119.24
1	B	508	GLY	C-N-CA	6.69	126.19	119.24
1	A	118	GLU	N-CA-C	6.68	120.40	110.52
1	A	23	THR	CA-C-N	-6.66	109.95	120.47
1	A	23	THR	C-N-CA	-6.66	109.95	120.47
1	B	571	ASN	CA-C-O	-6.64	113.30	121.67
1	A	115	PRO	CA-N-CD	-6.64	102.71	112.00
1	B	108	ASP	CA-CB-CG	6.63	119.23	112.60
1	A	94	GLN	CG-CD-NE2	6.62	126.33	116.40
1	A	135	GLN	CG-CD-NE2	6.62	126.33	116.40
1	A	465	LYS	N-CA-C	-6.61	104.48	112.54
1	A	259	GLN	CG-CD-NE2	6.61	126.31	116.40
1	A	63	GLN	CG-CD-NE2	6.60	126.31	116.40
1	B	140	PHE	CA-C-N	6.57	128.05	119.84
1	B	140	PHE	C-N-CA	6.57	128.05	119.84
1	B	259	GLN	CG-CD-NE2	6.53	126.19	116.40
1	A	165	PHE	CA-CB-CG	-6.51	107.29	113.80
1	B	135	GLN	CG-CD-NE2	6.50	126.15	116.40
1	A	173	GLN	CG-CD-NE2	6.49	126.13	116.40
1	B	573	PRO	CB-CA-C	-6.45	101.41	112.64
1	A	102	GLN	CG-CD-NE2	6.40	126.00	116.40
1	A	86	LEU	CB-CA-C	-6.40	98.63	109.38
1	A	289	GLY	N-CA-C	-6.40	105.03	112.83
1	A	48	PRO	CA-N-CD	-6.39	103.05	112.00
1	A	126	LYS	N-CA-C	-6.38	104.42	111.82
1	B	250	GLN	CG-CD-NE2	6.37	125.95	116.40
1	B	167	CYS	CA-C-N	-6.34	111.91	119.84
1	B	167	CYS	C-N-CA	-6.34	111.91	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	PRO	N-CA-C	-6.31	99.48	112.47
1	A	128	GLN	CG-CD-NE2	6.29	125.84	116.40
1	A	294	GLN	CG-CD-NE2	6.29	125.84	116.40
1	B	208	SER	N-CA-C	-6.27	100.53	109.62
1	A	169	THR	N-CA-C	-6.26	95.98	109.81
1	B	131	GLU	N-CA-C	6.26	121.20	113.50
1	A	154	GLN	CG-CD-NE2	6.25	125.77	116.40
1	B	93	ASP	N-CA-C	-6.25	101.17	110.23
1	B	121	SER	CA-C-N	6.24	133.45	121.54
1	B	121	SER	C-N-CA	6.24	133.45	121.54
1	B	345	PHE	N-CA-C	-6.14	106.39	114.31
1	A	225	ALA	N-CA-C	6.14	118.63	110.53
1	B	476	LEU	N-CA-C	-6.14	103.91	111.40
1	B	53	ASP	N-CA-C	-6.12	106.07	112.93
1	A	209	PRO	CA-C-O	6.11	126.99	118.86
1	B	120	GLY	CA-C-N	6.10	130.41	120.23
1	B	120	GLY	C-N-CA	6.10	130.41	120.23
1	B	441	CYS	N-CA-C	-6.08	104.58	111.14
1	A	594	GLU	N-CA-C	6.05	123.70	110.80
1	A	208	SER	CA-C-N	6.05	125.75	119.64
1	A	208	SER	C-N-CA	6.05	125.75	119.64
1	A	9	PRO	CA-N-CD	-6.05	103.53	112.00
1	A	310	ARG	CA-C-N	-6.04	112.76	122.24
1	A	310	ARG	C-N-CA	-6.04	112.76	122.24
1	B	305	GLN	CG-CD-NE2	6.01	125.41	116.40
1	A	173	GLN	CA-C-N	5.99	132.99	121.54
1	A	173	GLN	C-N-CA	5.99	132.99	121.54
1	B	124	HIS	CB-CG-CD2	5.99	138.98	131.20
1	A	407	MET	N-CA-C	-5.97	101.00	109.96
1	B	465	LYS	N-CA-C	-5.88	105.59	112.89
1	A	172	TYR	CB-CG-CD1	5.88	129.62	120.80
1	B	33	SER	CA-C-N	5.88	127.19	119.84
1	B	33	SER	C-N-CA	5.88	127.19	119.84
1	A	167	CYS	O-C-N	5.86	128.06	121.32
1	B	171	PRO	N-CA-C	5.86	124.55	112.47
1	A	293	TYR	CA-CB-CG	5.85	124.42	113.90
1	B	191	LEU	N-CA-C	-5.83	106.33	113.50
1	A	168	PRO	CA-C-N	5.82	136.01	121.80
1	A	168	PRO	C-N-CA	5.82	136.01	121.80
1	A	57	VAL	CA-C-O	5.81	122.58	119.15
1	A	62	THR	N-CA-C	-5.80	100.38	109.25
1	A	311	ASP	N-CA-C	-5.80	106.83	114.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	98	LEU	N-CA-C	-5.77	105.12	111.82
1	B	29	ASN	N-CA-C	-5.77	104.36	111.75
1	B	426	HIS	N-CA-C	5.75	118.85	110.17
1	A	172	TYR	CB-CG-CD2	-5.73	112.20	120.80
1	A	345	PHE	N-CA-C	-5.71	106.94	114.31
1	A	73	PRO	CA-N-CD	-5.70	104.03	112.00
1	B	137	ASP	CB-CA-C	-5.67	108.40	116.34
1	B	172	TYR	N-CA-C	5.66	122.85	110.80
1	A	87	ASP	CA-CB-CG	5.64	118.25	112.60
1	B	87	ASP	N-CA-C	5.61	117.44	108.41
1	B	205	ASN	N-CA-C	-5.61	99.66	108.63
1	B	10	VAL	N-CA-C	5.58	120.94	108.88
1	B	118	GLU	CB-CA-C	-5.58	103.31	111.23
1	B	16	ASP	N-CA-C	-5.57	98.94	110.80
1	A	137	ASP	N-CA-C	5.57	122.66	110.80
1	A	489	ASN	N-CA-C	5.54	120.04	113.28
1	A	117	THR	N-CA-C	5.53	122.58	110.80
1	A	335	VAL	N-CA-C	5.53	115.82	107.80
1	B	302	ALA	N-CA-C	-5.52	104.65	111.33
1	A	356	SER	N-CA-C	5.49	118.03	111.71
1	B	82	ILE	N-CA-C	5.49	120.75	109.34
1	B	295	GLU	N-CA-C	-5.48	105.46	111.82
1	B	60	GLY	N-CA-C	-5.47	107.03	116.01
1	A	44	ALA	N-CA-C	5.47	122.46	110.80
1	B	363	GLU	N-CA-C	5.44	122.39	110.80
1	A	154	GLN	N-CA-C	-5.44	105.76	112.72
1	B	562	VAL	CA-C-N	5.43	125.34	119.85
1	B	562	VAL	C-N-CA	5.43	125.34	119.85
1	A	208	SER	N-CA-C	5.42	113.10	108.22
1	B	151	LEU	N-CA-C	5.42	117.89	111.33
1	B	572	TYR	N-CA-CB	5.41	120.00	110.37
1	A	141	PRO	CA-C-N	-5.39	115.78	123.11
1	A	141	PRO	C-N-CA	-5.39	115.78	123.11
1	B	21	TYR	N-CA-C	5.38	118.33	109.72
1	B	308	THR	N-CA-C	5.38	117.14	111.28
1	A	427	LYS	N-CA-C	5.36	119.76	111.56
1	B	559	ILE	N-CA-C	-5.36	102.67	109.80
1	B	9	PRO	N-CA-C	-5.35	101.45	112.47
1	B	149	PRO	CB-CA-C	5.35	120.38	111.56
1	B	119	LEU	CA-C-N	-5.34	110.94	121.41
1	B	119	LEU	C-N-CA	-5.34	110.94	121.41
1	A	141	PRO	CA-N-CD	-5.33	104.53	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	LYS	N-CA-C	-5.32	102.47	110.24
1	A	198	SER	N-CA-C	5.31	117.07	111.28
1	B	207	SER	N-CA-C	-5.31	107.47	114.31
1	A	9	PRO	N-CA-C	5.30	123.38	112.47
1	B	275	ARG	CB-CG-CD	5.27	123.42	111.30
1	A	48	PRO	CA-C-N	-5.26	113.40	120.87
1	A	48	PRO	C-N-CA	-5.26	113.40	120.87
1	B	284	ASN	CA-C-N	5.26	124.71	119.24
1	B	284	ASN	C-N-CA	5.26	124.71	119.24
1	A	170	PRO	CA-C-N	5.25	126.40	119.84
1	A	170	PRO	C-N-CA	5.25	126.40	119.84
1	B	529	TRP	N-CA-C	-5.25	102.28	110.10
1	A	95	ASN	N-CA-C	5.22	116.71	110.44
1	A	116	GLU	CB-CG-CD	-5.21	103.74	112.60
1	B	521	GLN	N-CA-C	5.21	116.96	111.28
1	A	233	LYS	CB-CA-C	5.21	116.92	109.11
1	A	426	HIS	N-CA-C	5.20	118.41	110.36
1	B	412	MET	N-CA-C	5.18	117.93	111.24
1	A	464	LEU	N-CA-C	-5.17	105.72	111.36
1	A	312	TYR	N-CA-C	5.11	119.50	113.12
1	B	233	LYS	N-CA-C	-5.09	103.66	110.07
1	B	108	ASP	N-CA-C	-5.08	105.01	111.11
1	B	196	GLU	CA-C-N	5.08	126.19	119.84
1	B	196	GLU	C-N-CA	5.08	126.19	119.84
1	A	359	SER	N-CA-C	5.07	117.96	110.46
1	A	141	PRO	N-CA-C	5.06	119.18	111.34
1	B	8	ALA	CA-C-N	-5.05	113.53	119.84
1	B	8	ALA	C-N-CA	-5.05	113.53	119.84
1	B	564	LEU	N-CA-C	-5.05	105.48	111.69
1	B	428	ILE	N-CA-C	5.05	115.55	110.05
1	B	220	TRP	N-CA-C	5.04	116.65	109.14
1	A	43	LEU	CB-CA-C	-5.03	100.40	110.42
1	B	86	LEU	N-CA-C	5.03	119.44	112.45
1	B	154	GLN	N-CA-C	5.03	119.68	113.50
1	B	573	PRO	N-CD-CG	5.02	110.73	103.20
1	B	573	PRO	N-CA-CB	5.01	108.97	103.26

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	570	ASN	Mainchain

5.2 Too-close contacts ⓘ

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	4764	0	4684	328	0
1	B	4764	0	4683	319	0
2	C	28	0	25	4	0
2	D	28	0	25	2	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
3	E	39	0	34	7	0
3	H	39	0	34	1	0
4	A	14	0	13	0	0
4	B	14	0	13	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	4	0	0	0	0
6	B	4	0	0	0	0
7	A	43	0	30	10	0
7	B	43	0	30	12	0
8	A	9	0	4	2	0
8	B	9	0	3	1	0
9	A	86	0	0	10	0
9	B	62	0	0	4	0
All	All	10008	0	9628	658	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 34.

All (658) 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)
1:B:67:ARG:N	1:B:67:ARG:CA	1.83	1.40
1:A:108:ASP:OD2	7:A:605:HEM:HMD1	1.25	1.33
1:A:108:ASP:OD2	7:A:605:HEM:CMD	1.81	1.28
1:A:233:LYS:HB3	1:A:234:PRO:HD2	1.31	1.11
1:A:167:CYS:HB2	1:A:168:PRO:HD2	1.13	1.10
1:A:209:PRO:O	1:A:289:GLY:HA2	1.50	1.08
1:B:67:ARG:N	1:B:67:ARG:CB	2.17	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ARG:HH11	1:A:360:ARG:HB2	1.21	1.05
1:A:167:CYS:HB2	1:A:168:PRO:CD	1.85	1.04
1:A:294:GLN:HA	1:A:297:ARG:HB3	1.37	1.04
1:A:172:TYR:HB3	1:A:173:GLN:HG2	1.35	1.03
1:A:175:LEU:HD23	1:A:176:ALA:N	1.73	1.01
1:B:381:PHE:CZ	1:B:424:PRO:HG3	1.95	1.00
1:A:42:ALA:HB2	1:A:166:VAL:CG1	1.95	0.96
1:B:327:ARG:HD2	1:B:327:ARG:H	1.34	0.93
1:B:66:THR:C	1:B:67:ARG:CA	2.43	0.92
1:B:179:GLN:HG3	1:B:444:HIS:CE1	2.04	0.92
1:A:233:LYS:CB	1:A:234:PRO:HD2	1.97	0.91
1:B:551:ARG:HD3	1:B:583:ASP:O	1.71	0.90
1:A:24:ILE:HD13	1:A:200:ALA:HB1	1.53	0.90
1:B:179:GLN:HG3	1:B:444:HIS:ND1	1.87	0.89
1:A:169:THR:HB	1:A:170:PRO:CD	2.01	0.89
1:A:167:CYS:CB	1:A:168:PRO:HD2	1.98	0.88
1:A:108:ASP:OD2	7:A:605:HEM:HMD2	1.72	0.88
1:A:42:ALA:CB	1:A:166:VAL:HG11	2.04	0.87
1:A:123:GLU:HG3	1:A:125:SER:H	1.40	0.87
1:B:119:LEU:HD13	1:B:119:LEU:N	1.89	0.87
1:A:167:CYS:CB	1:A:168:PRO:CD	2.49	0.87
1:A:175:LEU:HD23	1:A:176:ALA:H	1.36	0.86
1:A:64:ARG:HA	1:A:71:ARG:NH2	1.90	0.86
1:B:158:MET:HE2	1:B:431:PHE:HA	1.55	0.86
1:A:42:ALA:CB	1:A:166:VAL:CG1	2.54	0.85
1:A:209:PRO:O	1:A:289:GLY:CA	2.24	0.84
1:A:344:THR:HB	7:A:605:HEM:O1D	1.76	0.84
1:A:119:LEU:CD2	1:A:120:GLY:H	1.90	0.84
1:A:260:ILE:HB	9:A:3060:HOH:O	1.78	0.83
1:A:119:LEU:HD23	1:A:120:GLY:H	1.42	0.82
1:B:42:ALA:HB2	1:B:166:VAL:CG1	2.10	0.82
1:B:362:ASP:HB3	1:B:368:ARG:HB3	1.61	0.82
1:B:588:SER:OG	1:B:589:PRO:HD3	1.78	0.82
1:A:42:ALA:HB3	1:A:166:VAL:HG11	1.58	0.82
1:A:432:ASP:O	1:A:436:ILE:HG12	1.79	0.82
1:B:130:GLU:HG3	1:B:159:PRO:HG3	1.62	0.82
1:B:505:GLY:O	1:B:506:ARG:HD3	1.81	0.81
1:B:595:ASN:N	1:B:595:ASN:HD22	1.78	0.81
1:A:144:PHE:HE2	1:A:158:MET:HE3	1.45	0.80
1:A:24:ILE:HD11	1:A:293:TYR:OH	1.82	0.80
1:B:572:TYR:CD2	1:B:573:PRO:HD3	2.15	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:THR:HG22	1:A:27:ASP:H	1.46	0.79
1:B:67:ARG:N	1:B:67:ARG:HB2	1.96	0.79
1:B:123:GLU:OE2	1:B:125:SER:HB2	1.82	0.79
1:B:360:ARG:NH1	1:B:372:ALA:HA	1.98	0.78
1:A:341:ASN:HB3	8:A:3002:FMT:H	1.66	0.78
1:A:279:GLU:HG2	1:A:587:LEU:HD12	1.66	0.77
1:A:30:ASN:HB3	1:A:33:SER:O	1.85	0.77
1:B:295:GLU:O	1:B:299:ILE:HG12	1.83	0.77
1:A:272:GLU:HA	1:A:272:GLU:OE1	1.83	0.77
1:A:96:ARG:NH2	1:A:315:ILE:HB	2.00	0.76
1:B:377:HIS:HB3	1:B:416:GLU:OE2	1.85	0.76
1:A:368:ARG:HG3	1:A:368:ARG:HH11	1.49	0.76
1:A:260:ILE:HD11	1:A:386:ILE:HG13	1.68	0.76
1:A:52:GLU:HB3	1:A:57:VAL:HG23	1.67	0.76
1:B:327:ARG:HD2	1:B:327:ARG:N	2.02	0.75
1:A:230:ASN:ND2	1:A:232:MET:HB3	2.01	0.75
7:B:605:HEM:HMC2	7:B:605:HEM:HBC2	1.69	0.74
1:A:169:THR:HB	1:A:170:PRO:HD2	1.69	0.74
1:B:8:ALA:HB1	1:B:10:VAL:CG1	2.17	0.74
1:B:204:ARG:HB3	1:B:206:LEU:HG	1.70	0.74
1:A:501:MET:HE2	1:A:506:ARG:HD2	1.70	0.73
1:B:300:LEU:O	1:B:303:PHE:HB3	1.88	0.73
1:A:170:PRO:HB3	1:A:171:PRO:HD2	1.70	0.73
1:A:425:THR:HB	1:A:426:HIS:CD2	2.24	0.73
1:B:16:ASP:O	1:B:17:GLU:HB2	1.88	0.73
1:B:173:GLN:HG3	1:B:174:SER:N	2.04	0.73
1:B:400:LEU:HD11	1:B:553:ILE:HD13	1.71	0.73
1:B:173:GLN:HG3	1:B:174:SER:H	1.54	0.73
1:A:230:ASN:HD21	1:A:232:MET:HB3	1.53	0.72
1:B:42:ALA:HB2	1:B:166:VAL:HG11	1.71	0.72
1:A:368:ARG:HG3	1:A:368:ARG:O	1.91	0.71
1:B:400:LEU:HD11	1:B:553:ILE:CD1	2.21	0.71
1:A:169:THR:HB	1:A:170:PRO:HD3	1.71	0.70
1:A:288:ASP:HB3	1:A:291:LYS:H	1.56	0.70
1:A:360:ARG:HB2	1:A:360:ARG:NH1	2.03	0.70
1:A:76:ARG:NH2	1:A:150:LYS:HG3	2.06	0.70
1:B:381:PHE:HZ	1:B:424:PRO:HG3	1.54	0.70
1:A:467:LEU:HG	1:A:471:LEU:HD22	1.73	0.70
1:A:464:LEU:O	1:A:468:GLN:HG3	1.92	0.69
1:A:260:ILE:CD1	1:A:386:ILE:HG13	2.22	0.69
1:B:544:LEU:O	1:B:547:VAL:HG22	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:TYR:N	1:B:172:TYR:CD1	2.60	0.69
1:A:56:ALA:O	1:A:58:PRO:HD3	1.93	0.68
1:A:362:ASP:HB2	1:A:368:ARG:CB	2.22	0.68
1:A:475:ILE:HD11	1:A:479:LYS:HE3	1.75	0.68
1:A:368:ARG:HG3	1:A:368:ARG:NH1	2.05	0.68
1:A:408:ASN:O	1:A:410:ASN:N	2.26	0.68
1:B:279:GLU:O	1:B:282:ARG:HG2	1.94	0.68
1:B:464:LEU:O	1:B:464:LEU:HD22	1.95	0.67
1:B:258:GLU:O	1:B:381:PHE:N	2.27	0.67
1:A:260:ILE:HD11	1:A:386:ILE:CG1	2.24	0.67
1:A:360:ARG:HH11	1:A:360:ARG:CB	2.05	0.67
1:A:298:LYS:HG2	1:A:536:PHE:CZ	2.30	0.67
1:A:137:ASP:OD1	1:A:138:ASN:N	2.28	0.67
1:A:327:ARG:H	1:A:327:ARG:HD2	1.60	0.67
1:B:8:ALA:HB1	1:B:10:VAL:HG13	1.76	0.67
1:B:257:SER:O	1:B:258:GLU:C	2.35	0.67
1:A:139:CYS:SG	1:A:141:PRO:HD3	2.34	0.66
1:A:360:ARG:HH12	1:A:372:ALA:HA	1.60	0.66
1:B:218:GLU:HG3	9:B:3060:HOH:O	1.94	0.66
1:B:10:VAL:HG23	1:B:10:VAL:O	1.94	0.66
1:A:588:SER:N	1:A:589:PRO:HD2	2.10	0.66
1:B:53:ASP:CG	1:B:57:VAL:HG13	2.21	0.66
1:A:191:LEU:HD23	1:A:191:LEU:H	1.61	0.66
1:A:24:ILE:HD13	1:A:200:ALA:CB	2.22	0.66
1:A:227:PRO:HG2	1:A:249:PHE:CG	2.31	0.66
1:B:123:GLU:HB3	1:B:126:LYS:HE3	1.77	0.66
1:A:549:PHE:CE1	1:A:552:LEU:HD23	2.30	0.65
1:B:370:PRO:HG3	3:H:3:MAN:H62	1.76	0.65
1:B:166:VAL:HG23	1:B:167:CYS:SG	2.37	0.65
1:A:445:GLY:O	1:A:447:PRO:HD3	1.95	0.65
1:B:64:ARG:O	1:B:64:ARG:HG2	1.96	0.65
1:A:123:GLU:HB3	1:A:126:LYS:NZ	2.12	0.65
1:A:391:GLY:HA3	9:A:3007:HOH:O	1.95	0.64
1:A:294:GLN:CA	1:A:297:ARG:HB3	2.21	0.64
1:A:377:HIS:CE1	1:A:378:THR:HG23	2.33	0.64
1:B:410:ASN:O	1:B:410:ASN:ND2	2.31	0.64
1:B:407:MET:HB3	1:B:501:MET:CE	2.27	0.64
1:B:258:GLU:O	1:B:380:PHE:HA	1.98	0.64
1:A:368:ARG:HH11	1:A:368:ARG:CG	2.10	0.64
1:B:432:ASP:O	1:B:436:ILE:HG13	1.97	0.64
1:A:362:ASP:HB2	1:A:368:ARG:HB3	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:579:CYS:O	1:B:582:ILE:HG23	1.98	0.63
1:B:2:TRP:HD1	1:B:4:VAL:CG1	2.12	0.63
1:A:73:PRO:HG3	1:A:483:LEU:O	1.99	0.63
1:A:259:GLN:OE1	1:A:261:LEU:HB2	1.98	0.63
1:B:113:PHE:HA	7:B:605:HEM:O2D	1.99	0.63
1:B:169:THR:N	1:B:170:PRO:CD	2.62	0.63
1:A:44:ALA:O	1:A:340:SER:HA	1.99	0.62
1:B:7:GLY:O	1:B:8:ALA:HB2	1.98	0.62
1:A:327:ARG:HD2	1:A:327:ARG:N	2.14	0.62
1:B:359:SER:HB3	1:B:373:GLU:HG2	1.81	0.62
1:A:324:TRP:O	1:A:326:PRO:HD3	1.99	0.62
1:A:504:ARG:NH2	2:C:2:NAG:H5	2.14	0.62
1:B:91:VAL:HG22	1:B:91:VAL:O	2.00	0.62
1:B:407:MET:HB3	1:B:501:MET:HE3	1.81	0.62
1:A:144:PHE:CE2	1:A:158:MET:HE3	2.32	0.62
1:A:382:ASN:HB3	9:A:3060:HOH:O	1.99	0.62
1:A:492:ILE:CG2	1:A:493:TRP:N	2.63	0.62
1:B:4:VAL:HG23	1:B:4:VAL:O	1.99	0.62
1:A:91:VAL:O	1:A:91:VAL:HG22	1.98	0.61
1:B:2:TRP:HD1	1:B:4:VAL:HG11	1.65	0.61
1:B:3:GLU:O	1:B:5:GLY:N	2.33	0.61
1:B:10:VAL:O	1:B:10:VAL:CG2	2.49	0.61
1:B:368:ARG:O	1:B:372:ALA:HB2	2.00	0.61
1:B:453:ARG:NH2	1:B:458:LEU:HB3	2.15	0.61
1:A:179:GLN:NE2	1:A:444:HIS:HA	2.15	0.61
1:B:123:GLU:HG3	1:B:124:HIS:N	2.15	0.61
1:B:377:HIS:C	1:B:377:HIS:CD2	2.79	0.61
1:A:46:TRP:CE2	1:A:340:SER:HB3	2.36	0.61
1:B:173:GLN:CG	1:B:174:SER:N	2.62	0.61
1:A:42:ALA:HB1	1:A:179:GLN:O	2.00	0.61
1:A:229:PHE:CG	1:A:247:PRO:HG2	2.36	0.61
1:B:343:PHE:CD1	1:B:518:GLN:HG2	2.36	0.61
1:A:123:GLU:HB3	1:A:126:LYS:HZ2	1.66	0.61
1:A:169:THR:CB	1:A:170:PRO:HD2	2.30	0.61
1:B:169:THR:OG1	1:B:170:PRO:HD3	2.01	0.61
1:A:362:ASP:O	1:A:364:ASN:N	2.34	0.60
1:B:53:ASP:C	1:B:55:LEU:H	2.09	0.60
1:B:467:LEU:HD23	1:B:477:ALA:HA	1.82	0.60
1:A:119:LEU:CD2	1:A:120:GLY:N	2.64	0.60
1:A:587:LEU:C	1:A:589:PRO:HD2	2.26	0.60
1:B:342:VAL:HB	1:B:446:MET:HE1	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:HIS:HB3	1:A:416:GLU:OE2	2.01	0.60
1:A:98:LEU:O	1:A:98:LEU:HD22	2.02	0.60
1:A:274:ASN:HB3	1:A:278:ARG:NH2	2.16	0.60
1:B:170:PRO:HB3	1:B:171:PRO:HD2	1.83	0.60
1:B:420:LYS:HA	1:B:429:HIS:O	2.01	0.60
1:A:293:TYR:C	1:A:294:GLN:HG2	2.27	0.60
1:A:362:ASP:N	1:A:368:ARG:HB3	2.17	0.59
1:A:407:MET:SD	1:A:408:ASN:N	2.75	0.59
1:A:310:ARG:O	1:A:310:ARG:HG2	2.02	0.59
1:B:361:LEU:HB3	1:B:365:TYR:HA	1.84	0.59
1:B:408:ASN:HB3	1:B:411:LYS:HB2	1.85	0.59
1:A:449:TYR:HB2	1:A:490:ILE:HG22	1.85	0.59
1:A:313:LEU:N	1:A:314:PRO:CD	2.64	0.59
1:B:119:LEU:N	1:B:119:LEU:CD1	2.61	0.59
1:A:108:ASP:CG	7:A:605:HEM:HMD1	2.18	0.59
1:A:142:ILE:O	1:A:157:CYS:HB2	2.02	0.58
1:A:450:ASN:HD21	1:A:487:PRO:HB2	1.68	0.58
1:A:61:TRP:CD1	1:A:61:TRP:C	2.81	0.58
1:B:98:LEU:HD13	1:B:399:LEU:HD23	1.85	0.58
1:A:537:THR:OG1	1:A:540:GLN:HG3	2.03	0.58
1:A:58:PRO:HG3	1:A:443:ASP:HA	1.84	0.58
1:A:17:GLU:O	1:A:17:GLU:HG3	2.02	0.58
1:B:187:LEU:HB3	1:B:305:GLN:HG3	1.84	0.58
1:A:351:HIS:CE1	1:A:433:LEU:HD21	2.38	0.58
1:A:221:ASP:O	1:A:224:LEU:HB2	2.04	0.57
1:B:105:GLN:NE2	7:B:605:HEM:C4B	2.71	0.57
1:A:27:ASP:O	1:A:28:CYS:HB2	2.03	0.57
1:A:202:ARG:NH2	1:A:250:GLN:HE22	2.03	0.57
1:B:2:TRP:C	1:B:4:VAL:H	2.11	0.57
1:B:481:LEU:C	1:B:483:LEU:H	2.10	0.57
1:A:504:ARG:HH21	2:C:2:NAG:H5	1.69	0.57
1:A:150:LYS:HD3	1:A:154:GLN:OE1	2.04	0.57
3:E:2:NAG:H5	3:E:3:MAN:H2	1.85	0.57
3:E:2:NAG:H61	3:E:2:NAG:H2	1.85	0.57
1:B:287:TRP:CE2	1:B:291:LYS:HE3	2.39	0.57
1:B:417:LEU:HD11	7:B:605:HEM:HBB2	1.87	0.57
1:A:169:THR:CB	1:A:170:PRO:CD	2.74	0.56
1:B:323:LYS:HD3	1:B:324:TRP:NE1	2.20	0.56
1:A:229:PHE:CD1	1:A:247:PRO:HG2	2.40	0.56
1:B:9:PRO:HG2	1:B:10:VAL:H	1.70	0.56
1:B:9:PRO:CG	1:B:10:VAL:H	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:GLN:OE1	1:A:261:LEU:HB3	2.05	0.56
1:B:362:ASP:O	1:B:364:ASN:N	2.38	0.56
1:B:480:LEU:O	1:B:483:LEU:HB2	2.04	0.56
1:A:473:ASN:OD1	1:A:475:ILE:HG23	2.06	0.56
1:B:191:LEU:H	1:B:191:LEU:HD23	1.70	0.56
1:A:492:ILE:HG23	1:A:493:TRP:N	2.21	0.56
1:B:158:MET:CE	1:B:431:PHE:HA	2.33	0.56
1:A:137:ASP:OD1	1:A:137:ASP:C	2.48	0.56
1:B:283:LEU:C	1:B:285:PRO:HD3	2.31	0.56
1:B:327:ARG:H	1:B:327:ARG:CD	2.14	0.56
1:B:299:ILE:HD11	1:B:590:TRP:NE1	2.20	0.56
1:A:356:SER:HA	1:A:414:THR:HG21	1.86	0.56
9:A:3023:HOH:O	2:D:1:NAG:H61	2.06	0.56
1:B:276:LEU:O	1:B:280:LEU:HG	2.06	0.55
1:B:362:ASP:O	1:B:362:ASP:OD2	2.24	0.55
1:B:494:ILE:CG2	1:B:495:GLY:N	2.69	0.55
1:A:551:ARG:CA	1:A:582:ILE:HD11	2.36	0.55
1:A:377:HIS:CE1	1:A:378:THR:CG2	2.89	0.55
1:A:158:MET:HB2	1:A:431:PHE:CE2	2.42	0.55
1:B:168:PRO:HB3	1:B:170:PRO:HD2	1.89	0.55
1:B:481:LEU:C	1:B:483:LEU:N	2.59	0.55
1:B:595:ASN:N	1:B:595:ASN:ND2	2.50	0.55
1:A:542:ASP:HA	1:A:545:GLN:NE2	2.21	0.54
1:B:362:ASP:OD2	1:B:362:ASP:C	2.49	0.54
1:A:481:LEU:C	1:A:483:LEU:N	2.57	0.54
1:A:494:ILE:HG23	1:A:495:GLY:H	1.73	0.54
1:B:66:THR:O	1:B:67:ARG:CA	2.55	0.54
1:A:119:LEU:N	1:A:119:LEU:HD13	2.22	0.54
1:B:216:ASN:ND2	9:B:3060:HOH:O	2.41	0.54
1:A:556:ASN:C	1:A:557:THR:HG23	2.33	0.54
1:B:354:VAL:HG11	7:B:605:HEM:HBB1	1.88	0.54
1:A:564:LEU:HD22	9:A:3079:HOH:O	2.07	0.54
1:B:27:ASP:OD1	1:B:35:ALA:HA	2.08	0.54
1:A:408:ASN:C	1:A:410:ASN:H	2.15	0.54
1:A:217:GLN:NE2	2:D:1:NAG:H83	2.23	0.54
1:A:412:MET:HB2	9:A:3062:HOH:O	2.07	0.54
1:B:106:ILE:HD11	1:B:265:VAL:HB	1.89	0.54
1:B:165:PHE:CG	1:B:177:ARG:HD2	2.43	0.54
1:B:75:ALA:HB2	1:B:439:GLN:HG3	1.89	0.53
1:B:106:ILE:HG23	1:B:191:LEU:HD11	1.90	0.53
1:B:260:ILE:HG21	1:B:379:LEU:HD22	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ARG:NH1	1:A:372:ALA:HA	2.23	0.53
1:A:588:SER:N	1:A:589:PRO:CD	2.70	0.53
1:A:85:TYR:CD2	1:A:411:LYS:HB3	2.43	0.53
1:A:162:ARG:HB3	1:A:443:ASP:OD1	2.08	0.53
1:A:481:LEU:C	1:A:483:LEU:H	2.14	0.53
1:B:142:ILE:HA	1:B:439:GLN:HE22	1.72	0.53
1:A:148:ASP:OD1	1:A:148:ASP:C	2.52	0.53
1:B:242:THR:O	1:B:245:ARG:HG3	2.09	0.53
1:B:360:ARG:HH11	1:B:372:ALA:HA	1.73	0.53
1:A:189:ALA:HB2	1:A:304:ILE:HG13	1.90	0.53
1:B:66:THR:HB	1:B:70:PHE:N	2.24	0.53
1:B:360:ARG:HH11	1:B:360:ARG:HB2	1.74	0.53
1:A:432:ASP:OD1	1:A:432:ASP:C	2.51	0.53
1:A:201:SER:HA	9:A:3072:HOH:O	2.09	0.53
1:A:258:GLU:O	1:A:380:PHE:HA	2.08	0.53
1:A:584:LYS:O	1:A:585:LEU:HB3	2.07	0.53
1:B:77:GLU:OE2	1:B:81:LYS:NZ	2.42	0.53
1:B:108:ASP:OD2	7:B:605:HEM:HH1	2.07	0.53
1:A:226:TYR:CE1	1:A:387:ILE:HG12	2.43	0.53
1:A:309:PHE:CD1	1:A:529:TRP:HH2	2.27	0.53
1:B:328:TYR:OH	1:B:532:ASN:HB2	2.09	0.53
1:A:131:GLU:HB2	1:A:132:TYR:CE1	2.44	0.53
1:A:322:GLN:HE21	1:A:327:ARG:HH21	1.57	0.53
1:A:25:THR:HG21	1:A:27:ASP:OD2	2.09	0.52
1:B:16:ASP:OD2	1:B:19:SER:HB2	2.10	0.52
1:B:267:THR:O	1:B:271:ARG:HG3	2.08	0.52
3:E:1:NAG:H62	3:E:2:NAG:N2	2.24	0.52
1:A:494:ILE:HG23	1:A:495:GLY:N	2.23	0.52
1:B:98:LEU:HD13	1:B:399:LEU:CD2	2.38	0.52
1:B:142:ILE:HA	1:B:439:GLN:NE2	2.24	0.52
1:A:113:PHE:HB3	1:A:182:SER:OG	2.10	0.52
1:B:360:ARG:HH12	1:B:372:ALA:HA	1.73	0.52
1:A:362:ASP:CA	1:A:368:ARG:HB3	2.39	0.52
1:A:501:MET:CE	1:A:506:ARG:HD2	2.37	0.52
1:B:9:PRO:O	1:B:10:VAL:HG22	2.10	0.52
1:A:52:GLU:OE1	1:A:65:LYS:HD2	2.10	0.52
1:A:321:MET:HE3	1:A:325:ILE:HD12	1.92	0.52
1:B:165:PHE:CE2	1:B:172:TYR:HB3	2.44	0.52
1:B:213:MET:HG2	1:B:273:HIS:CD2	2.45	0.52
1:B:386:ILE:HG23	1:B:392:ILE:HG22	1.91	0.52
1:B:410:ASN:O	1:B:411:LYS:HG3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:HIS:HB3	9:A:3080:HOH:O	2.09	0.52
1:A:242:THR:HB	1:A:245:ARG:NH2	2.25	0.52
1:B:52:GLU:HB3	1:B:57:VAL:HG22	1.90	0.52
1:B:173:GLN:HG3	1:B:174:SER:OG	2.10	0.52
1:A:132:TYR:N	1:A:132:TYR:CD1	2.78	0.51
1:A:193:TYR:OH	1:A:300:LEU:HD23	2.10	0.51
1:B:142:ILE:HG12	1:B:439:GLN:NE2	2.25	0.51
1:B:351:HIS:HD1	1:B:437:ASN:HD21	1.58	0.51
1:A:501:MET:HE2	1:A:506:ARG:HA	1.92	0.51
1:B:61:TRP:O	1:B:135:GLN:NE2	2.43	0.51
1:B:362:ASP:O	1:B:365:TYR:N	2.43	0.51
1:A:16:ASP:O	1:A:17:GLU:CB	2.59	0.51
1:B:186:PHE:O	1:B:188:ASP:N	2.42	0.51
1:A:381:PHE:CZ	1:A:424:PRO:HB3	2.46	0.51
1:B:118:GLU:C	1:B:119:LEU:HD13	2.36	0.51
1:B:253:ASP:OD2	1:B:255:ARG:HD3	2.11	0.51
1:B:393:ASP:N	1:B:394:PRO:HD2	2.25	0.51
1:B:95:ASN:O	1:B:96:ARG:HD3	2.09	0.51
1:B:342:VAL:CB	1:B:446:MET:HE1	2.41	0.51
1:A:175:LEU:CD2	1:A:176:ALA:N	2.62	0.51
1:B:13:VAL:HG21	1:B:25:THR:HG21	1.92	0.51
1:A:362:ASP:CB	1:A:368:ARG:HB3	2.40	0.51
1:A:396:VAL:CG1	1:A:553:ILE:HD12	2.40	0.51
1:B:208:SER:HB2	1:B:210:LEU:HD21	1.92	0.51
1:A:101:MET:SD	1:A:101:MET:C	2.94	0.50
1:A:255:ARG:O	1:A:257:SER:N	2.43	0.50
1:A:571:ASN:N	1:A:575:ASP:OD2	2.41	0.50
1:B:101:MET:HE3	1:B:354:VAL:HG22	1.93	0.50
1:B:471:LEU:HD12	1:B:499:GLU:HA	1.92	0.50
1:B:475:ILE:HD11	1:B:479:LYS:NZ	2.26	0.50
1:B:453:ARG:CZ	1:B:458:LEU:HB3	2.41	0.50
1:B:66:THR:HB	1:B:70:PHE:H	1.77	0.50
1:B:298:LYS:HG2	1:B:536:PHE:CZ	2.46	0.50
1:B:464:LEU:HD22	1:B:464:LEU:C	2.35	0.50
1:A:585:LEU:HG	1:A:585:LEU:O	2.11	0.50
1:B:112:ASP:HA	1:B:183:VAL:CG2	2.41	0.50
1:B:258:GLU:OE1	7:B:605:HEM:C2B	2.64	0.50
1:B:375:PRO:O	1:B:376:LEU:C	2.54	0.50
1:A:522:ILE:HD12	9:A:3064:HOH:O	2.12	0.50
1:B:30:ASN:HB3	1:B:33:SER:O	2.12	0.50
1:A:400:LEU:HD13	1:A:563:PRO:HD3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:564:LEU:HB2	1:A:565:HIS:CD2	2.47	0.50
1:B:368:ARG:O	1:B:368:ARG:HG3	2.11	0.50
1:B:165:PHE:HA	1:B:179:GLN:HA	1.92	0.50
1:B:169:THR:H	1:B:170:PRO:CD	2.25	0.50
1:A:88:GLU:OE1	1:A:356:SER:HB3	2.12	0.50
1:A:341:ASN:CB	8:A:3002:FMT:H	2.39	0.50
1:A:505:GLY:O	1:A:506:ARG:HD3	2.11	0.50
1:B:187:LEU:HD13	1:B:305:GLN:HA	1.93	0.50
1:B:208:SER:HB2	1:B:210:LEU:CD2	2.42	0.50
2:C:1:NAG:O3	2:C:1:NAG:C7	2.60	0.50
1:A:42:ALA:HB2	1:A:166:VAL:HG12	1.90	0.49
1:B:82:ILE:HD13	1:B:480:LEU:HD13	1.94	0.49
1:B:93:ASP:CA	1:B:406:LEU:HB2	2.43	0.49
1:B:245:ARG:HD2	9:B:3017:HOH:O	2.13	0.49
1:B:130:GLU:CD	1:B:426:HIS:HD1	2.20	0.49
1:B:454:GLY:O	1:B:456:CYS:N	2.45	0.49
1:A:406:LEU:HD12	1:A:407:MET:H	1.76	0.49
1:A:230:ASN:HD21	1:A:232:MET:CB	2.22	0.49
1:A:351:HIS:CD2	1:A:433:LEU:HD21	2.48	0.49
1:A:42:ALA:HB2	1:A:166:VAL:HG13	1.90	0.49
1:A:148:ASP:O	1:A:151:LEU:HB2	2.13	0.49
1:A:365:TYR:CE1	1:A:397:ARG:HB3	2.48	0.49
1:A:420:LYS:HA	1:A:429:HIS:O	2.13	0.49
1:B:62:THR:O	1:B:63:GLN:C	2.55	0.49
1:B:170:PRO:CB	1:B:171:PRO:HD2	2.42	0.49
1:B:546:LYS:HD3	1:B:583:ASP:OD2	2.13	0.49
1:A:116:GLU:HA	1:A:163:ALA:HA	1.94	0.49
1:B:52:GLU:C	1:B:54:GLY:H	2.21	0.49
1:B:365:TYR:CE2	1:B:397:ARG:HD3	2.48	0.49
1:B:579:CYS:O	1:B:582:ILE:CG2	2.60	0.49
1:A:132:TYR:HB2	1:A:134:ILE:HG13	1.95	0.48
1:A:119:LEU:HD22	1:A:119:LEU:H	1.78	0.48
1:A:170:PRO:CB	1:A:171:PRO:HD2	2.35	0.48
1:A:244:ALA:HB2	3:E:1:NAG:O6	2.12	0.48
1:A:362:ASP:HB2	1:A:368:ARG:HB2	1.95	0.48
1:A:556:ASN:O	1:A:557:THR:CG2	2.62	0.48
1:A:589:PRO:HG2	1:A:590:TRP:CE3	2.49	0.48
1:B:123:GLU:HG2	1:B:126:LYS:H	1.76	0.48
1:B:168:PRO:CB	1:B:170:PRO:HD2	2.42	0.48
1:B:282:ARG:HG3	1:B:283:LEU:N	2.29	0.48
1:A:556:ASN:C	1:A:557:THR:CG2	2.87	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:ASP:O	1:B:224:LEU:HB2	2.12	0.48
1:B:280:LEU:CD2	1:B:587:LEU:HD13	2.44	0.48
1:A:137:ASP:OD1	1:A:138:ASN:HB2	2.14	0.48
1:A:422:PHE:HE1	1:A:427:LYS:O	1.95	0.48
1:B:62:THR:O	1:B:62:THR:HG22	2.13	0.48
1:B:180:ILE:CG2	1:B:181:ASN:N	2.76	0.48
1:B:343:PHE:CG	1:B:518:GLN:HG2	2.49	0.48
1:B:66:THR:HB	1:B:70:PHE:O	2.12	0.48
1:B:82:ILE:HD13	1:B:480:LEU:CD1	2.44	0.48
1:B:283:LEU:O	1:B:285:PRO:HD3	2.14	0.48
1:A:541:ARG:O	1:A:544:LEU:N	2.46	0.48
1:B:16:ASP:CG	1:B:17:GLU:N	2.69	0.48
1:B:365:TYR:CZ	1:B:397:ARG:HD3	2.49	0.48
1:B:560:THR:HB	8:B:3004:FMT:O1	2.14	0.48
1:A:274:ASN:HB3	1:A:278:ARG:HH22	1.79	0.48
1:A:300:LEU:O	1:A:303:PHE:HB3	2.12	0.48
1:A:351:HIS:NE2	1:A:433:LEU:HD21	2.28	0.48
1:B:64:ARG:O	1:B:64:ARG:CG	2.60	0.48
1:B:97:SER:OG	1:B:99:LEU:HB3	2.14	0.48
1:B:150:LYS:HE3	1:B:430:GLY:O	2.14	0.47
1:B:302:ALA:O	1:B:306:ILE:HG13	2.14	0.47
1:A:116:GLU:H	1:A:116:GLU:HG3	1.58	0.47
1:A:119:LEU:O	1:A:126:LYS:HE3	2.15	0.47
1:A:396:VAL:HG11	1:A:553:ILE:HD12	1.96	0.47
1:B:99:LEU:HG	1:B:567:PHE:CE1	2.49	0.47
1:A:167:CYS:CB	1:A:168:PRO:HD3	2.40	0.47
1:A:175:LEU:O	1:A:176:ALA:C	2.57	0.47
1:A:408:ASN:C	1:A:410:ASN:N	2.73	0.47
1:B:139:CYS:SG	1:B:141:PRO:HD3	2.54	0.47
1:B:475:ILE:O	1:B:475:ILE:HG13	2.14	0.47
1:B:560:THR:HG23	1:B:578:ASP:OD1	2.14	0.47
1:B:98:LEU:HB3	1:B:399:LEU:HD23	1.95	0.47
1:B:165:PHE:CZ	1:B:172:TYR:HB3	2.49	0.47
1:B:332:ASN:O	1:B:334:SER:N	2.47	0.47
1:B:484:TYR:O	1:B:485:LYS:HB2	2.14	0.47
1:B:43:LEU:HD23	1:B:181:ASN:HB2	1.95	0.47
1:B:274:ASN:O	1:B:275:ARG:C	2.56	0.47
1:A:96:ARG:NH2	1:A:315:ILE:O	2.48	0.47
1:A:362:ASP:OD1	1:A:363:GLU:N	2.48	0.47
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.49	0.47
1:B:159:PRO:HD2	9:B:3023:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:450:ASN:HD21	1:B:487:PRO:HB2	1.80	0.47
1:B:560:THR:HG22	1:B:561:LYS:CD	2.45	0.47
1:B:91:VAL:HG13	1:B:92:LEU:HD12	1.94	0.47
1:B:467:LEU:HG	1:B:471:LEU:HD22	1.97	0.47
1:B:35:ALA:HB1	1:B:41:ARG:CZ	2.45	0.47
7:B:605:HEM:CBB	7:B:605:HEM:HMB1	2.45	0.47
1:A:165:PHE:CD1	1:A:165:PHE:N	2.83	0.47
1:B:123:GLU:HG2	1:B:126:LYS:HB2	1.96	0.47
1:B:160:PHE:CD1	1:B:160:PHE:C	2.93	0.47
1:A:239:PHE:O	1:A:240:ILE:C	2.58	0.46
1:B:471:LEU:HD12	1:B:471:LEU:HA	1.75	0.46
1:A:16:ASP:O	1:A:17:GLU:HB3	2.14	0.46
1:A:406:LEU:HD12	1:A:407:MET:N	2.29	0.46
1:A:368:ARG:O	1:A:368:ARG:CG	2.62	0.46
1:A:462:LYS:NZ	1:A:488:ASP:OD2	2.48	0.46
2:C:1:NAG:O3	2:C:1:NAG:O7	2.34	0.46
1:A:131:GLU:HB2	1:A:132:TYR:CD1	2.50	0.46
1:A:199:LEU:HD12	1:A:199:LEU:O	2.15	0.46
1:A:350:GLY:HA3	7:A:605:HEM:HBC2	1.97	0.46
1:A:407:MET:SD	1:A:407:MET:C	2.98	0.46
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.96	0.46
1:B:132:TYR:HB3	1:B:134:ILE:CD1	2.45	0.46
1:B:158:MET:HE1	1:B:432:ASP:N	2.30	0.46
1:A:85:TYR:HD1	1:A:87:ASP:O	1.98	0.46
1:B:258:GLU:OE2	7:B:605:HEM:HHB	2.14	0.46
1:B:376:LEU:HD13	1:B:380:PHE:CZ	2.49	0.46
1:A:130:GLU:O	1:A:130:GLU:HG2	2.16	0.46
1:A:475:ILE:HD12	1:A:475:ILE:O	2.16	0.46
1:A:510:LEU:O	1:A:513:CYS:HB3	2.15	0.46
1:B:53:ASP:O	1:B:55:LEU:N	2.48	0.46
1:B:130:GLU:CD	1:B:426:HIS:ND1	2.73	0.46
1:B:272:GLU:OE2	1:B:276:LEU:HD13	2.15	0.46
1:A:272:GLU:OE1	1:A:272:GLU:CA	2.56	0.46
1:A:492:ILE:CG2	1:A:493:TRP:H	2.28	0.46
1:B:349:PHE:CD1	1:B:350:GLY:N	2.84	0.46
1:A:203:LEU:HD21	1:A:252:GLY:HA2	1.97	0.46
1:A:528:PHE:O	1:A:529:TRP:C	2.58	0.46
1:B:130:GLU:OE1	1:B:426:HIS:ND1	2.49	0.46
1:B:170:PRO:O	1:B:172:TYR:N	2.41	0.46
1:B:212:LEU:HD22	1:B:274:ASN:OD1	2.16	0.46
1:A:91:VAL:HG22	1:A:405:LYS:HG2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:LEU:HD23	1:B:187:LEU:HA	1.66	0.46
1:B:287:TRP:NE1	1:B:291:LYS:HE3	2.31	0.45
1:A:45:ARG:NE	1:A:178:ASP:OD1	2.39	0.45
1:B:253:ASP:OD1	1:B:255:ARG:NH1	2.48	0.45
1:B:358:VAL:HB	1:B:379:LEU:HD11	1.98	0.45
1:A:249:PHE:CZ	1:A:387:ILE:HD11	2.51	0.45
1:A:257:SER:O	1:A:381:PHE:HA	2.16	0.45
1:A:343:PHE:CG	1:A:518:GLN:HG2	2.51	0.45
1:B:93:ASP:HA	1:B:406:LEU:HB2	1.98	0.45
1:A:88:GLU:CD	1:A:356:SER:HB3	2.42	0.45
1:A:376:LEU:C	1:A:376:LEU:CD2	2.89	0.45
1:A:484:TYR:C	1:A:486:THR:H	2.23	0.45
1:B:99:LEU:HD23	1:B:566:ALA:HB1	1.98	0.45
1:B:261:LEU:HD12	1:B:261:LEU:HA	1.83	0.45
1:B:288:ASP:O	1:B:291:LYS:HB3	2.17	0.45
1:B:312:TYR:CD2	1:B:312:TYR:C	2.94	0.45
1:B:480:LEU:HD12	1:B:480:LEU:HA	1.84	0.45
1:B:494:ILE:HG23	1:B:495:GLY:N	2.31	0.45
1:A:520:GLN:HG2	1:A:521:GLN:N	2.29	0.45
1:B:342:VAL:HG13	1:B:343:PHE:N	2.31	0.45
1:B:467:LEU:HD23	1:B:477:ALA:CA	2.46	0.45
1:A:350:GLY:HA3	7:A:605:HEM:CBC	2.47	0.45
1:B:503:GLU:HG2	1:B:504:ARG:HG3	1.98	0.45
1:A:85:TYR:CD1	1:A:87:ASP:O	2.70	0.45
1:A:226:TYR:N	9:A:3063:HOH:O	2.50	0.45
1:A:362:ASP:O	1:A:365:TYR:N	2.47	0.45
1:B:305:GLN:HG2	1:B:529:TRP:CZ3	2.51	0.45
1:B:418:ARG:O	1:B:418:ARG:HG2	2.17	0.45
1:A:106:ILE:HG23	1:A:191:LEU:HD11	1.99	0.45
1:B:2:TRP:C	1:B:4:VAL:N	2.75	0.45
1:B:57:VAL:HA	1:B:58:PRO:HD3	1.72	0.45
1:B:377:HIS:CD2	1:B:378:THR:HG22	2.52	0.45
1:A:88:GLU:OE2	1:A:414:THR:HG23	2.17	0.45
1:A:464:LEU:HD13	1:A:468:GLN:OE1	2.17	0.45
1:B:478:LYS:O	1:B:479:LYS:C	2.60	0.45
1:B:180:ILE:HG22	1:B:181:ASN:N	2.32	0.45
1:B:9:PRO:C	1:B:10:VAL:HG13	2.42	0.44
1:B:58:PRO:HG3	1:B:162:ARG:NH2	2.32	0.44
1:A:492:ILE:HG22	1:A:493:TRP:H	1.83	0.44
1:A:233:LYS:HB3	1:A:234:PRO:CD	2.23	0.44
1:B:165:PHE:CD1	1:B:165:PHE:N	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:TRP:C	1:A:386:ILE:N	2.75	0.44
1:B:7:GLY:O	1:B:8:ALA:CB	2.62	0.44
1:B:91:VAL:C	1:B:92:LEU:HD12	2.43	0.44
1:B:358:VAL:HB	1:B:379:LEU:CD1	2.48	0.44
1:B:102:GLN:OE1	1:B:259:GLN:NE2	2.47	0.44
1:B:306:ILE:O	1:B:310:ARG:HB3	2.17	0.44
1:A:67:ARG:HD3	1:A:489:ASN:OD1	2.18	0.44
1:A:173:GLN:HB2	1:A:174:SER:H	1.38	0.44
1:B:458:LEU:HD21	1:B:509:PRO:HB3	2.00	0.44
1:B:368:ARG:O	1:B:368:ARG:CG	2.66	0.44
1:B:299:ILE:HD11	1:B:590:TRP:CE2	2.52	0.44
1:A:106:ILE:HG23	1:A:191:LEU:CD1	2.48	0.44
1:A:255:ARG:O	1:A:256:ALA:C	2.61	0.44
1:B:53:ASP:C	1:B:55:LEU:N	2.72	0.44
1:A:35:ALA:O	1:A:36:LEU:C	2.60	0.43
1:A:433:LEU:O	1:A:437:ASN:ND2	2.51	0.43
1:A:551:ARG:N	1:A:582:ILE:HD11	2.32	0.43
1:B:62:THR:O	1:B:64:ARG:N	2.51	0.43
1:B:454:GLY:C	1:B:456:CYS:N	2.76	0.43
1:B:541:ARG:O	1:B:544:LEU:N	2.51	0.43
1:A:78:VAL:HG13	1:A:82:ILE:HD12	2.00	0.43
1:B:246:VAL:HG11	1:B:387:ILE:HG13	2.00	0.43
1:B:258:GLU:OE2	7:B:605:HEM:CHB	2.66	0.43
1:B:493:TRP:O	1:B:497:ASN:ND2	2.50	0.43
1:A:25:THR:HG22	1:A:27:ASP:N	2.25	0.43
1:B:251:ALA:O	1:B:253:ASP:N	2.52	0.43
1:B:584:LYS:O	1:B:585:LEU:HB3	2.18	0.43
1:B:588:SER:OG	1:B:589:PRO:CD	2.59	0.43
1:A:194:GLY:HA2	1:A:252:GLY:O	2.18	0.43
1:B:91:VAL:O	1:B:91:VAL:HG13	2.17	0.43
1:A:249:PHE:CE1	1:A:387:ILE:HD11	2.54	0.43
1:A:319:SER:OG	1:A:503:GLU:HB3	2.18	0.43
1:A:428:ILE:HG12	1:A:430:GLY:H	1.83	0.43
1:A:436:ILE:HG12	1:A:436:ILE:H	1.71	0.43
1:B:335:VAL:O	1:B:337:PRO:HD3	2.18	0.43
1:A:260:ILE:HG23	1:A:261:LEU:N	2.32	0.43
1:B:233:LYS:CB	1:B:234:PRO:HD2	2.48	0.43
1:A:85:TYR:CG	1:A:411:LYS:HB3	2.53	0.43
1:A:294:GLN:O	1:A:298:LYS:HB2	2.18	0.43
1:B:14:LYS:O	1:B:15:CYS:HB2	2.18	0.43
1:A:136:GLY:O	1:A:138:ASN:N	2.41	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:556:ASN:O	1:A:557:THR:HG22	2.19	0.43
1:B:191:LEU:H	1:B:191:LEU:CD2	2.32	0.43
1:A:64:ARG:HA	1:A:71:ARG:HH22	1.77	0.43
1:A:130:GLU:C	1:A:131:GLU:HG2	2.44	0.43
1:A:226:TYR:CE1	1:A:387:ILE:CG1	3.01	0.43
1:A:417:LEU:HD11	7:A:605:HEM:HBB2	2.01	0.43
1:B:187:LEU:HD11	1:B:308:THR:HG21	2.00	0.43
1:A:24:ILE:CD1	1:A:200:ALA:HB1	2.35	0.42
1:A:227:PRO:HG2	1:A:249:PHE:CD2	2.53	0.42
1:A:448:GLY:O	1:A:452:TRP:CD1	2.73	0.42
1:B:158:MET:O	1:B:159:PRO:C	2.61	0.42
1:B:454:GLY:O	1:B:455:PHE:C	2.62	0.42
1:B:569:ALA:O	1:B:570:ASN:ND2	2.52	0.42
1:A:108:ASP:CG	7:A:605:HEM:CMD	2.78	0.42
1:B:16:ASP:O	1:B:17:GLU:CB	2.64	0.42
1:B:476:LEU:O	1:B:480:LEU:HB2	2.19	0.42
1:A:39:ALA:O	1:A:40:ASN:HB2	2.18	0.42
1:A:92:LEU:HD22	1:A:404:SER:O	2.19	0.42
1:B:101:MET:HE2	7:B:605:HEM:HMC1	2.01	0.42
1:B:584:LYS:O	1:B:584:LYS:HG3	2.19	0.42
1:A:475:ILE:HD11	1:A:479:LYS:CE	2.46	0.42
1:B:88:GLU:O	1:B:91:VAL:HG12	2.20	0.42
1:B:126:LYS:HE3	1:B:126:LYS:HB2	1.85	0.42
1:B:153:THR:HG22	1:B:154:GLN:HG3	2.02	0.42
1:A:170:PRO:CB	1:A:171:PRO:CD	2.96	0.42
1:B:85:TYR:CD2	1:B:411:LYS:HA	2.55	0.42
1:B:98:LEU:HB3	1:B:399:LEU:HA	2.02	0.42
1:B:99:LEU:HD21	1:B:549:PHE:CD2	2.54	0.42
3:E:2:NAG:H61	3:E:2:NAG:C2	2.47	0.42
1:A:30:ASN:OD1	1:A:30:ASN:C	2.63	0.42
1:A:179:GLN:N	1:A:179:GLN:OE1	2.53	0.42
1:A:283:LEU:C	1:A:285:PRO:HD3	2.45	0.42
1:A:532:ASN:HA	1:A:533:PRO:HD3	1.92	0.42
1:B:346:ALA:HB1	1:B:511:LEU:HD12	2.01	0.42
1:B:351:HIS:CE1	1:B:433:LEU:HD21	2.55	0.42
1:B:139:CYS:C	1:B:141:PRO:HD3	2.45	0.42
1:B:170:PRO:CB	1:B:171:PRO:CD	2.97	0.42
1:B:260:ILE:HG13	1:B:386:ILE:HD11	2.01	0.42
1:A:260:ILE:HD11	1:A:386:ILE:HG12	2.00	0.42
1:A:572:TYR:HA	1:A:573:PRO:HA	1.86	0.42
1:B:51:TYR:CD2	1:B:56:ALA:HA	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:345:PHE:HZ	1:B:444:HIS:CD2	2.38	0.42
1:B:63:GLN:C	1:B:65:LYS:H	2.28	0.41
1:B:251:ALA:O	1:B:252:GLY:C	2.63	0.41
1:B:299:ILE:CD1	1:B:590:TRP:CE2	3.03	0.41
1:A:433:LEU:HG	1:A:437:ASN:HD21	1.85	0.41
1:B:140:PHE:N	1:B:141:PRO:HD3	2.36	0.41
1:B:251:ALA:C	1:B:253:ASP:N	2.78	0.41
1:B:332:ASN:C	1:B:332:ASN:OD1	2.61	0.41
1:B:351:HIS:HD1	1:B:437:ASN:ND2	2.18	0.41
1:B:362:ASP:HB3	1:B:368:ARG:CB	2.40	0.41
1:B:410:ASN:O	1:B:410:ASN:CG	2.63	0.41
1:B:586:ASP:OD1	1:B:586:ASP:C	2.63	0.41
1:A:36:LEU:HA	1:A:36:LEU:HD12	1.79	0.41
1:B:293:TYR:CD2	1:B:294:GLN:HG2	2.55	0.41
1:B:571:ASN:O	1:B:576:PHE:HE1	2.03	0.41
1:A:118:GLU:HG3	1:A:119:LEU:H	1.85	0.41
1:A:502:VAL:HG13	1:A:507:VAL:O	2.20	0.41
1:A:535:VAL:HG12	1:A:536:PHE:CD1	2.55	0.41
1:B:472:LYS:HE3	1:B:472:LYS:HB2	1.76	0.41
1:A:99:LEU:HA	1:A:399:LEU:HD22	2.02	0.41
1:A:349:PHE:HA	1:A:497:ASN:HD21	1.85	0.41
1:A:426:HIS:CD2	1:A:426:HIS:N	2.88	0.41
1:A:578:ASP:C	1:A:580:SER:N	2.78	0.41
1:B:233:LYS:HB3	1:B:234:PRO:CD	2.50	0.41
1:A:83:VAL:O	1:A:412:MET:HB2	2.20	0.41
1:A:260:ILE:CD1	1:A:386:ILE:CG1	2.92	0.41
1:A:484:TYR:CD2	1:A:490:ILE:HG13	2.55	0.41
1:B:108:ASP:OD2	7:B:605:HEM:CHD	2.69	0.41
1:B:370:PRO:C	1:B:371:GLU:HG3	2.45	0.41
1:B:475:ILE:HD11	1:B:479:LYS:HZ1	1.83	0.41
1:A:95:ASN:HA	1:A:569:ALA:HB2	2.03	0.41
1:A:229:PHE:HE1	1:A:387:ILE:CD1	2.34	0.41
1:A:233:LYS:HD2	1:A:233:LYS:HA	1.84	0.41
1:A:551:ARG:HA	1:A:582:ILE:HD11	2.02	0.41
1:B:264:THR:O	1:B:267:THR:HB	2.20	0.41
1:A:165:PHE:CG	1:A:177:ARG:HD2	2.55	0.41
1:A:450:ASN:C	1:A:452:TRP:N	2.79	0.41
1:A:561:LYS:HA	1:A:577:VAL:O	2.20	0.41
1:B:85:TYR:CG	1:B:411:LYS:HG2	2.55	0.41
1:A:120:GLY:O	1:A:121:SER:HB2	2.21	0.41
1:A:140:PHE:O	1:A:160:PHE:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:ILE:HD12	1:A:386:ILE:HG13	2.02	0.41
1:A:352:MET:SD	1:A:412:MET:HG2	2.61	0.41
1:A:400:LEU:HD13	1:A:563:PRO:CD	2.50	0.41
1:A:544:LEU:O	1:A:547:VAL:HG22	2.20	0.41
1:B:19:SER:HA	1:B:20:PRO:HD3	1.76	0.41
1:B:123:GLU:HG3	1:B:125:SER:H	1.85	0.41
1:B:511:LEU:HD13	1:B:511:LEU:HA	1.82	0.41
3:E:2:NAG:H2	3:E:2:NAG:C6	2.50	0.41
1:A:549:PHE:O	1:A:553:ILE:HG12	2.21	0.41
1:A:572:TYR:CE2	1:A:573:PRO:HB3	2.56	0.41
1:B:124:HIS:C	1:B:127:VAL:H	2.29	0.41
1:B:150:LYS:HG2	1:B:154:GLN:OE1	2.21	0.41
1:B:438:LEU:HD23	1:B:438:LEU:HA	1.77	0.41
1:A:19:SER:C	1:A:21:TYR:H	2.29	0.40
1:A:258:GLU:O	1:A:258:GLU:HG2	2.21	0.40
1:A:348:ARG:NH1	7:A:605:HEM:O2A	2.48	0.40
1:A:414:THR:HG22	1:A:416:GLU:H	1.86	0.40
1:A:568:GLN:H	1:A:568:GLN:HG2	1.69	0.40
1:B:172:TYR:H	1:B:172:TYR:HD1	1.66	0.40
1:B:542:ASP:O	1:B:545:GLN:HG3	2.20	0.40
1:A:260:ILE:CG2	1:A:261:LEU:N	2.84	0.40
1:A:384:TRP:C	1:A:386:ILE:H	2.29	0.40
1:A:434:ALA:HA	1:A:437:ASN:HD22	1.86	0.40
1:A:464:LEU:CD1	1:A:468:GLN:OE1	2.69	0.40
1:A:537:THR:O	1:A:541:ARG:HG3	2.21	0.40
1:A:576:PHE:N	1:A:576:PHE:CD1	2.89	0.40
1:B:134:ILE:O	1:B:134:ILE:HG22	2.22	0.40
1:B:160:PHE:HE2	1:B:436:ILE:HG23	1.86	0.40
3:E:2:NAG:C2	3:E:2:NAG:C6	2.99	0.40
1:A:545:GLN:HE21	1:A:545:GLN:HB2	1.67	0.40
1:A:85:TYR:CE2	1:A:411:LYS:HB3	2.56	0.40
1:B:46:TRP:NE1	1:B:340:SER:HB3	2.36	0.40
1:A:216:ASN:HB2	1:A:227:PRO:O	2.21	0.40
1:B:95:ASN:HA	1:B:569:ALA:CB	2.52	0.40
1:B:95:ASN:HA	1:B:569:ALA:HB2	2.03	0.40
1:B:541:ARG:O	1:B:542:ASP:C	2.64	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 X-ray entries followed by that with respect to entries of similar resolution.

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	593/595 (100%)	494 (83%)	75 (13%)	24 (4%)	2	14
1	B	593/595 (100%)	492 (83%)	70 (12%)	31 (5%)	1	10
All	All	1186/1190 (100%)	986 (83%)	145 (12%)	55 (5%)	2	11

All (55) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	117	THR
1	A	137	ASP
1	A	169	THR
1	A	174	SER
1	A	234	PRO
1	A	256	ALA
1	A	409	GLN
1	A	594	GLU
1	B	4	VAL
1	B	9	PRO
1	B	17	GLU
1	B	63	GLN
1	B	123	GLU
1	B	167	CYS
1	B	168	PRO
1	B	169	THR
1	B	171	PRO
1	B	172	TYR
1	B	174	SER
1	B	234	PRO
1	B	363	GLU
1	A	170	PRO
1	A	232	MET
1	A	363	GLU

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Mol	Chain	Res	Type
1	A	504	ARG
1	A	583	ASP
1	B	5	GLY
1	B	8	ALA
1	B	64	ARG
1	B	211	GLY
1	B	231	ASN
1	B	232	MET
1	B	455	PHE
1	B	594	GLU
1	A	9	PRO
1	A	168	PRO
1	B	10	VAL
1	B	54	GLY
1	B	170	PRO
1	B	333	ASN
1	B	473	ASN
1	A	63	GLN
1	A	566	ALA
1	B	411	LYS
1	A	17	GLU
1	A	121	SER
1	A	167	CYS
1	A	542	ASP
1	B	34	PRO
1	B	173	GLN
1	B	252	GLY
1	A	166	VAL
1	A	34	PRO
1	B	260	ILE

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 X-ray entries followed by that with respect to entries of similar resolution.

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	518/518 (100%)	422 (82%)	96 (18%)	1 7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	518/518 (100%)	445 (86%)	73 (14%)	3 15
All	All	1036/1036 (100%)	867 (84%)	169 (16%)	2 10

All (169) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	3	GLU
1	A	4	VAL
1	A	9	PRO
1	A	12	LEU
1	A	13	VAL
1	A	24	ILE
1	A	25	THR
1	A	32	ARG
1	A	48	PRO
1	A	61	TRP
1	A	71	ARG
1	A	72	VAL
1	A	73	PRO
1	A	79	SER
1	A	92	LEU
1	A	98	LEU
1	A	106	ILE
1	A	115	PRO
1	A	116	GLU
1	A	117	THR
1	A	118	GLU
1	A	119	LEU
1	A	124	HIS
1	A	132	TYR
1	A	134	ILE
1	A	135	GLN
1	A	137	ASP
1	A	138	ASN
1	A	139	CYS
1	A	141	PRO
1	A	143	MET
1	A	149	PRO
1	A	151	LEU
1	A	159	PRO
1	A	169	THR

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Mol	Chain	Res	Type
1	A	173	GLN
1	A	174	SER
1	A	177	ARG
1	A	187	LEU
1	A	197	PRO
1	A	202	ARG
1	A	203	LEU
1	A	207	SER
1	A	208	SER
1	A	209	PRO
1	A	226	TYR
1	A	233	LYS
1	A	234	PRO
1	A	255	ARG
1	A	257	SER
1	A	260	ILE
1	A	261	LEU
1	A	266	HIS
1	A	269	LEU
1	A	276	LEU
1	A	292	LEU
1	A	294	GLN
1	A	298	LYS
1	A	317	LEU
1	A	321	MET
1	A	337	PRO
1	A	359	SER
1	A	360	ARG
1	A	363	GLU
1	A	368	ARG
1	A	373	GLU
1	A	376	LEU
1	A	392	ILE
1	A	405	LYS
1	A	420	LYS
1	A	428	ILE
1	A	447	PRO
1	A	451	SER
1	A	458	LEU
1	A	464	LEU
1	A	471	LEU
1	A	472	LYS

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Mol	Chain	Res	Type
1	A	475	ILE
1	A	478	LYS
1	A	480	LEU
1	A	482	ASP
1	A	483	LEU
1	A	490	ILE
1	A	492	ILE
1	A	507	VAL
1	A	520	GLN
1	A	521	GLN
1	A	522	ILE
1	A	523	ARG
1	A	538	GLU
1	A	542	ASP
1	A	564	LEU
1	A	573	PRO
1	A	582	ILE
1	A	592	SER
1	B	3	GLU
1	B	10	VAL
1	B	11	PRO
1	B	13	VAL
1	B	16	ASP
1	B	17	GLU
1	B	18	ASN
1	B	25	THR
1	B	33	SER
1	B	57	VAL
1	B	64	ARG
1	B	98	LEU
1	B	119	LEU
1	B	124	HIS
1	B	134	ILE
1	B	137	ASP
1	B	150	LYS
1	B	165	PHE
1	B	169	THR
1	B	170	PRO
1	B	172	TYR
1	B	175	LEU
1	B	177	ARG
1	B	179	GLN

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Mol	Chain	Res	Type
1	B	185	SER
1	B	191	LEU
1	B	198	SER
1	B	207	SER
1	B	210	LEU
1	B	215	VAL
1	B	216	ASN
1	B	232	MET
1	B	234	PRO
1	B	242	THR
1	B	246	VAL
1	B	255	ARG
1	B	261	LEU
1	B	265	VAL
1	B	278	ARG
1	B	288	ASP
1	B	304	ILE
1	B	314	PRO
1	B	317	LEU
1	B	329	GLN
1	B	333	ASN
1	B	359	SER
1	B	360	ARG
1	B	373	GLU
1	B	375	PRO
1	B	376	LEU
1	B	377	HIS
1	B	389	ASP
1	B	392	ILE
1	B	404	SER
1	B	410	ASN
1	B	451	SER
1	B	462	LYS
1	B	464	LEU
1	B	471	LEU
1	B	474	LYS
1	B	475	ILE
1	B	494	ILE
1	B	511	LEU
1	B	520	GLN
1	B	523	ARG
1	B	537	THR

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Mol	Chain	Res	Type
1	B	545	GLN
1	B	547	VAL
1	B	550	SER
1	B	561	LYS
1	B	573	PRO
1	B	582	ILE
1	B	595	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	94	GLN
1	A	135	GLN
1	A	147	ASN
1	A	230	ASN
1	A	250	GLN
1	A	322	GLN
1	A	364	ASN
1	A	377	HIS
1	A	450	ASN
1	A	460	GLN
1	A	497	ASN
1	A	520	GLN
1	A	545	GLN
1	A	565	HIS
1	A	570	ASN
1	B	18	ASN
1	B	40	ASN
1	B	63	GLN
1	B	138	ASN
1	B	217	GLN
1	B	322	GLN
1	B	329	GLN
1	B	333	ASN
1	B	341	ASN
1	B	410	ASN
1	B	439	GLN
1	B	460	GLN
1	B	468	GLN
1	B	521	GLN
1	B	595	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

14 monosaccharides 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
2	NAG	C	1	1,2	14,14,15	1.14	1 (7%)	17,19,21	2.66	5 (29%)
2	NAG	C	2	2	14,14,15	1.28	1 (7%)	17,19,21	2.03	5 (29%)
2	NAG	D	1	1,2	14,14,15	0.67	0	17,19,21	0.96	1 (5%)
2	NAG	D	2	2	14,14,15	0.78	0	17,19,21	0.78	0
3	NAG	E	1	1,3	14,14,15	0.55	0	17,19,21	0.85	0
3	NAG	E	2	3	14,14,15	0.88	0	17,19,21	0.87	1 (5%)
3	MAN	E	3	3	11,11,12	0.65	0	15,15,17	0.36	0
2	NAG	F	1	1,2	14,14,15	0.66	0	17,19,21	0.72	0
2	NAG	F	2	2	14,14,15	0.66	0	17,19,21	1.51	2 (11%)
2	NAG	G	1	1,2	14,14,15	0.67	0	17,19,21	0.88	1 (5%)
2	NAG	G	2	2	14,14,15	0.88	1 (7%)	17,19,21	0.82	0
3	NAG	H	1	1,3	14,14,15	0.77	1 (7%)	17,19,21	1.37	2 (11%)
3	NAG	H	2	3	14,14,15	0.84	0	17,19,21	1.61	3 (17%)
3	MAN	H	3	3	11,11,12	0.82	0	15,15,17	1.60	2 (13%)

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
2	NAG	C	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
3	NAG	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	1/6/23/26	0/1/1/1
3	MAN	E	3	3	-	0/2/19/22	1/1/1/1
2	NAG	F	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	H	2	3	-	3/6/23/26	0/1/1/1
3	MAN	H	3	3	-	0/2/19/22	1/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	NAG	C3-C2	2.83	1.58	1.52
2	C	1	NAG	C1-C2	2.58	1.55	1.52
2	G	2	NAG	C1-C2	2.52	1.55	1.52
3	H	1	NAG	C4-C5	2.20	1.57	1.53

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C2-N2-C7	-5.72	115.23	122.90
2	C	1	NAG	C4-C3-C2	-5.27	103.30	111.02
2	C	1	NAG	C1-C2-N2	5.23	118.67	110.43
2	C	2	NAG	C3-C4-C5	5.22	119.69	110.23
3	H	3	MAN	C1-O5-C5	5.19	119.15	112.19
3	H	2	NAG	C2-N2-C7	-4.62	116.71	122.90
2	F	2	NAG	C4-C3-C2	4.33	117.36	111.02
2	C	2	NAG	O5-C1-C2	-3.33	106.13	111.29
2	C	1	NAG	C3-C4-C5	-3.25	104.34	110.23
2	F	2	NAG	C3-C4-C5	3.23	116.10	110.23
3	H	2	NAG	C6-C5-C4	2.77	119.83	113.02
3	H	1	NAG	C2-N2-C7	-2.73	119.24	122.90
3	H	3	MAN	C1-C2-C3	2.72	113.61	109.64
3	H	2	NAG	C1-O5-C5	2.58	115.64	112.19
2	C	2	NAG	O5-C5-C6	-2.46	102.87	107.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1	NAG	O4-C4-C3	-2.37	104.80	110.38
2	C	2	NAG	C4-C3-C2	2.30	114.39	111.02
2	C	1	NAG	C8-C7-N2	2.24	119.83	116.12
3	E	2	NAG	C1-O5-C5	2.14	115.05	112.19
2	D	1	NAG	C2-N2-C7	-2.12	120.06	122.90
2	C	2	NAG	C2-N2-C7	-2.09	120.10	122.90
2	G	1	NAG	C2-N2-C7	-2.06	120.13	122.90

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	H	2	NAG	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6
2	D	1	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	C	1	NAG	C3-C2-N2-C7
2	F	1	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	H	1	NAG	C1-C2-N2-C7
3	H	2	NAG	C1-C2-N2-C7

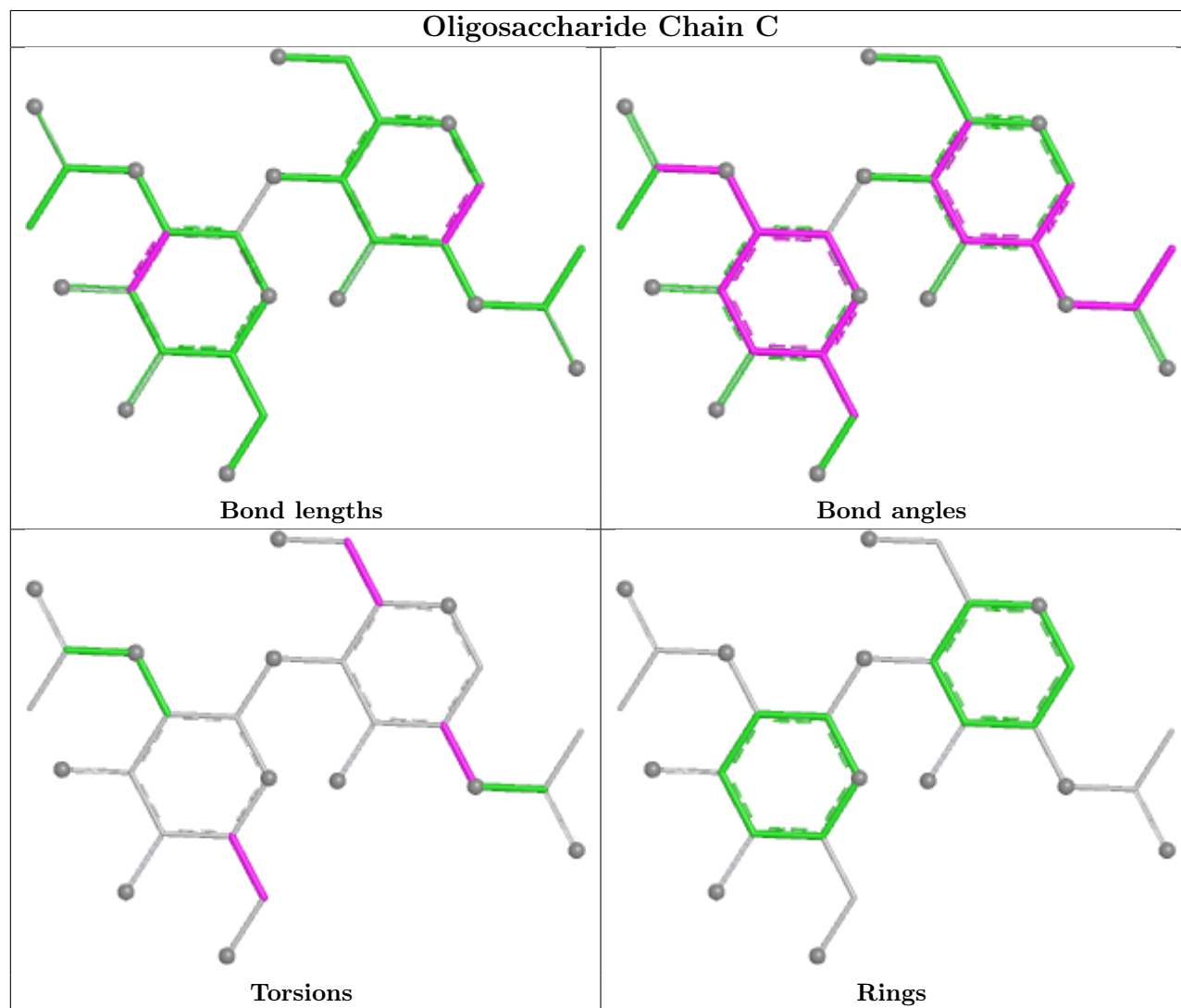
All (2) ring outliers are listed below:

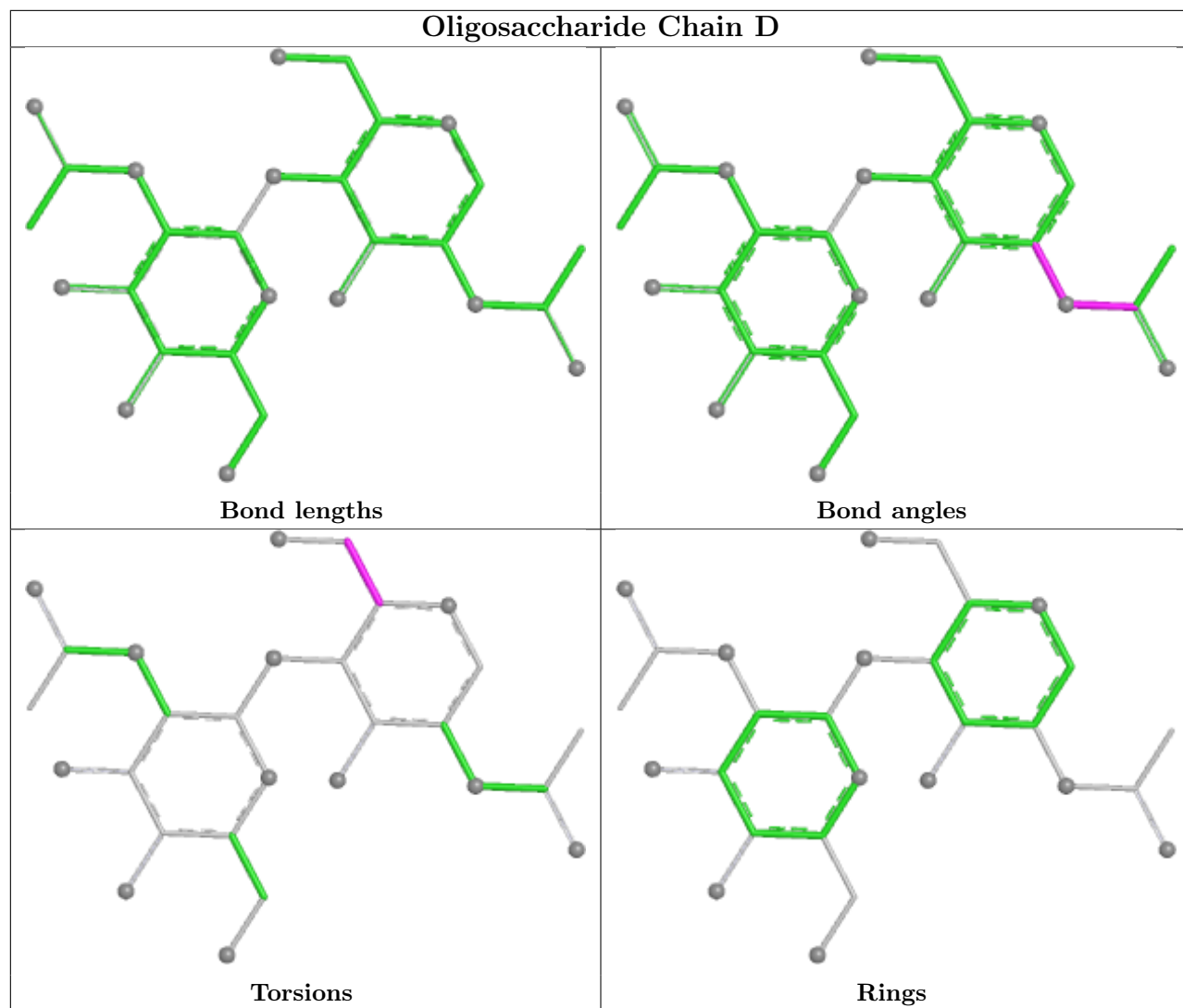
Mol	Chain	Res	Type	Atoms
3	H	3	MAN	C1-C2-C3-C4-C5-O5
3	E	3	MAN	C1-C2-C3-C4-C5-O5

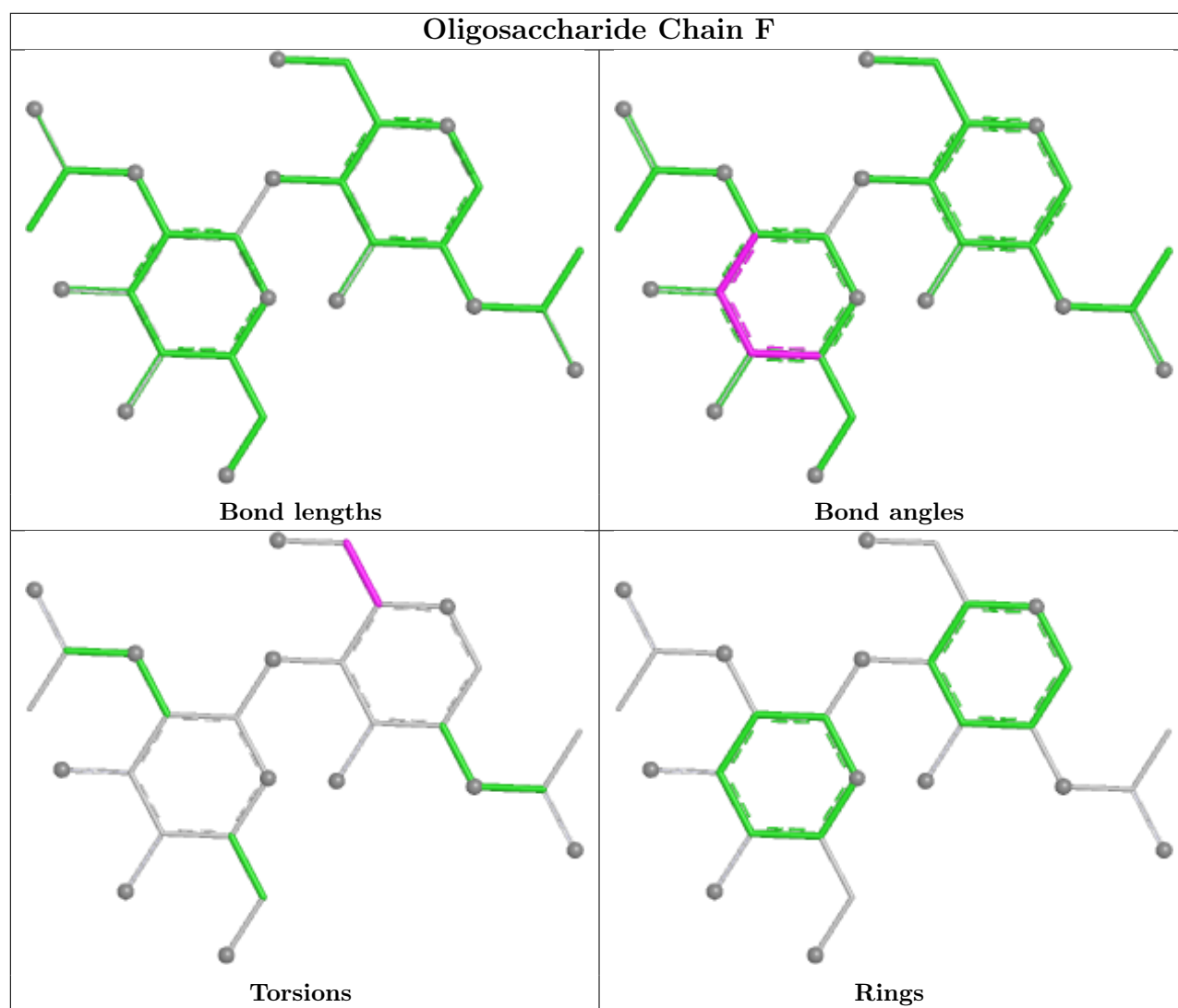
7 monomers are involved in 14 short contacts:

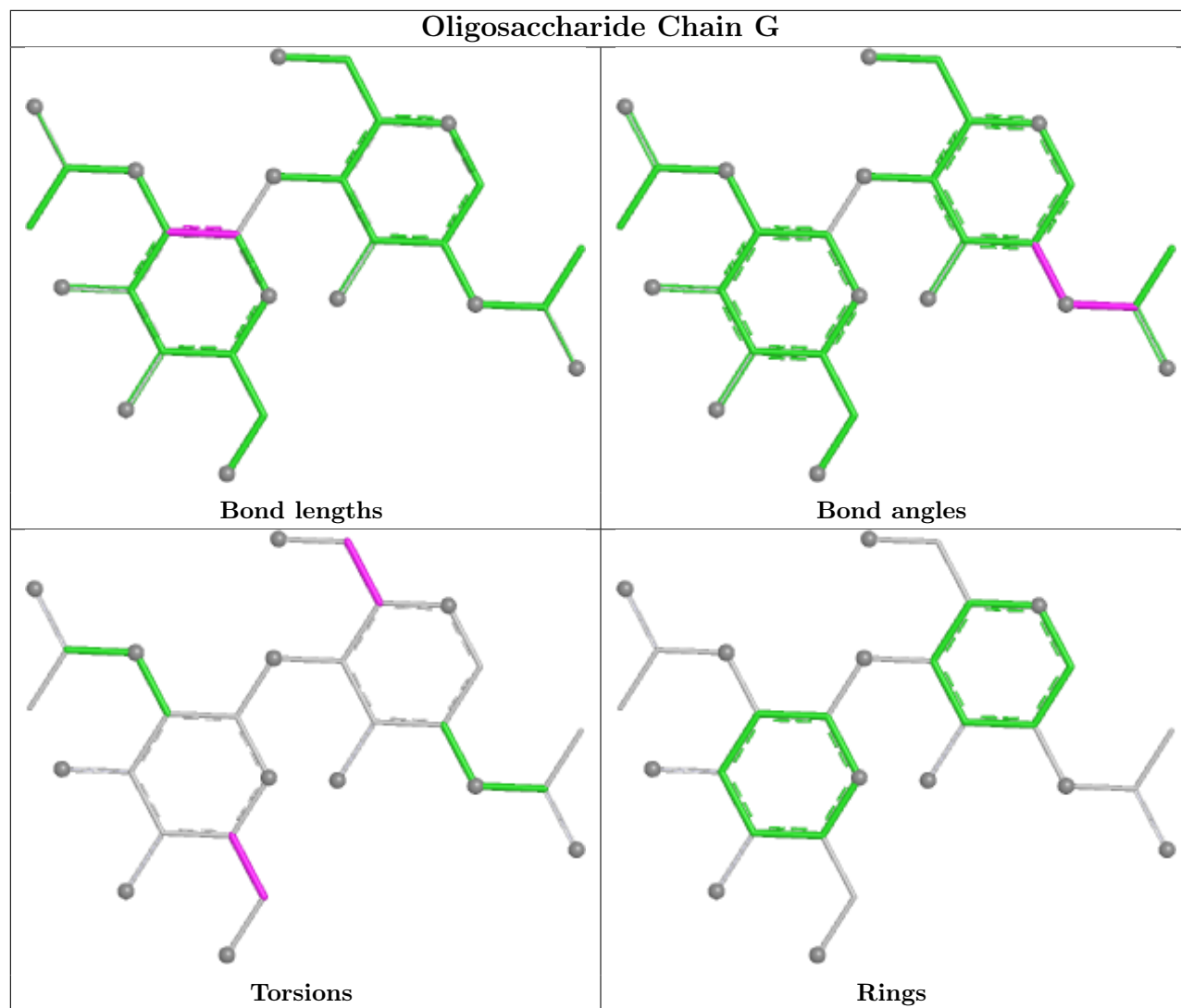
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1	NAG	2	0
3	E	2	NAG	6	0
3	E	3	MAN	1	0
2	C	1	NAG	2	0
2	D	1	NAG	2	0
2	C	2	NAG	2	0
3	H	3	MAN	1	0

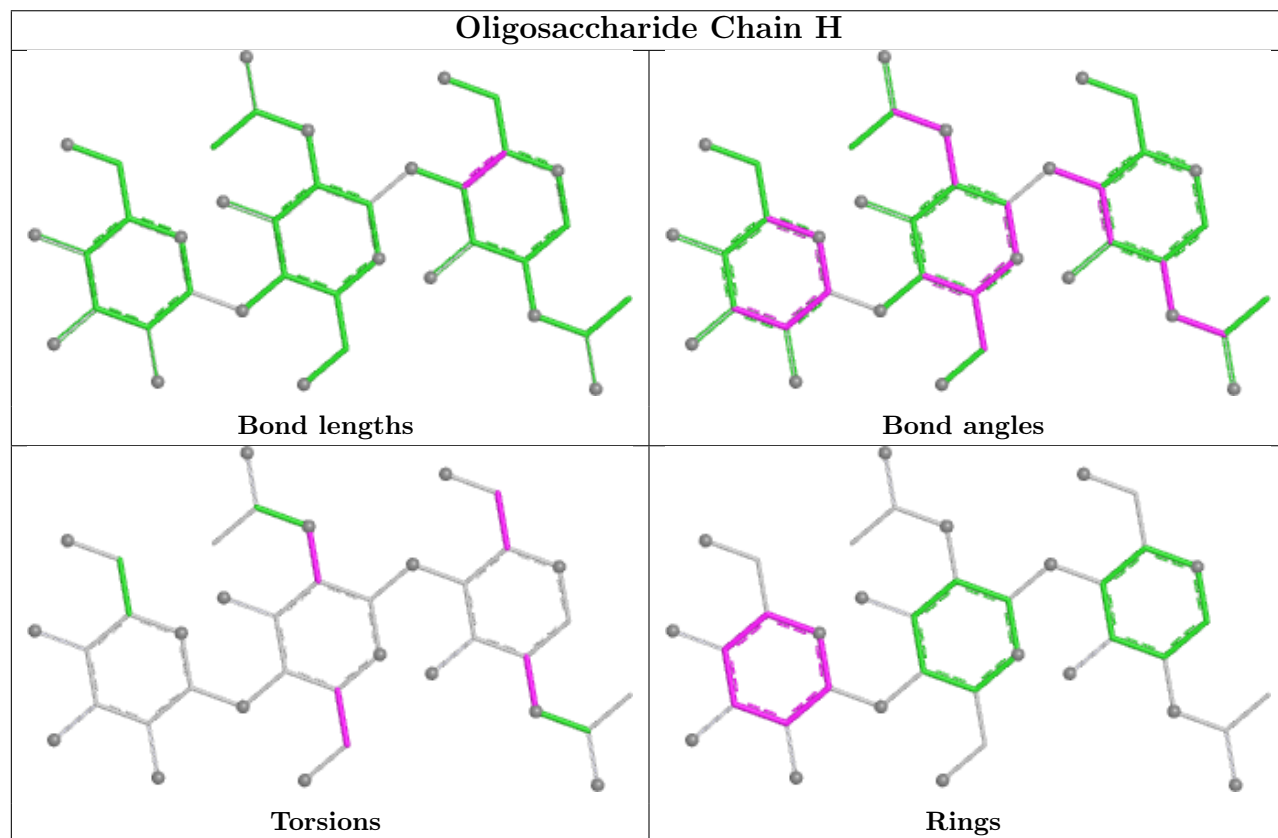
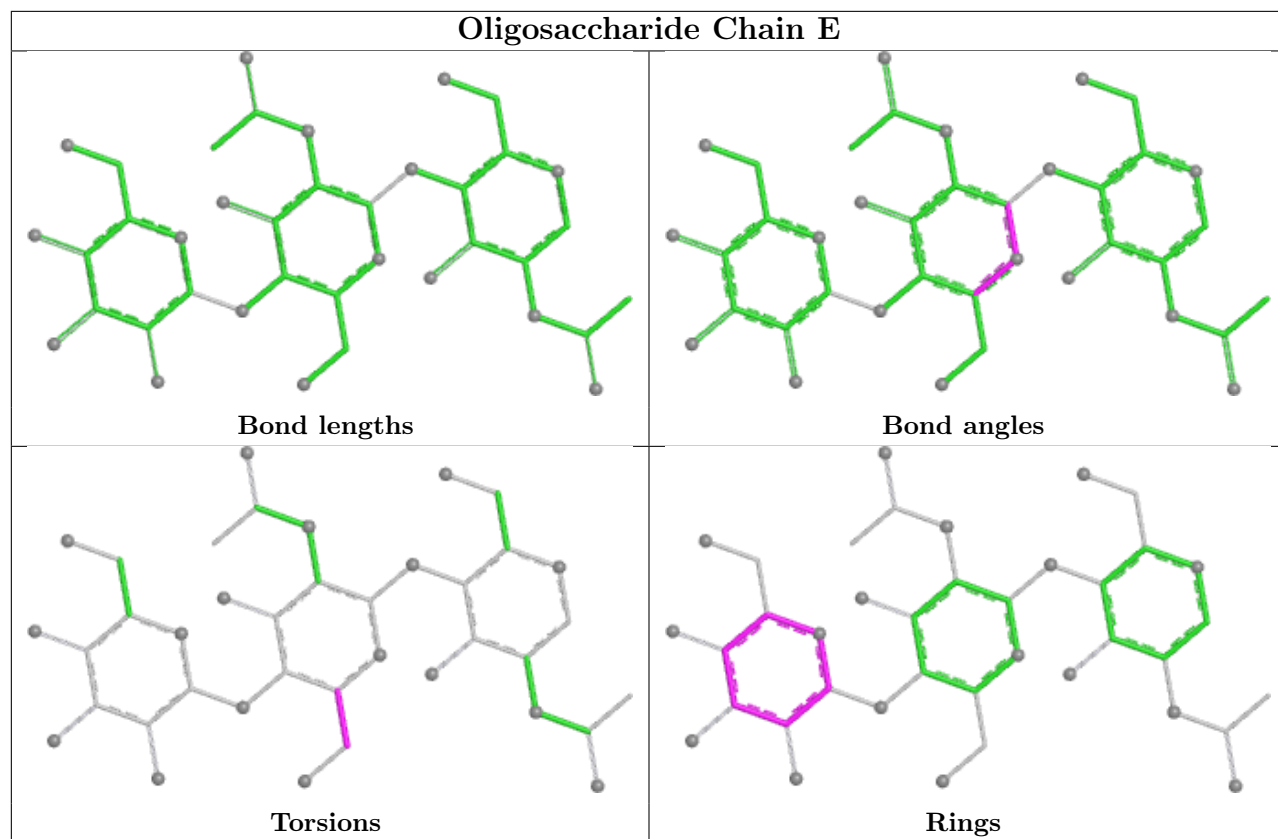
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry

Of 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 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
8	FMT	A	3001	-	2,2,2	0.85	0	1,1,1	0.57	0
6	CO3	B	2002	-	3,3,3	0.88	0	2,3,3	0.16	0
4	NAG	B	603	1	14,14,15	0.83	1 (7%)	17,19,21	1.40	2 (11%)
6	CO3	A	2001	-	3,3,3	0.86	0	2,3,3	0.17	0
7	HEM	A	605	1	50,50,50	2.50	15 (30%)	67,82,82	2.32	26 (38%)
8	FMT	B	3006	-	2,2,2	0.97	0	1,1,1	0.52	0
8	FMT	A	3002	-	2,2,2	0.87	0	1,1,1	0.59	0
7	HEM	B	605	1	50,50,50	2.53	16 (32%)	67,82,82	2.34	26 (38%)
4	NAG	A	603	1	14,14,15	0.81	0	17,19,21	1.51	2 (11%)
8	FMT	A	3003	-	2,2,2	1.13	0	1,1,1	0.50	0
8	FMT	B	3005	-	2,2,2	0.76	0	1,1,1	0.63	0
8	FMT	B	3004	-	2,2,2	0.90	0	1,1,1	0.56	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
7	HEM	B	605	1	-	7/14/54/54	-
4	NAG	B	603	1	-	0/6/23/26	0/1/1/1
4	NAG	A	603	1	-	0/6/23/26	0/1/1/1
7	HEM	A	605	1	-	6/14/54/54	-

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	605	HEM	FE-NA	8.32	2.22	1.95
7	A	605	HEM	C3D-C2D	7.89	1.53	1.36
7	B	605	HEM	FE-NA	7.89	2.21	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	605	HEM	C3D-C2D	7.80	1.53	1.36
7	B	605	HEM	FE-NC	7.56	2.20	1.95
7	A	605	HEM	FE-NC	6.25	2.15	1.95
7	A	605	HEM	FE-NB	4.85	2.09	1.94
7	B	605	HEM	CAC-C3C	3.99	1.58	1.47
7	A	605	HEM	CAC-C3C	3.55	1.56	1.47
7	B	605	HEM	CAB-C3B	3.49	1.56	1.47
7	A	605	HEM	CMC-C2C	3.47	1.57	1.50
7	A	605	HEM	C1A-NA	-3.13	1.33	1.39
7	B	605	HEM	C1A-NA	-3.08	1.33	1.39
7	A	605	HEM	CAB-C3B	3.08	1.55	1.47
7	B	605	HEM	CMC-C2C	3.07	1.57	1.50
7	B	605	HEM	FE-NB	2.90	2.03	1.94
7	B	605	HEM	FE-ND	2.86	2.03	1.94
7	B	605	HEM	C1C-C2C	2.78	1.50	1.45
7	B	605	HEM	CAD-C3D	2.69	1.58	1.51
7	A	605	HEM	CHB-C1B	2.68	1.43	1.38
7	B	605	HEM	CHB-C1B	2.66	1.43	1.38
7	A	605	HEM	C1C-C2C	2.51	1.50	1.45
7	B	605	HEM	CBA-CGA	2.46	1.56	1.50
7	A	605	HEM	O1A-CGA	2.35	1.29	1.22
7	A	605	HEM	CHC-C1C	2.23	1.42	1.38
7	B	605	HEM	CHC-C1C	2.15	1.42	1.38
7	A	605	HEM	CAD-C3D	2.14	1.56	1.51
7	A	605	HEM	CMA-C3A	2.14	1.55	1.50
7	B	605	HEM	O1A-CGA	2.07	1.28	1.22
4	B	603	NAG	C4-C3	2.06	1.57	1.52
7	A	605	HEM	C2A-C3A	-2.02	1.33	1.38
7	B	605	HEM	CAA-C2A	2.01	1.56	1.51

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	605	HEM	C3B-C4B-NB	-6.91	104.51	109.47
7	A	605	HEM	C3B-C4B-NB	-6.69	104.67	109.47
7	B	605	HEM	C1B-NB-C4B	5.43	111.63	105.21
7	A	605	HEM	C1B-NB-C4B	5.31	111.49	105.21
4	A	603	NAG	C4-C3-C2	4.80	118.05	111.02
7	B	605	HEM	CHD-C1D-ND	4.68	129.46	124.42
7	B	605	HEM	CHC-C4B-NB	4.53	129.29	124.42
7	A	605	HEM	C2B-C1B-NB	-4.46	104.71	109.84
7	A	605	HEM	C4D-ND-C1D	4.41	110.43	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	605	HEM	C2B-C1B-NB	-4.33	104.86	109.84
4	B	603	NAG	C4-C3-C2	4.24	117.23	111.02
7	A	605	HEM	CHD-C1D-ND	4.24	128.99	124.42
7	B	605	HEM	C4D-ND-C1D	4.19	110.17	105.21
7	A	605	HEM	CHC-C4B-NB	4.00	128.73	124.42
7	B	605	HEM	CMA-C3A-C4A	3.88	131.32	125.42
7	A	605	HEM	CMA-C3A-C4A	3.77	131.16	125.42
7	A	605	HEM	CHB-C1B-NB	3.72	128.97	124.37
7	B	605	HEM	CHB-C1B-NB	3.63	128.85	124.37
7	A	605	HEM	CAD-CBD-CGD	-3.61	104.09	113.67
7	A	605	HEM	C1A-CHA-C4D	3.43	134.32	126.25
7	B	605	HEM	C1A-CHA-C4D	3.36	134.15	126.25
7	A	605	HEM	C3B-C2B-C1B	3.35	108.92	106.41
7	B	605	HEM	C3B-C2B-C1B	3.31	108.90	106.41
7	A	605	HEM	C4B-C3B-C2B	3.18	110.20	107.28
7	B	605	HEM	CAD-CBD-CGD	-3.17	105.25	113.67
7	B	605	HEM	CAD-C3D-C4D	3.11	130.12	124.70
7	B	605	HEM	C4C-CHD-C1D	-3.05	119.53	126.02
7	A	605	HEM	CBA-CAA-C2A	-3.05	104.10	112.53
7	B	605	HEM	C4B-C3B-C2B	3.03	110.07	107.28
7	B	605	HEM	C1C-CHC-C4B	2.93	132.26	126.02
4	A	603	NAG	C3-C4-C5	2.83	115.36	110.23
7	B	605	HEM	C4C-C3C-C2C	-2.82	104.37	106.81
7	A	605	HEM	CHB-C4A-NA	-2.74	118.89	123.86
7	B	605	HEM	CHA-C4D-ND	2.73	127.75	124.37
7	B	605	HEM	CHB-C4A-NA	-2.72	118.93	123.86
7	A	605	HEM	CBB-CAB-C3B	-2.68	114.15	127.53
7	A	605	HEM	CMB-C2B-C3B	-2.66	122.00	128.43
7	B	605	HEM	CHD-C1D-C2D	-2.62	120.90	125.03
7	A	605	HEM	C4C-CHD-C1D	-2.59	120.51	126.02
7	A	605	HEM	CMB-C2B-C1B	2.59	129.08	125.03
7	A	605	HEM	CAD-C3D-C4D	2.57	129.18	124.70
7	A	605	HEM	C4C-C3C-C2C	-2.57	104.59	106.81
7	A	605	HEM	CHD-C1D-C2D	-2.52	121.05	125.03
7	B	605	HEM	C3C-C2C-C1C	2.48	109.39	107.05
7	A	605	HEM	C1C-CHC-C4B	2.39	131.10	126.02
7	B	605	HEM	CMB-C2B-C3B	-2.30	122.86	128.43
7	A	605	HEM	C3C-C2C-C1C	2.27	109.19	107.05
4	B	603	NAG	C3-C4-C5	2.24	114.30	110.23
7	A	605	HEM	CAB-C3B-C2B	-2.19	121.31	128.43
7	B	605	HEM	CBA-CAA-C2A	-2.13	106.64	112.53
7	A	605	HEM	O2D-CGD-CBD	2.13	120.72	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	605	HEM	CBB-CAB-C3B	-2.08	117.15	127.53
7	B	605	HEM	CMB-C2B-C1B	2.06	128.25	125.03
7	B	605	HEM	O2D-CGD-CBD	2.02	120.39	114.00
7	B	605	HEM	C4D-C3D-C2D	-2.01	103.97	106.89
7	A	605	HEM	C4D-C3D-C2D	-2.00	103.97	106.89

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	605	HEM	C2B-C3B-CAB-CBB
7	A	605	HEM	C4B-C3B-CAB-CBB
7	B	605	HEM	C2B-C3B-CAB-CBB
7	B	605	HEM	C4B-C3B-CAB-CBB
7	B	605	HEM	CAD-CBD-CGD-O1D
7	A	605	HEM	CAD-CBD-CGD-O1D
7	B	605	HEM	CAD-CBD-CGD-O2D
7	A	605	HEM	CAD-CBD-CGD-O2D
7	B	605	HEM	C2D-C3D-CAD-CBD
7	B	605	HEM	CAA-CBA-CGA-O2A
7	B	605	HEM	CAA-CBA-CGA-O1A
7	A	605	HEM	CAA-CBA-CGA-O2A
7	A	605	HEM	CAA-CBA-CGA-O1A

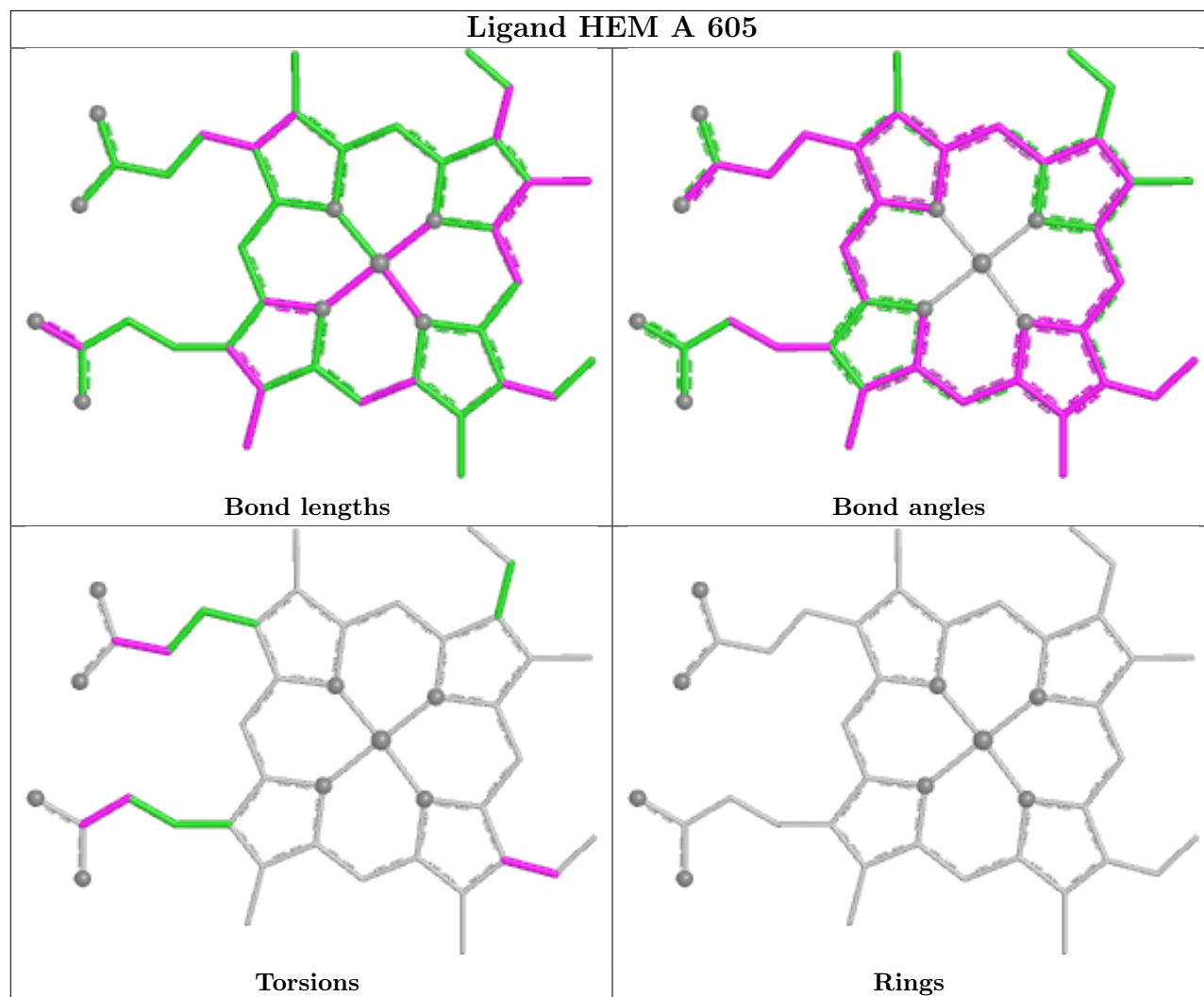
There are no ring outliers.

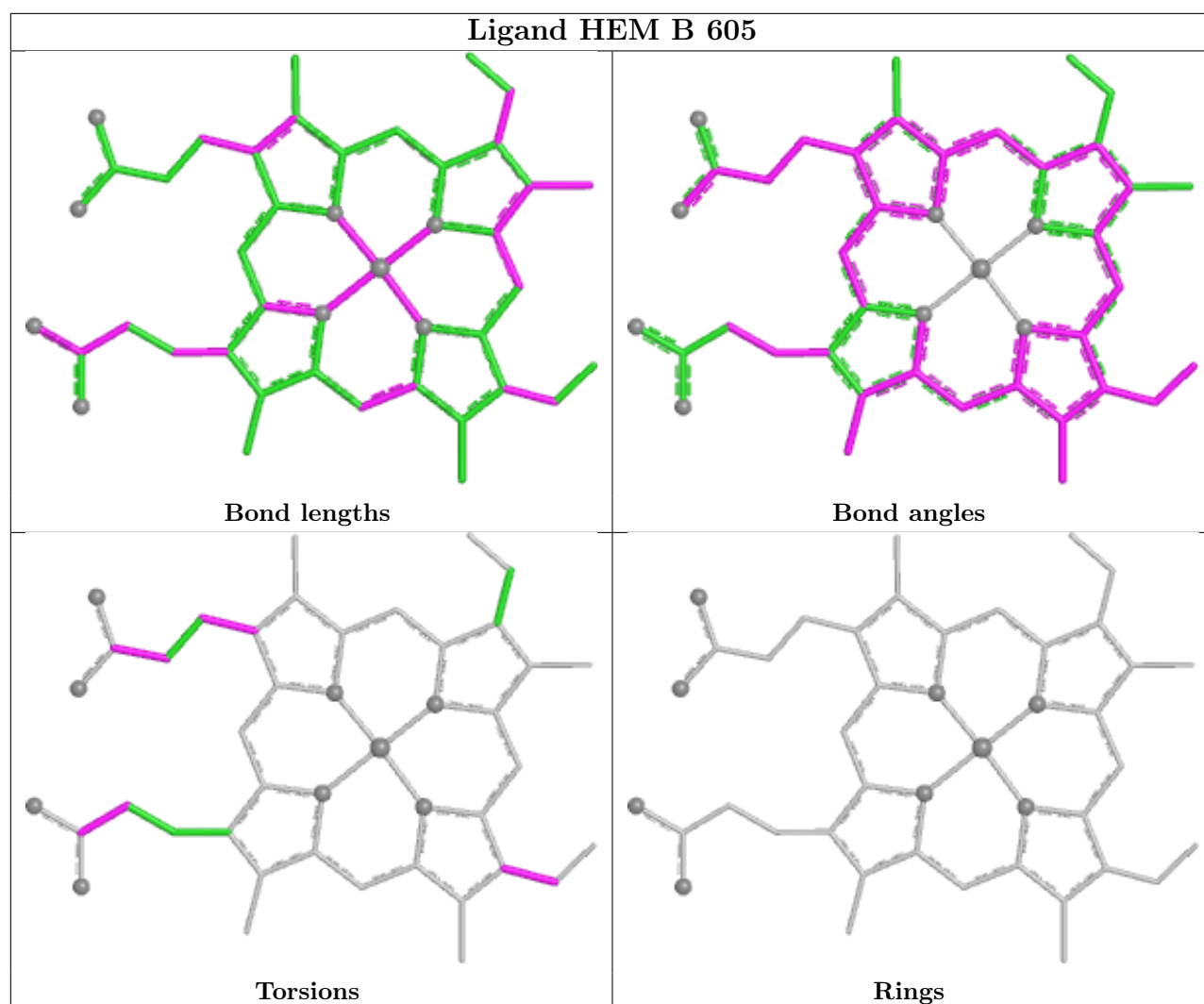
4 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	605	HEM	10	0
8	A	3002	FMT	2	0
7	B	605	HEM	12	0
8	B	3004	FMT	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 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.