



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 02:49 PM UTC

PDB ID : 8FPF / pdb_00008fpf
EMDB ID : EMD-29362
Title : Heterodimeric ABC transporter BmrCD in the inward-facing conformation bound to ATP: BmrCD_IF-ATP
Authors : Tang, Q.; Mchaourab, H.
Deposited on : 2023-01-04
Resolution : 3.27 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.dev132
Mogul : 2022.3.0, CSD as543be (2022)
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) : 2.49

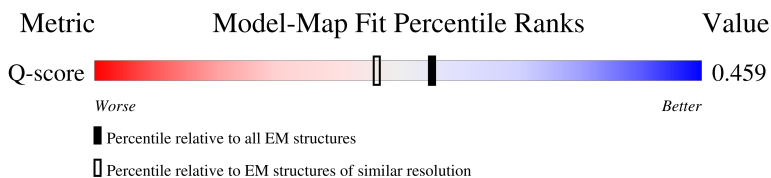
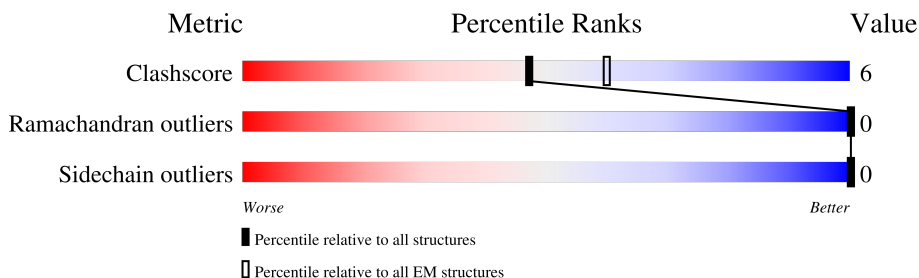
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.27 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	14508 (2.77 - 3.77)

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	D	681	 79% 18%
2	C	607	 82% 9% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21749 atoms, of which 11125 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 Probable multidrug resistance ABC transporter ATP-binding/permease protein YheH.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	D	666	10708	3423	5387	907	974	17	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	154	ALA	CYS	engineered mutation	UNP O07549
D	256	ALA	CYS	engineered mutation	UNP O07549
D	351	ALA	CYS	engineered mutation	UNP O07549
D	592	GLN	GLU	engineered mutation	UNP O07549
D	674	LEU	-	expression tag	UNP O07549
D	675	GLU	-	expression tag	UNP O07549
D	676	HIS	-	expression tag	UNP O07549
D	677	HIS	-	expression tag	UNP O07549
D	678	HIS	-	expression tag	UNP O07549
D	679	HIS	-	expression tag	UNP O07549
D	680	HIS	-	expression tag	UNP O07549
D	681	HIS	-	expression tag	UNP O07549

- Molecule 2 is a protein called Probable multidrug resistance ABC transporter ATP-binding/permease protein YheI.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	C	558	8861	2844	4467	732	796	22	0	0

There are 24 discrepancies between the modelled and reference sequences:

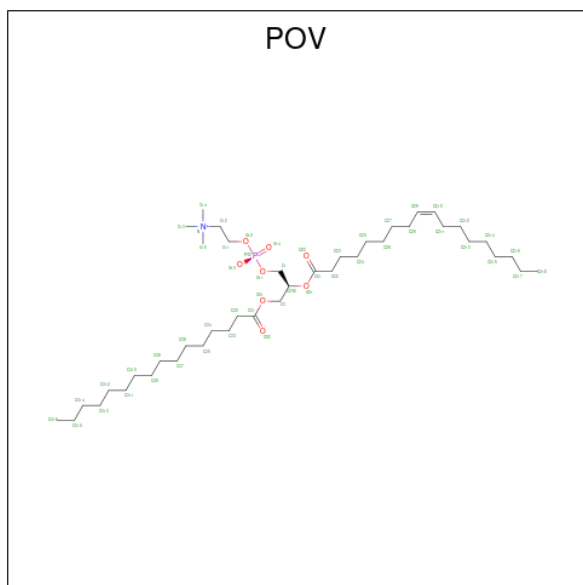
Chain	Residue	Modelled	Actual	Comment	Reference
C	-21	MET	-	initiating methionine	UNP O07550
C	-20	GLY	-	expression tag	UNP O07550
C	-19	SER	-	expression tag	UNP O07550

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-18	SER	-	expression tag	UNP O07550
C	-17	HIS	-	expression tag	UNP O07550
C	-16	HIS	-	expression tag	UNP O07550
C	-15	HIS	-	expression tag	UNP O07550
C	-14	HIS	-	expression tag	UNP O07550
C	-13	HIS	-	expression tag	UNP O07550
C	-12	HIS	-	expression tag	UNP O07550
C	-11	SER	-	expression tag	UNP O07550
C	-10	SER	-	expression tag	UNP O07550
C	-9	GLY	-	expression tag	UNP O07550
C	-8	LEU	-	expression tag	UNP O07550
C	-7	VAL	-	expression tag	UNP O07550
C	-6	PRO	-	expression tag	UNP O07550
C	-5	ARG	-	expression tag	UNP O07550
C	-4	GLY	-	expression tag	UNP O07550
C	-3	SER	-	expression tag	UNP O07550
C	-2	HIS	-	expression tag	UNP O07550
C	-1	MET	-	expression tag	UNP O07550
C	0	LEU	-	expression tag	UNP O07550
C	1	GLU	-	expression tag	UNP O07550
C	500	GLN	ASP	engineered mutation	UNP O07550

- Molecule 3 is (2S)-3-(hexadecanoyloxy)-2-[(9Z)-octadec-9-enoyloxy]propyl 2-(trimethylammonio)ethyl phosphate (CCD ID: POV) (formula: $C_{42}H_{82}NO_8P$) (labeled as "Ligand of Interest" by depositor).



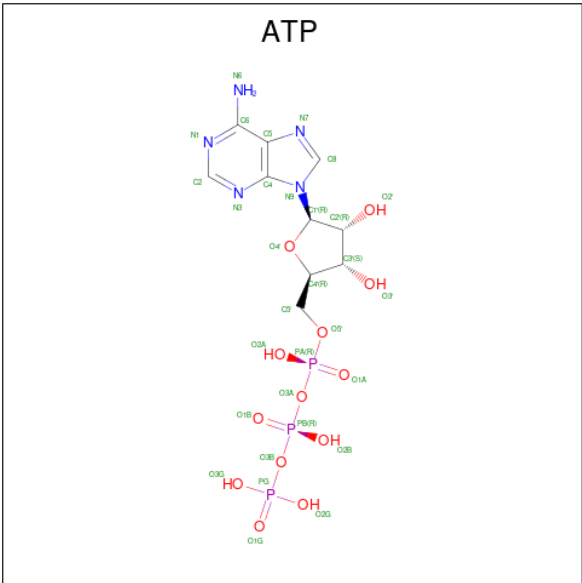
Mol	Chain	Residues	Atoms					AltConf
3	D	1	Total	C	H	O	P	0
			63	21	33	8	1	
3	D	1	Total	C	H			0
			39	15	24			
3	D	1	Total	C	H	O	P	0
			100	33	58	8	1	
3	D	1	Total	C	H			0
			71	25	46			
3	D	1	Total	C	H	O		0
			53	18	33	2		
3	D	1	Total	C	H	N	O	P
			119	38	71	1	8	1
3	D	1	Total	C	H	N	O	P
			101	33	58	1	8	1
3	D	1	Total	C	H	N	O	P
			73	23	40	1	8	1
3	D	1	Total	C	H	N	O	P
			103	32	61	1	8	1
3	D	1	Total	C	H			0
			51	18	33			
3	D	1	Total	C	H	N	O	P
			113	36	67	1	8	1
3	D	1	Total	C	H			0
			71	25	46			
3	D	1	Total	C	H			0
			71	25	46			
3	D	1	Total	C	H	N	O	P
			125	40	75	1	8	1
3	C	1	Total	C	H	O	P	0
			76	26	41	8	1	
3	C	1	Total	C	H	N	O	P
			90	29	51	1	8	1
3	C	1	Total	C	H	O		0
			39	14	24	1		
3	C	1	Total	C	H	N	O	P
			101	33	58	1	8	1
3	C	1	Total	C	H	O	P	0
			115	37	69	8	1	
3	C	1	Total	C	H			0
			98	34	64			
3	C	1	Total	C	H	N	O	P
			95	31	54	1	8	1
3	C	1	Total	C	H	O		0
			81	28	49	4		

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Mol	Chain	Residues	Atoms					AltConf
3	C	1	Total	C	H			0
			98	34	64			
3	C	1	Total	C	H	O	P	0
			84	27	48	8	1	
3	C	1	Total	C	H	N	O	P
			64	20	34	1	8	1

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃) (labeled as "Ligand of Interest" by depositor).

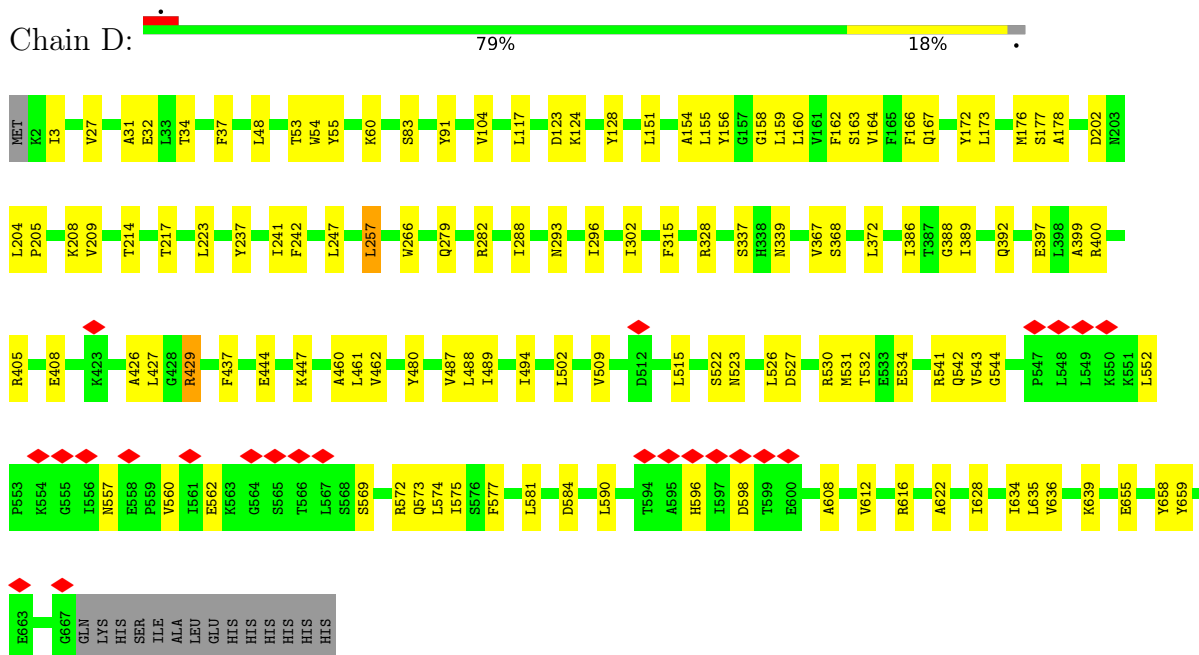


Mol	Chain	Residues	Atoms					AltConf
4	D	1	Total	C	H	N	O	P
			43	10	12	5	13	3
4	C	1	Total	C	H	N	O	P
			43	10	12	5	13	3

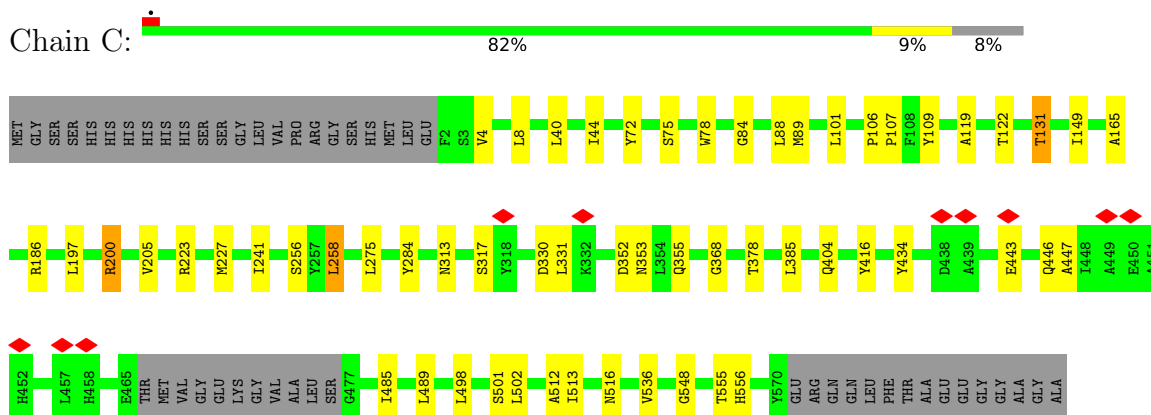
3 Residue-property plots

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 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: Probable multidrug resistance ABC transporter ATP-binding/permease protein YheH



- Molecule 2: Probable multidrug resistance ABC transporter ATP-binding/permease protein YheH



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	255381	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	54	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1600	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	5.439	Depositor
Minimum map value	-0.368	Depositor
Average map value	0.027	Depositor
Map value standard deviation	0.067	Depositor
Recommended contour level	0.43	Depositor
Map size ($\text{\AA}$)	330.0, 330.0, 330.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	Depositor

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: POV, ATP

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	D	0.32	0/5422	0.50	3/7319 (0.0%)
2	C	0.33	1/4486 (0.0%)	0.51	2/6072 (0.0%)
All	All	0.32	1/9908 (0.0%)	0.50	5/13391 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	200	ARG	C-N	6.62	1.42	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	257	LEU	N-CA-C	-6.79	104.47	112.89
2	C	131	THR	CB-CA-C	-5.78	101.02	110.85
2	C	258	LEU	CA-C-O	-5.41	114.81	120.55
1	D	337	SER	N-CA-C	5.31	117.50	111.02
1	D	429	ARG	CB-CA-C	5.00	118.00	109.84

There are no chirality outliers.

There are no planarity outliers.

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	D	5321	5387	5387	91	0
2	C	4394	4467	4465	39	0
3	C	385	556	556	7	0
3	D	462	691	691	9	0
4	C	31	12	12	0	0
4	D	31	12	12	0	0
All	All	10624	11125	11123	131	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 6.

All (131) 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:D:429:ARG:O	1:D:489:ILE:HD12	2.03	0.59
2:C:197:LEU:HA	2:C:227:MET:HE2	1.88	0.55
1:D:515:LEU:HD13	1:D:523:ASN:HD21	1.73	0.55
1:D:628:ILE:HG22	1:D:628:ILE:O	2.07	0.54
1:D:577:PHE:CE2	1:D:590:LEU:HD11	2.43	0.54
1:D:405:ARG:O	1:D:408:GLU:HG3	2.08	0.54
2:C:258:LEU:HD23	3:C:604:POV:H212	1.88	0.54
1:D:48:LEU:HD11	2:C:275:LEU:HD11	1.90	0.53
1:D:522:SER:O	1:D:526:LEU:N	2.42	0.53
1:D:209:VAL:HG21	2:C:205:VAL:HG21	1.90	0.53
1:D:569:SER:O	1:D:573:GLN:OE1	2.27	0.53
1:D:31:ALA:O	1:D:34:THR:HG22	2.08	0.53
2:C:313:ASN:O	2:C:317:SER:OG	2.23	0.53
1:D:447:LYS:NZ	1:D:639:LYS:O	2.41	0.52
1:D:489:ILE:HG23	1:D:489:ILE:O	2.09	0.52
2:C:443:GLU:HA	2:C:446:GLN:OE1	2.09	0.52
1:D:562:GLU:N	1:D:562:GLU:OE1	2.43	0.52
2:C:165:ALA:O	2:C:256:SER:OG	2.27	0.52
1:D:397:GLU:OE2	1:D:400:ARG:NH2	2.41	0.51
1:D:462:VAL:HG12	1:D:622:ALA:O	2.10	0.51
2:C:330:ASP:C	2:C:331:LEU:HD12	2.36	0.51
1:D:460:ALA:CB	1:D:628:ILE:HG21	2.41	0.51
1:D:437:PHE:N	1:D:444:GLU:OE2	2.44	0.51
1:D:214:THR:O	1:D:217:THR:OG1	2.28	0.51
1:D:531:MET:HE2	1:D:534:GLU:HB2	1.93	0.51
1:D:123:ASP:OD1	1:D:124:LYS:N	2.43	0.50
1:D:163:SER:O	1:D:167:GLN:OE1	2.30	0.50
1:D:328:ARG:HH21	3:D:708:POV:H14B	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:242:PHE:CD2	1:D:257:LEU:HD12	2.47	0.50
2:C:8:LEU:HD22	2:C:89:MET:HE1	1.93	0.50
2:C:101:LEU:O	2:C:109:TYR:OH	2.29	0.49
1:D:560:VAL:O	1:D:560:VAL:HG12	2.12	0.49
2:C:485:ILE:O	2:C:489:LEU:HD23	2.12	0.49
1:D:542:GLN:NE2	1:D:574:LEU:HD21	2.27	0.49
3:D:711:POV:H34	3:D:711:POV:H3A	1.95	0.49
1:D:509:VAL:HB	1:D:590:LEU:HD13	1.94	0.49
1:D:242:PHE:CE2	1:D:257:LEU:HD12	2.47	0.49
2:C:149:ILE:HD11	2:C:284:TYR:CD1	2.48	0.49
2:C:446:GLN:O	2:C:447:ALA:C	2.55	0.49
1:D:572:ARG:O	1:D:575:ILE:HG22	2.14	0.48
2:C:385:LEU:HD11	2:C:416:TYR:CE1	2.48	0.48
1:D:204:LEU:HD21	1:D:208:LYS:HB2	1.95	0.48
3:C:607:POV:H2	3:C:607:POV:H35	1.95	0.48
1:D:91:TYR:OH	1:D:104:VAL:HG21	2.14	0.47
1:D:155:LEU:O	1:D:156:TYR:C	2.58	0.47
2:C:200:ARG:NH1	2:C:223:ARG:NH1	2.63	0.46
1:D:573:GLN:O	1:D:577:PHE:HD1	1.97	0.46
2:C:75:SER:HA	2:C:78:TRP:CD1	2.51	0.46
1:D:489:ILE:HB	1:D:494:ILE:HD11	1.98	0.46
1:D:634:ILE:HG21	1:D:658:TYR:OH	2.15	0.46
1:D:3:ILE:HD12	1:D:3:ILE:H	1.80	0.46
1:D:151:LEU:CD1	3:D:710:POV:H210	2.46	0.46
1:D:27:VAL:HG11	1:D:166:PHE:HE2	1.80	0.45
1:D:462:VAL:HG23	1:D:636:VAL:HG12	1.97	0.45
2:C:502:LEU:HD11	2:C:513:ILE:HG21	1.98	0.45
1:D:37:PHE:CD2	3:D:702:POV:H25A	2.52	0.45
1:D:172:TYR:O	1:D:173:LEU:C	2.58	0.45
1:D:552:LEU:O	1:D:557:ASN:N	2.49	0.45
1:D:37:PHE:CD2	3:D:702:POV:C25	3.00	0.45
1:D:163:SER:O	1:D:164:VAL:C	2.60	0.45
1:D:388:GLY:O	1:D:392:GLN:OE1	2.34	0.45
1:D:655:GLU:HB3	1:D:659:TYR:HB2	1.98	0.44
2:C:352:ASP:O	2:C:353:ASN:HB2	2.17	0.44
1:D:388:GLY:O	1:D:389:ILE:C	2.60	0.44
2:C:555:THR:HG22	2:C:556:HIS:N	2.33	0.44
3:C:601:POV:H3	3:C:601:POV:H22	1.99	0.44
2:C:512:ALA:O	2:C:516:ASN:ND2	2.47	0.44
1:D:339:ASN:HB3	2:C:72:TYR:CD1	2.53	0.43
1:D:426:ALA:O	1:D:427:LEU:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:526:LEU:O	1:D:527:ASP:OD1	2.35	0.43
2:C:404:GLN:N	2:C:404:GLN:OE1	2.51	0.43
1:D:628:ILE:HG22	1:D:634:ILE:HD11	2.00	0.43
1:D:608:ALA:O	1:D:612:VAL:HG12	2.18	0.43
1:D:34:THR:HB	3:D:702:POV:H27A	2.00	0.43
1:D:60:LYS:O	1:D:60:LYS:HG2	2.19	0.43
1:D:596:HIS:CE1	1:D:598:ASP:HA	2.54	0.43
2:C:355:GLN:N	2:C:548:GLY:O	2.51	0.43
1:D:527:ASP:OD1	1:D:527:ASP:C	2.61	0.43
2:C:106:PRO:N	2:C:107:PRO:CD	2.82	0.43
2:C:258:LEU:HD23	3:C:604:POV:C212	2.49	0.43
1:D:32:GLU:OE2	1:D:167:GLN:NE2	2.52	0.43
1:D:279:GLN:O	1:D:282:ARG:HG2	2.19	0.43
1:D:530:ARG:O	1:D:530:ARG:HD2	2.19	0.43
1:D:584:ASP:OD1	1:D:616:ARG:NH1	2.52	0.42
1:D:237:TYR:CE1	1:D:241:ILE:HD11	2.54	0.42
1:D:302:ILE:HA	2:C:434:TYR:OH	2.19	0.42
2:C:378:THR:HG22	2:C:378:THR:O	2.19	0.42
1:D:162:PHE:O	1:D:163:SER:C	2.62	0.42
1:D:543:VAL:HG13	1:D:544:GLY:N	2.34	0.42
1:D:159:LEU:O	1:D:160:LEU:C	2.63	0.42
1:D:55:TYR:CE1	1:D:83:SER:O	2.72	0.42
1:D:288:ILE:HG23	1:D:315:PHE:HE1	1.84	0.42
1:D:223:LEU:HB2	1:D:399:ALA:HB1	2.02	0.42
1:D:202:ASP:OD2	1:D:480:TYR:OH	2.31	0.42
1:D:204:LEU:HD23	1:D:205:PRO:O	2.19	0.42
1:D:177:SER:O	1:D:178:ALA:C	2.63	0.42
1:D:266:TRP:CZ3	1:D:386:ILE:HG21	2.55	0.41
1:D:293:ASN:O	1:D:296:ILE:HG22	2.20	0.41
1:D:515:LEU:HD13	1:D:523:ASN:ND2	2.35	0.41
2:C:4:VAL:HG11	2:C:131:THR:HG21	2.02	0.41
1:D:176:MET:O	1:D:177:SER:C	2.61	0.41
1:D:581:LEU:C	1:D:581:LEU:HD23	2.45	0.41
1:D:247:LEU:HD22	3:D:711:POV:H3	2.02	0.41
3:C:603:POV:H35	3:C:603:POV:H32A	1.89	0.41
1:D:223:LEU:C	1:D:223:LEU:HD23	2.45	0.41
2:C:200:ARG:HD3	2:C:223:ARG:NH2	2.35	0.41
2:C:498:LEU:HD22	2:C:501:SER:CB	2.50	0.41
3:D:704:POV:H31G	3:D:704:POV:H218	2.03	0.41
2:C:75:SER:HA	2:C:78:TRP:NE1	2.35	0.41
1:D:27:VAL:HG11	1:D:166:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:607:POV:H28A	3:C:607:POV:H33	2.03	0.41
1:D:53:THR:HG22	1:D:54:TRP:N	2.36	0.41
1:D:372:LEU:CD2	2:C:44:ILE:HD11	2.50	0.41
1:D:487:VAL:C	1:D:488:LEU:HD22	2.46	0.41
1:D:541:ARG:HH12	1:D:581:LEU:HD13	1.85	0.41
1:D:209:VAL:CG2	2:C:205:VAL:HG21	2.51	0.41
1:D:461:LEU:HD13	1:D:635:LEU:HB2	2.02	0.41
2:C:119:ALA:HA	2:C:122:THR:HG22	2.03	0.41
1:D:154:ALA:O	1:D:155:LEU:C	2.64	0.40
1:D:158:GLY:O	1:D:159:LEU:C	2.62	0.40
1:D:532:THR:O	1:D:534:GLU:N	2.50	0.40
1:D:367:VAL:O	1:D:368:SER:OG	2.30	0.40
1:D:494:ILE:HG23	1:D:502:LEU:HD22	2.04	0.40
2:C:84:GLY:O	2:C:88:LEU:HD13	2.21	0.40
2:C:186:ARG:HD3	3:C:608:POV:O31	2.21	0.40
1:D:117:LEU:N	1:D:128:TYR:O	2.53	0.40
2:C:368:GLY:HA3	2:C:536:VAL:HG21	2.02	0.40
3:D:701:POV:H33A	3:D:701:POV:H2	2.02	0.40
1:D:634:ILE:HG22	1:D:635:LEU:N	2.36	0.40
2:C:40:LEU:HD23	2:C:44:ILE:HD12	2.04	0.40
2:C:186:ARG:HG3	2:C:241:ILE:HD12	2.04	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	D	664/681 (98%)	605 (91%)	59 (9%)	0	100	100
2	C	554/607 (91%)	529 (96%)	25 (4%)	0	100	100
All	All	1218/1288 (95%)	1134 (93%)	84 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	D	567/581 (98%)	567 (100%)	0	100	100
2	C	471/508 (93%)	471 (100%)	0	100	100
All	All	1038/1089 (95%)	1038 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	63	ASN
1	D	143	GLN
1	D	194	GLN
1	D	303	GLN
1	D	454	GLN
1	D	464	HIS
1	D	596	HIS
1	D	648	HIS
2	C	80	HIS
2	C	100	HIS
2	C	192	ASN
2	C	445	GLN

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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

27 ligands 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	$\# Z > 2$
3	POV	C	609	-	32,32,51	0.26	0	30,30,59	0.14	0
3	POV	D	702	-	14,14,51	0.33	0	13,13,59	0.15	0
3	POV	C	605	-	45,45,51	0.41	0	48,50,59	0.53	1 (2%)
3	POV	D	708	-	32,32,51	0.41	0	38,40,59	0.41	0
3	POV	C	610	-	35,35,51	0.44	0	38,40,59	0.53	1 (2%)
4	ATP	D	715	-	32,33,33	0.51	0	48,52,52	0.67	0
3	POV	C	606	-	32,32,51	0.26	0	30,30,59	0.15	0
3	POV	D	711	-	45,45,51	0.33	0	51,53,59	0.33	0
3	POV	D	704	-	23,23,51	0.23	0	21,21,59	0.10	0
3	POV	D	713	-	23,23,51	0.25	0	21,21,59	0.14	0
3	POV	C	611	-	29,29,51	0.38	0	35,37,59	0.59	0
3	POV	C	602	-	38,38,51	0.37	0	44,46,59	0.55	1 (2%)
3	POV	C	603	-	14,14,51	0.50	0	13,13,59	0.15	0
3	POV	D	701	-	29,29,51	0.46	0	32,34,59	0.68	1 (3%)
4	ATP	C	612	-	32,33,33	0.51	0	48,52,52	0.60	0
3	POV	D	710	-	17,17,51	0.19	0	16,16,59	0.19	0
3	POV	C	608	-	30,30,51	0.34	0	30,30,59	0.27	0
3	POV	D	703	-	41,41,51	0.43	0	44,46,59	0.47	1 (2%)
3	POV	D	705	-	19,19,51	0.40	0	19,19,59	0.28	0
3	POV	C	607	-	40,40,51	0.34	0	46,48,59	0.39	0
3	POV	D	707	-	42,42,51	0.36	0	48,50,59	0.47	0
3	POV	C	601	-	34,34,51	0.44	0	37,39,59	0.70	1 (2%)
3	POV	C	604	-	42,42,51	0.44	0	48,50,59	0.48	0
3	POV	D	709	-	41,41,51	0.38	0	47,49,59	0.43	0
3	POV	D	714	-	49,49,51	0.34	0	55,57,59	0.37	0
3	POV	D	712	-	23,23,51	0.28	0	21,21,59	0.13	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	POV	D	706	-	47,47,51	0.33	0	53,55,59	0.41	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
3	POV	C	609	-	-	17/28/28/55	-
3	POV	D	702	-	-	4/12/12/55	-
3	POV	C	605	-	-	22/47/47/55	-
3	POV	D	708	-	-	11/36/36/55	-
3	POV	C	610	-	-	11/37/37/55	-
4	ATP	D	715	-	-	2/22/38/38	0/3/3/3
3	POV	C	606	-	-	7/28/28/55	-
3	POV	D	711	-	-	18/49/49/55	-
3	POV	D	704	-	-	6/19/19/55	-
3	POV	D	713	-	-	2/19/19/55	-
3	POV	C	611	-	-	11/33/33/55	-
3	POV	C	602	-	-	15/42/42/55	-
3	POV	C	603	-	-	4/12/12/55	-
3	POV	D	701	-	-	8/31/31/55	-
4	ATP	C	612	-	-	2/22/38/38	0/3/3/3
3	POV	D	710	-	-	4/15/15/55	-
3	POV	C	608	-	-	5/26/26/55	-
3	POV	D	703	-	-	7/43/43/55	-
3	POV	D	705	-	-	4/17/17/55	-
3	POV	C	607	-	-	15/44/44/55	-
3	POV	D	707	-	-	13/46/46/55	-
3	POV	C	601	-	-	11/36/36/55	-
3	POV	C	604	-	-	10/46/46/55	-
3	POV	D	709	-	-	18/45/45/55	-
3	POV	D	714	-	-	13/53/53/55	-
3	POV	D	712	-	-	5/19/19/55	-
3	POV	D	706	-	-	15/51/51/55	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	610	POV	O13-P-O14	2.50	120.56	110.83
3	C	605	POV	O13-P-O14	2.37	120.07	110.83
3	D	701	POV	O13-P-O14	2.35	119.97	110.83
3	C	602	POV	O21-C21-C22	2.33	116.52	111.48
3	C	601	POV	O13-P-O14	2.20	119.39	110.83
3	D	703	POV	O13-P-O14	2.15	119.23	110.83

There are no chirality outliers.

All (260) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	701	POV	O22-C21-O21-C2
3	D	703	POV	C22-C21-O21-C2
3	D	707	POV	C1-O11-P-O12
3	D	707	POV	C2-C1-O11-P
3	D	707	POV	C3-C2-O21-C21
3	D	707	POV	C22-C21-O21-C2
3	D	708	POV	C1-O11-P-O14
3	D	708	POV	C22-C21-O21-C2
3	D	709	POV	C11-O12-P-O11
3	D	711	POV	C1-O11-P-O12
3	D	711	POV	C1-O11-P-O13
3	D	711	POV	O12-C11-C12-N
3	D	711	POV	C22-C21-O21-C2
3	D	712	POV	C26-C27-C28-C29
3	D	714	POV	C11-O12-P-O14
3	C	601	POV	C22-C21-O21-C2
3	C	601	POV	O22-C21-O21-C2
3	C	602	POV	C1-O11-P-O12
3	C	602	POV	C11-O12-P-O11
3	C	602	POV	C11-O12-P-O14
3	C	602	POV	C22-C21-O21-C2
3	C	604	POV	C11-O12-P-O14
3	C	604	POV	C22-C21-O21-C2
3	C	605	POV	C1-O11-P-O12
3	C	605	POV	C1-O11-P-O13
3	C	605	POV	C1-O11-P-O14
3	C	607	POV	C1-O11-P-O14
3	C	611	POV	C1-O11-P-O12
3	C	611	POV	C1-O11-P-O13

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Mol	Chain	Res	Type	Atoms
3	C	611	POV	C11-O12-P-O13
3	C	611	POV	O12-C11-C12-N
3	C	611	POV	C12-C11-O12-P
3	C	611	POV	C22-C21-O21-C2
3	C	611	POV	O22-C21-O21-C2
3	D	707	POV	O32-C31-O31-C3
3	D	707	POV	C32-C31-O31-C3
3	C	610	POV	C32-C31-O31-C3
3	D	709	POV	O32-C31-O31-C3
3	C	601	POV	O32-C31-O31-C3
3	C	604	POV	O32-C31-O31-C3
3	C	610	POV	O32-C31-O31-C3
3	C	611	POV	O32-C31-O31-C3
3	D	703	POV	O22-C21-O21-C2
3	D	707	POV	O22-C21-O21-C2
3	D	711	POV	O22-C21-O21-C2
3	C	602	POV	O22-C21-O21-C2
3	D	709	POV	C32-C31-O31-C3
3	C	601	POV	C32-C31-O31-C3
3	C	604	POV	C32-C31-O31-C3
3	D	701	POV	C22-C21-O21-C2
3	C	611	POV	C32-C31-O31-C3
3	D	704	POV	C211-C210-C29-C28
3	D	711	POV	C211-C210-C29-C28
3	C	605	POV	C211-C210-C29-C28
3	D	708	POV	O22-C21-O21-C2
3	C	604	POV	O22-C21-O21-C2
4	D	715	ATP	O4'-C4'-C5'-O5'
4	D	715	ATP	C3'-C4'-C5'-O5'
3	C	602	POV	C2-C1-O11-P
3	D	708	POV	C32-C31-O31-C3
3	D	701	POV	C32-C31-O31-C3
3	D	711	POV	C32-C31-O31-C3
3	D	701	POV	O32-C31-O31-C3
3	C	604	POV	O11-C1-C2-O21
3	C	605	POV	O11-C1-C2-O21
3	D	706	POV	C32-C31-O31-C3
3	C	601	POV	C21-C22-C23-C24
3	D	708	POV	O32-C31-O31-C3
3	D	711	POV	O32-C31-O31-C3
3	D	712	POV	C211-C210-C29-C28
3	C	602	POV	C32-C31-O31-C3

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Mol	Chain	Res	Type	Atoms
3	D	710	POV	C211-C210-C29-C28
3	D	706	POV	O32-C31-O31-C3
3	C	605	POV	C310-C311-C312-C313
3	C	607	POV	C32-C33-C34-C35
3	C	605	POV	C212-C213-C214-C215
3	C	602	POV	O32-C31-O31-C3
3	C	605	POV	C311-C312-C313-C314
3	D	710	POV	C25-C26-C27-C28
3	C	609	POV	C33-C34-C35-C36
3	C	605	POV	C36-C37-C38-C39
3	C	610	POV	C22-C23-C24-C25
3	C	611	POV	C33-C34-C35-C36
3	D	704	POV	C39-C310-C311-C312
3	D	711	POV	C34-C35-C36-C37
3	D	712	POV	C212-C213-C214-C215
3	D	714	POV	C22-C23-C24-C25
3	D	711	POV	C21-C22-C23-C24
3	D	704	POV	C212-C213-C214-C215
3	D	709	POV	C34-C35-C36-C37
3	C	609	POV	C37-C38-C39-C310
3	D	711	POV	C311-C312-C313-C314
3	C	601	POV	C36-C37-C38-C39
3	C	605	POV	C32-C31-O31-C3
3	D	705	POV	C211-C212-C213-C214
3	D	705	POV	C212-C213-C214-C215
3	C	606	POV	C37-C38-C39-C310
3	C	610	POV	C32-C33-C34-C35
3	C	608	POV	C37-C38-C39-C310
3	D	709	POV	C37-C38-C39-C310
3	C	609	POV	C26-C27-C28-C29
3	D	714	POV	C37-C38-C39-C310
3	C	601	POV	C25-C26-C27-C28
3	C	605	POV	O32-C31-O31-C3
3	C	601	POV	C211-C210-C29-C28
3	C	609	POV	C214-C215-C216-C217
3	C	606	POV	C39-C310-C311-C312
3	C	604	POV	O11-C1-C2-C3
3	C	605	POV	O11-C1-C2-C3
3	C	609	POV	C34-C35-C36-C37
3	C	602	POV	C21-C22-C23-C24
3	C	603	POV	C36-C37-C38-C39
3	D	707	POV	C25-C26-C27-C28

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Mol	Chain	Res	Type	Atoms
3	C	602	POV	C23-C24-C25-C26
3	C	609	POV	C27-C28-C29-C210
3	C	611	POV	C32-C33-C34-C35
3	C	609	POV	C310-C311-C312-C313
3	C	604	POV	C22-C23-C24-C25
3	C	607	POV	C311-C310-C39-C38
3	C	607	POV	C31-C32-C33-C34
3	C	609	POV	C211-C210-C29-C28
3	C	609	POV	C24-C25-C26-C27
3	D	709	POV	C22-C23-C24-C25
3	D	708	POV	O21-C2-C3-O31
3	C	609	POV	C32-C33-C34-C35
3	D	706	POV	O31-C31-C32-C33
3	D	710	POV	C213-C214-C215-C216
3	C	610	POV	C36-C37-C38-C39
3	C	607	POV	C32-C31-O31-C3
3	D	714	POV	C25-C26-C27-C28
3	D	702	POV	C25-C26-C27-C28
3	D	704	POV	C310-C311-C312-C313
3	D	706	POV	C22-C23-C24-C25
3	C	604	POV	C33-C34-C35-C36
3	C	606	POV	C214-C215-C216-C217
3	C	607	POV	C211-C210-C29-C28
3	D	712	POV	C213-C214-C215-C216
3	D	708	POV	C1-C2-C3-O31
3	C	609	POV	C312-C313-C314-C315
3	D	713	POV	C215-C216-C217-C218
3	C	605	POV	C2-C1-O11-P
3	C	607	POV	C2-C1-O11-P
3	D	701	POV	C32-C33-C34-C35
3	C	605	POV	C23-C24-C25-C26
3	C	605	POV	C213-C214-C215-C216
3	D	714	POV	C210-C211-C212-C213
3	D	714	POV	O11-C1-C2-C3
3	D	704	POV	C29-C210-C211-C212
3	C	606	POV	C24-C25-C26-C27
3	C	607	POV	O32-C31-O31-C3
3	C	601	POV	C1-O11-P-O13
3	D	709	POV	C22-C21-O21-C2
3	C	607	POV	O31-C31-C32-C33
3	C	605	POV	C22-C23-C24-C25
3	C	601	POV	C1-C2-O21-C21

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Mol	Chain	Res	Type	Atoms
3	D	702	POV	C26-C27-C28-C29
3	D	714	POV	O11-C1-C2-O21
3	C	610	POV	O11-C1-C2-O21
3	C	605	POV	C27-C28-C29-C210
3	C	602	POV	C12-C11-O12-P
3	D	709	POV	O21-C2-C3-O31
3	D	708	POV	C23-C24-C25-C26
3	D	706	POV	C23-C24-C25-C26
3	D	706	POV	O12-C11-C12-N
3	D	707	POV	O12-C11-C12-N
3	D	708	POV	O12-C11-C12-N
3	D	714	POV	O12-C11-C12-N
3	C	607	POV	O12-C11-C12-N
3	C	610	POV	C39-C310-C311-C312
3	D	711	POV	C37-C38-C39-C310
3	D	709	POV	C310-C311-C312-C313
3	D	708	POV	C26-C27-C28-C29
3	C	609	POV	C311-C312-C313-C314
3	C	610	POV	O11-C1-C2-C3
3	D	709	POV	O22-C21-O21-C2
3	D	703	POV	C310-C311-C312-C313
3	C	607	POV	O11-C1-C2-O21
3	D	714	POV	C212-C213-C214-C215
3	D	709	POV	C1-C2-C3-O31
3	D	711	POV	C313-C314-C315-C316
3	D	702	POV	C27-C28-C29-C210
3	D	707	POV	C1-O11-P-O14
3	D	709	POV	C1-O11-P-O14
3	D	709	POV	C11-O12-P-O14
3	C	602	POV	C1-O11-P-O14
3	D	711	POV	C310-C311-C312-C313
3	D	701	POV	C2-C1-O11-P
3	D	711	POV	C2-C1-O11-P
3	C	610	POV	C21-C22-C23-C24
3	D	707	POV	C27-C28-C29-C210
3	D	706	POV	C1-C2-O21-C21
3	C	610	POV	C310-C311-C312-C313
3	D	709	POV	C311-C310-C39-C38
3	C	607	POV	O11-C1-C2-C3
3	C	605	POV	C312-C313-C314-C315
3	D	714	POV	O32-C31-O31-C3
3	D	709	POV	C25-C26-C27-C28

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Mol	Chain	Res	Type	Atoms
3	C	603	POV	C35-C36-C37-C38
3	C	609	POV	C25-C26-C27-C28
3	D	706	POV	O21-C2-C3-O31
3	D	714	POV	C32-C31-O31-C3
3	C	606	POV	C211-C210-C29-C28
3	D	706	POV	C22-C21-O21-C2
3	D	705	POV	C210-C211-C212-C213
3	C	609	POV	C36-C37-C38-C39
3	D	704	POV	C311-C310-C39-C38
3	D	702	POV	C23-C24-C25-C26
3	D	710	POV	C26-C27-C28-C29
3	C	603	POV	C32-C33-C34-C35
3	D	703	POV	C39-C310-C311-C312
3	D	706	POV	O32-C31-C32-C33
3	D	706	POV	C33-C34-C35-C36
3	D	706	POV	C35-C36-C37-C38
3	C	606	POV	C311-C312-C313-C314
3	C	605	POV	C214-C215-C216-C217
3	C	607	POV	C33-C34-C35-C36
3	D	714	POV	C27-C28-C29-C210
3	C	601	POV	C33-C34-C35-C36
3	D	711	POV	O31-C31-C32-C33
3	C	609	POV	C21-C22-C23-C24
3	D	703	POV	C25-C26-C27-C28
3	D	707	POV	C22-C23-C24-C25
3	D	706	POV	O22-C21-O21-C2
3	D	706	POV	C34-C35-C36-C37
3	C	608	POV	O21-C21-C22-C23
3	C	605	POV	C215-C216-C217-C218
3	D	701	POV	C34-C35-C36-C37
3	D	701	POV	C310-C311-C312-C313
3	D	703	POV	C311-C312-C313-C314
3	C	610	POV	C1-O11-P-O13
3	C	604	POV	C29-C210-C211-C212
3	D	705	POV	C213-C214-C215-C216
3	D	708	POV	C2-C1-O11-P
3	C	602	POV	O11-C1-C2-O21
3	C	609	POV	C215-C216-C217-C218
3	C	608	POV	C35-C36-C37-C38
3	C	605	POV	C35-C36-C37-C38
3	D	713	POV	C213-C214-C215-C216
4	C	612	ATP	PB-O3A-PA-O1A

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Mol	Chain	Res	Type	Atoms
4	C	612	ATP	PB-O3A-PA-O2A
3	C	602	POV	C31-C32-C33-C34
3	C	605	POV	C313-C314-C315-C316
3	D	709	POV	C21-C22-C23-C24
3	C	608	POV	O22-C21-C22-C23
3	D	706	POV	C1-C2-C3-O31
3	D	709	POV	O21-C21-C22-C23
3	C	609	POV	C22-C23-C24-C25
3	C	607	POV	O32-C31-C32-C33
3	D	707	POV	C24-C25-C26-C27
3	D	711	POV	C3-C2-O21-C21
3	D	714	POV	C39-C310-C311-C312
3	C	603	POV	C37-C38-C39-C310
3	C	606	POV	C36-C37-C38-C39
3	D	703	POV	C33-C34-C35-C36
3	D	709	POV	O22-C21-C22-C23
3	D	712	POV	C313-C314-C315-C316
3	C	602	POV	C11-C12-N-C15
3	D	711	POV	C23-C24-C25-C26
3	C	608	POV	C27-C28-C29-C210
3	C	607	POV	C24-C25-C26-C27

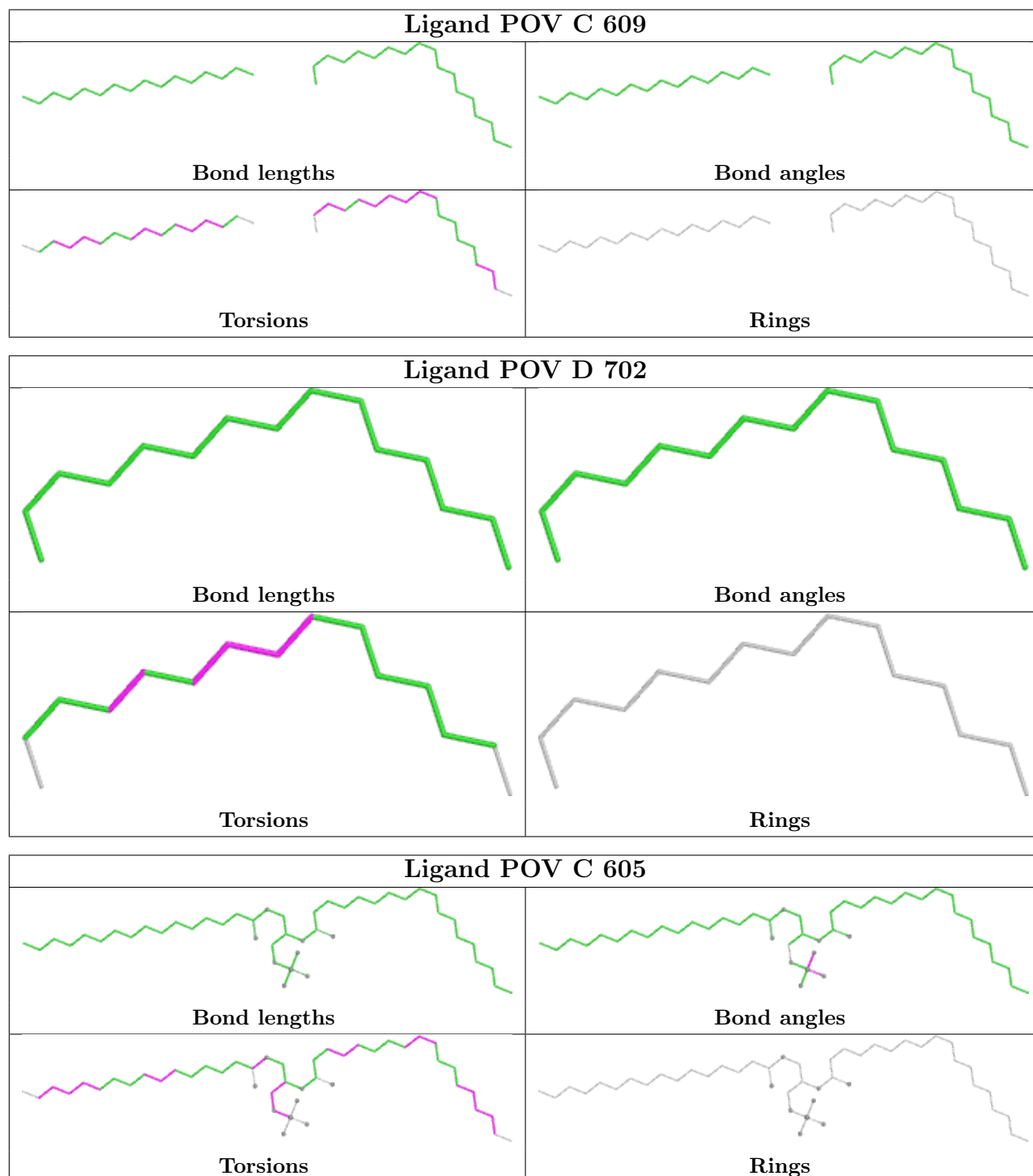
There are no ring outliers.

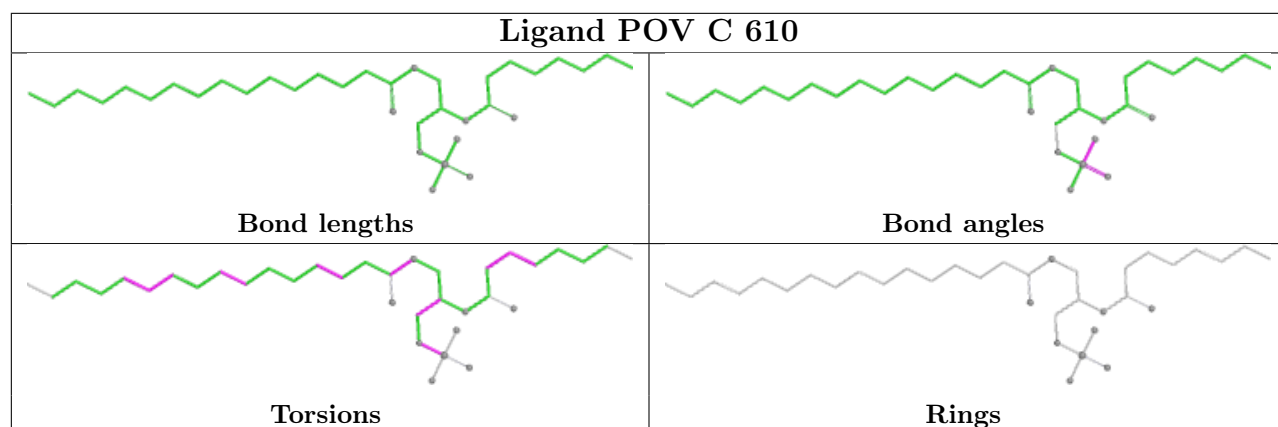
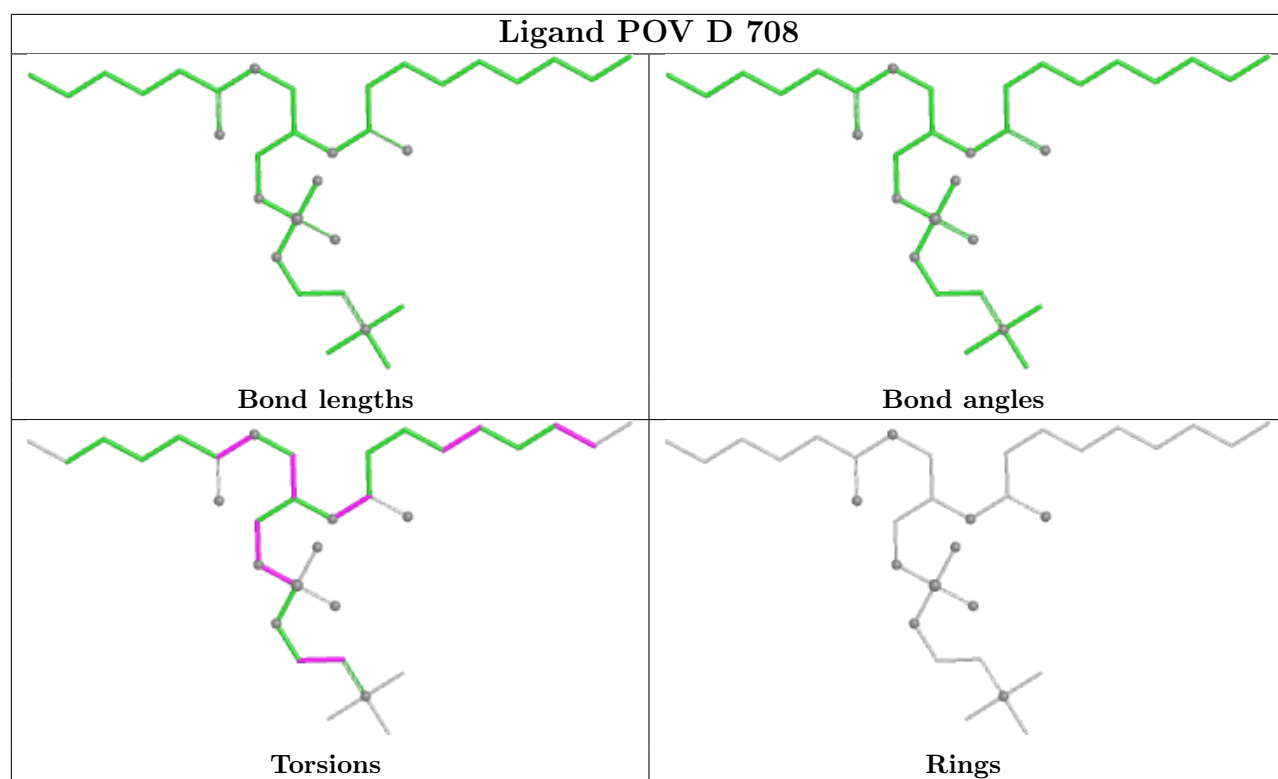
11 monomers are involved in 16 short contacts:

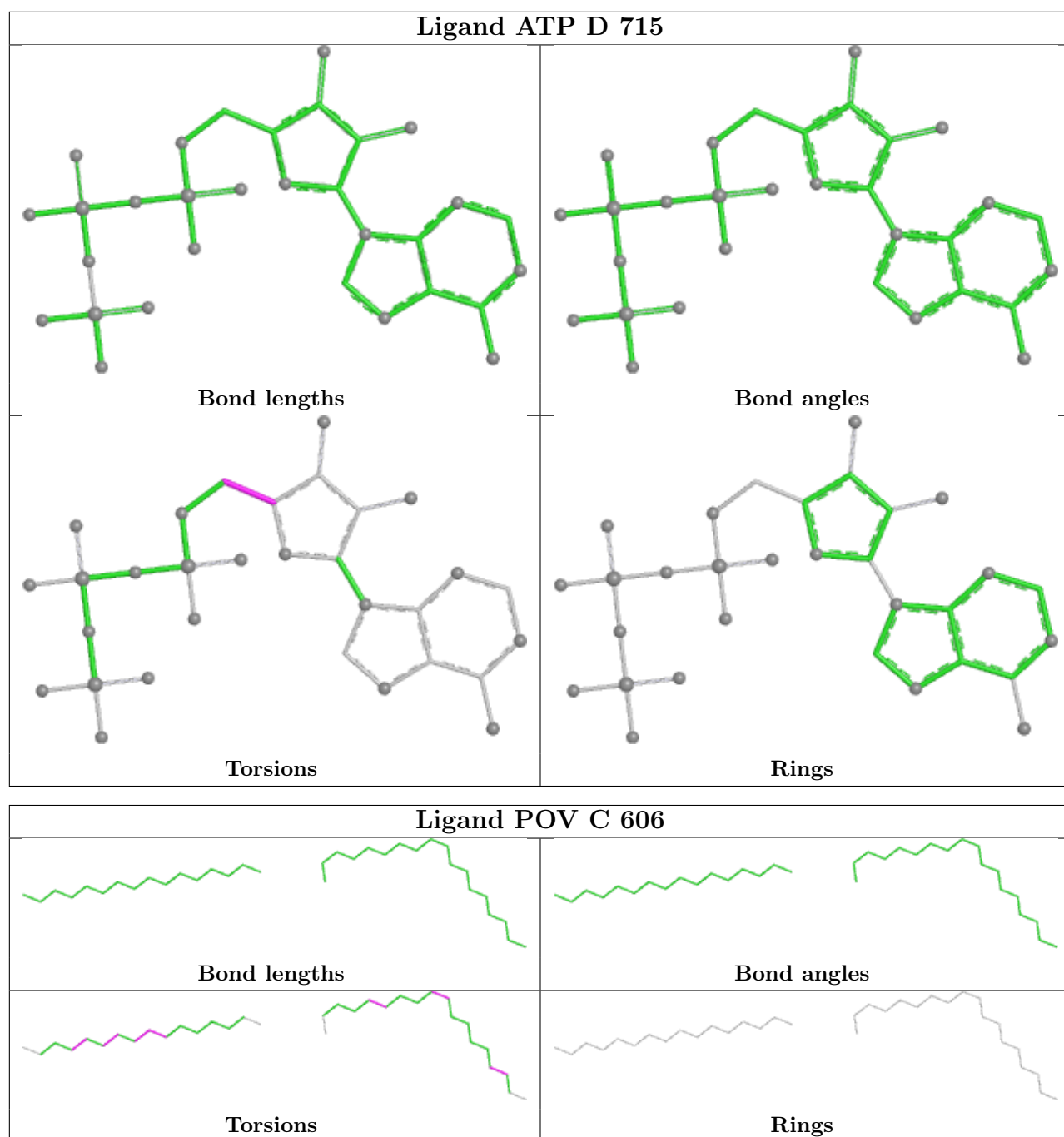
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	702	POV	3	0
3	D	708	POV	1	0
3	D	711	POV	2	0
3	D	704	POV	1	0
3	C	603	POV	1	0
3	D	701	POV	1	0
3	D	710	POV	1	0
3	C	608	POV	1	0
3	C	607	POV	2	0
3	C	601	POV	1	0
3	C	604	POV	2	0

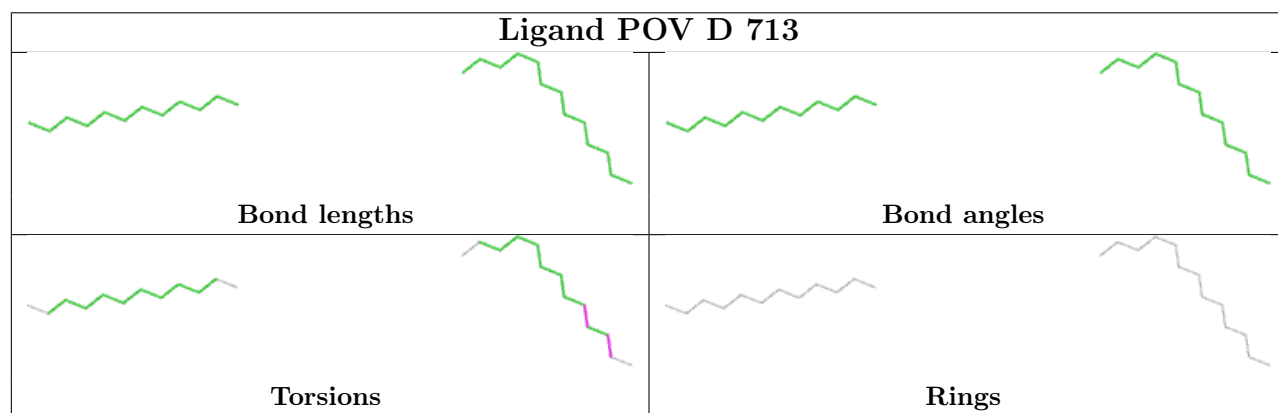
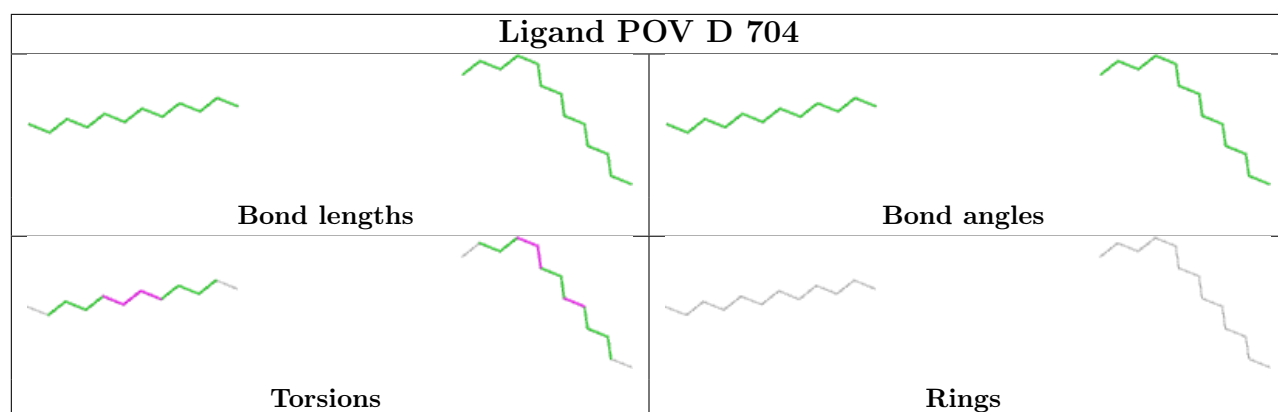
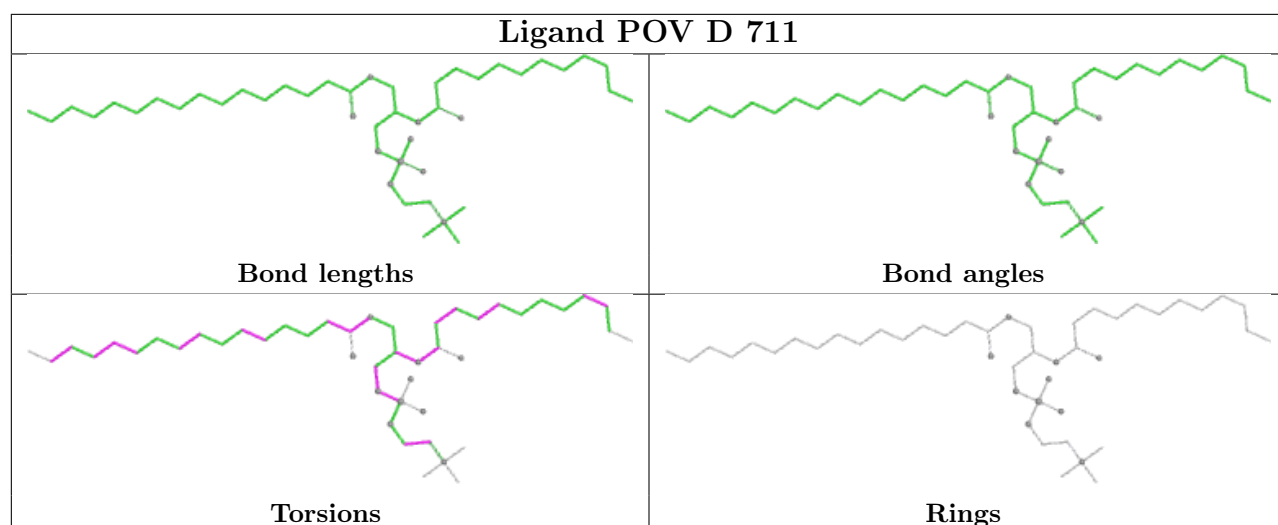
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

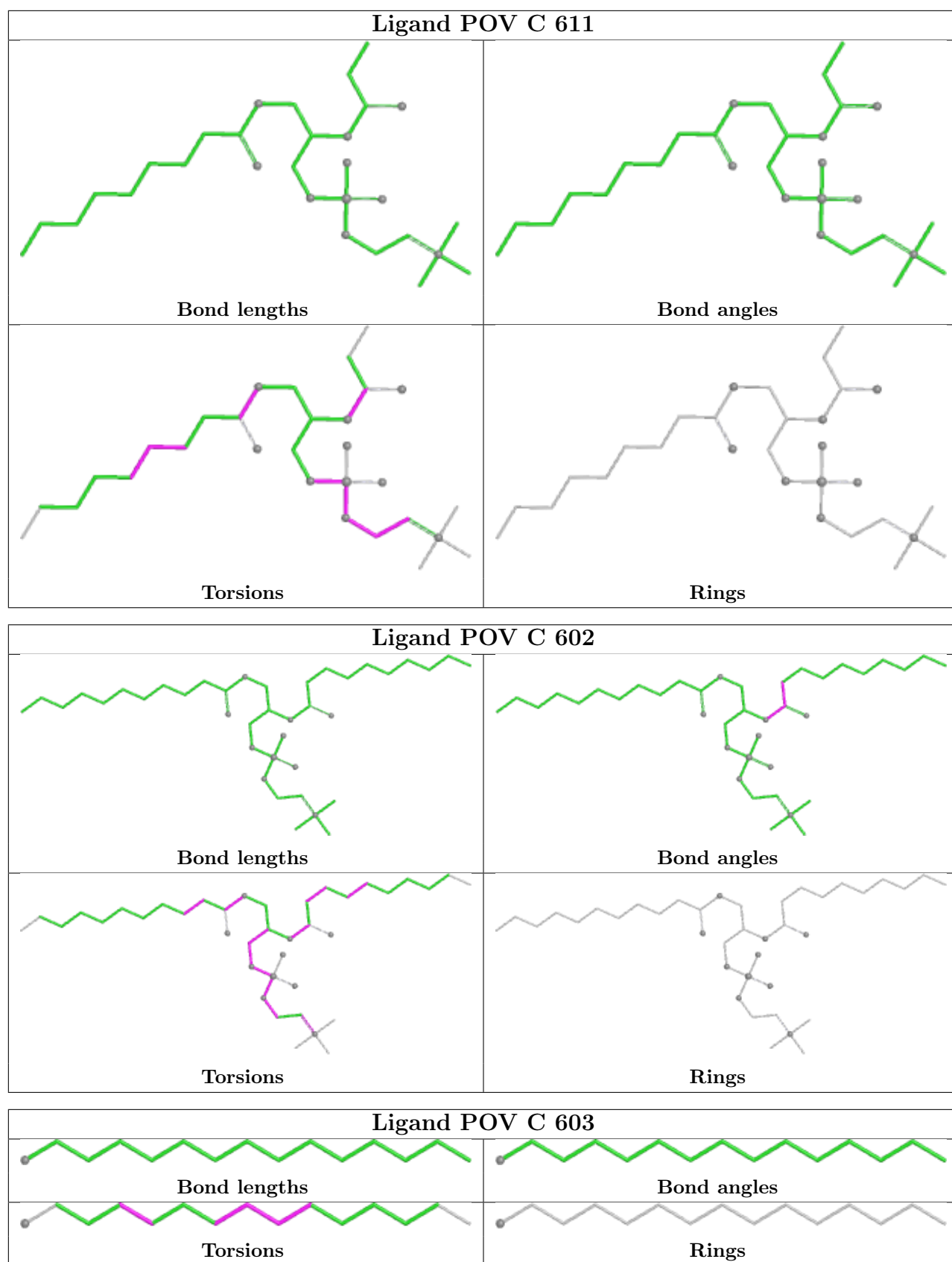
within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question 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.

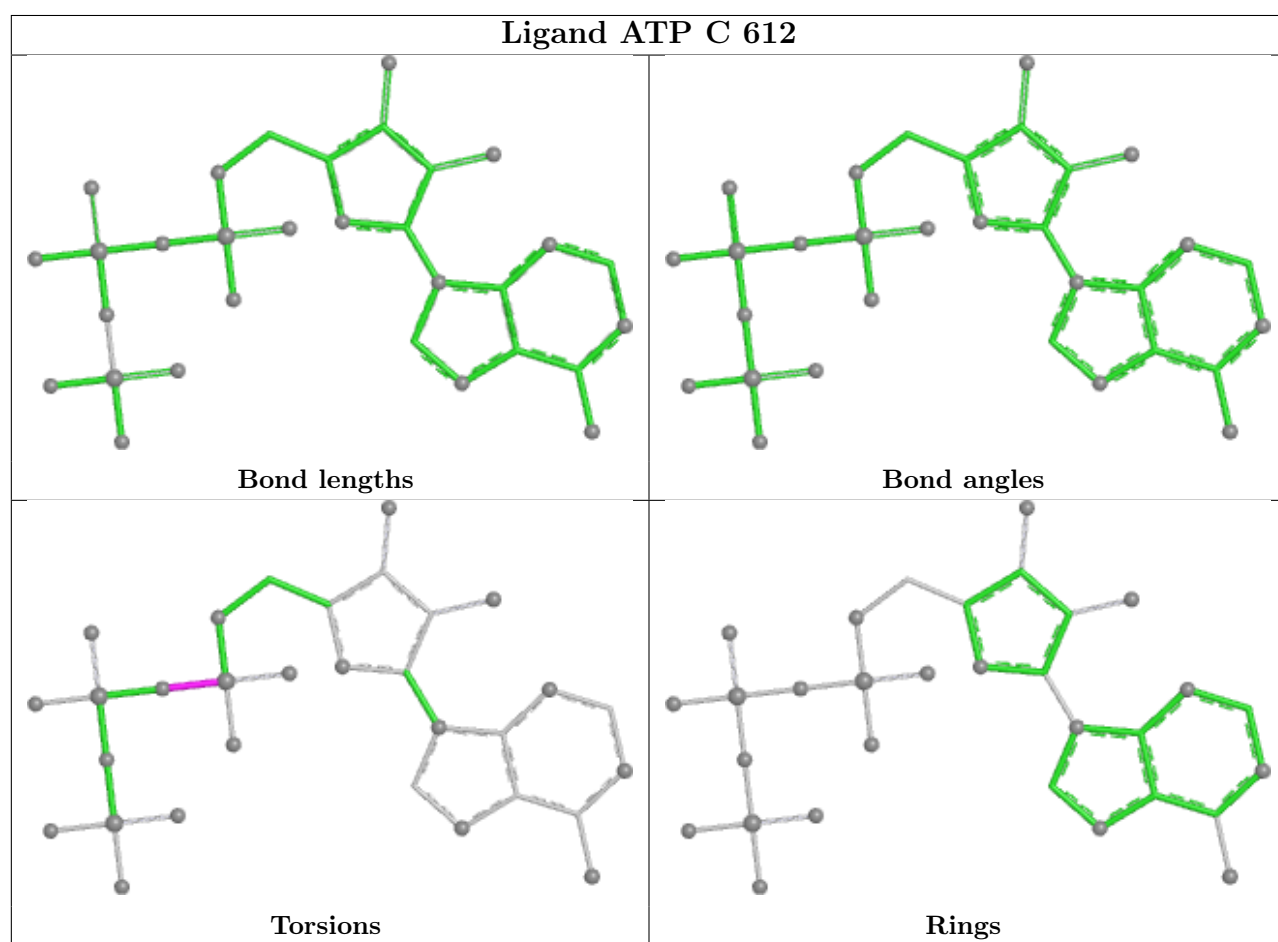
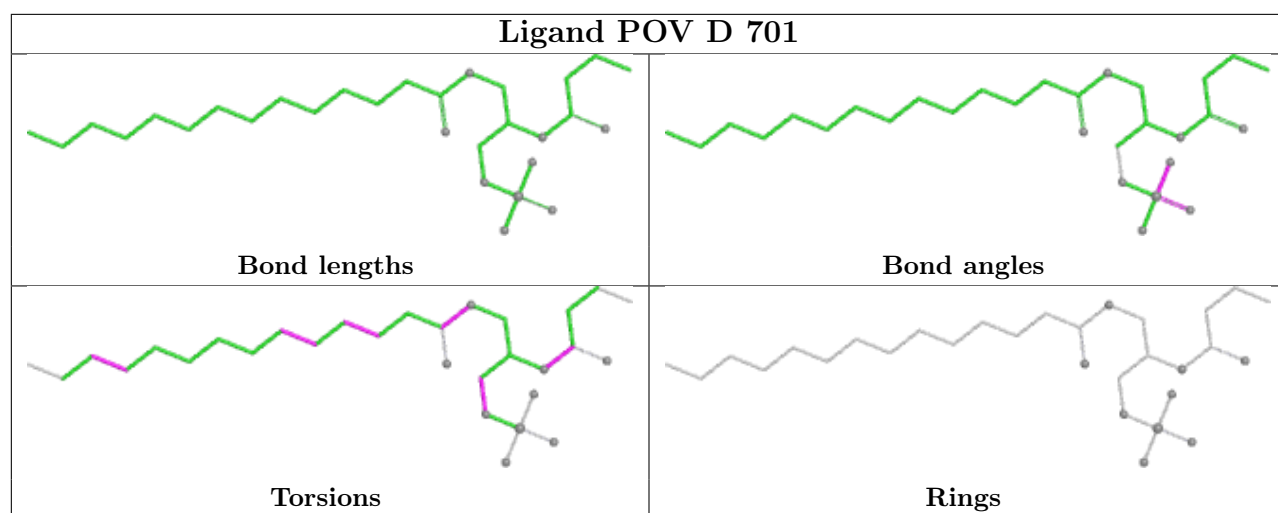


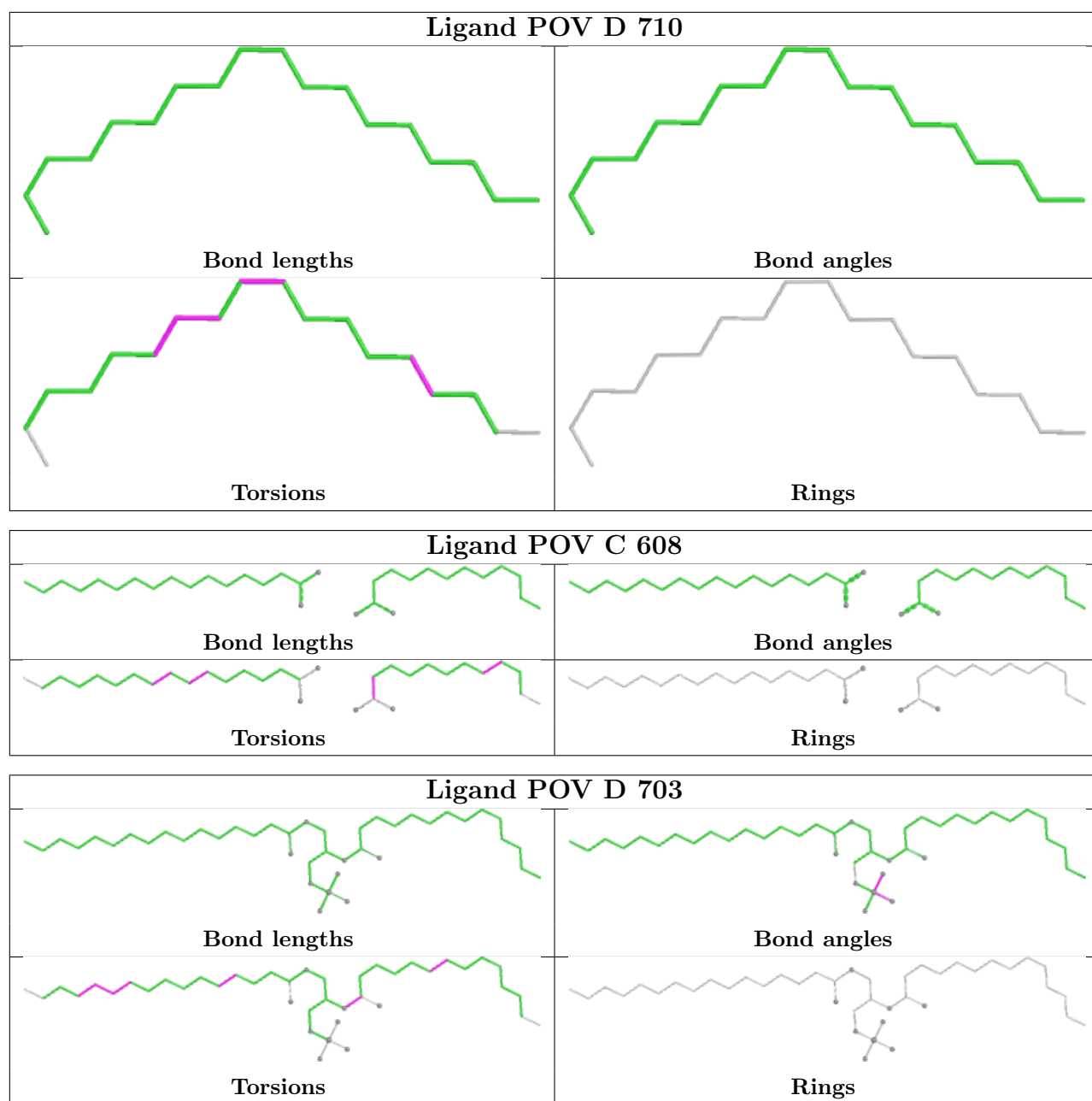


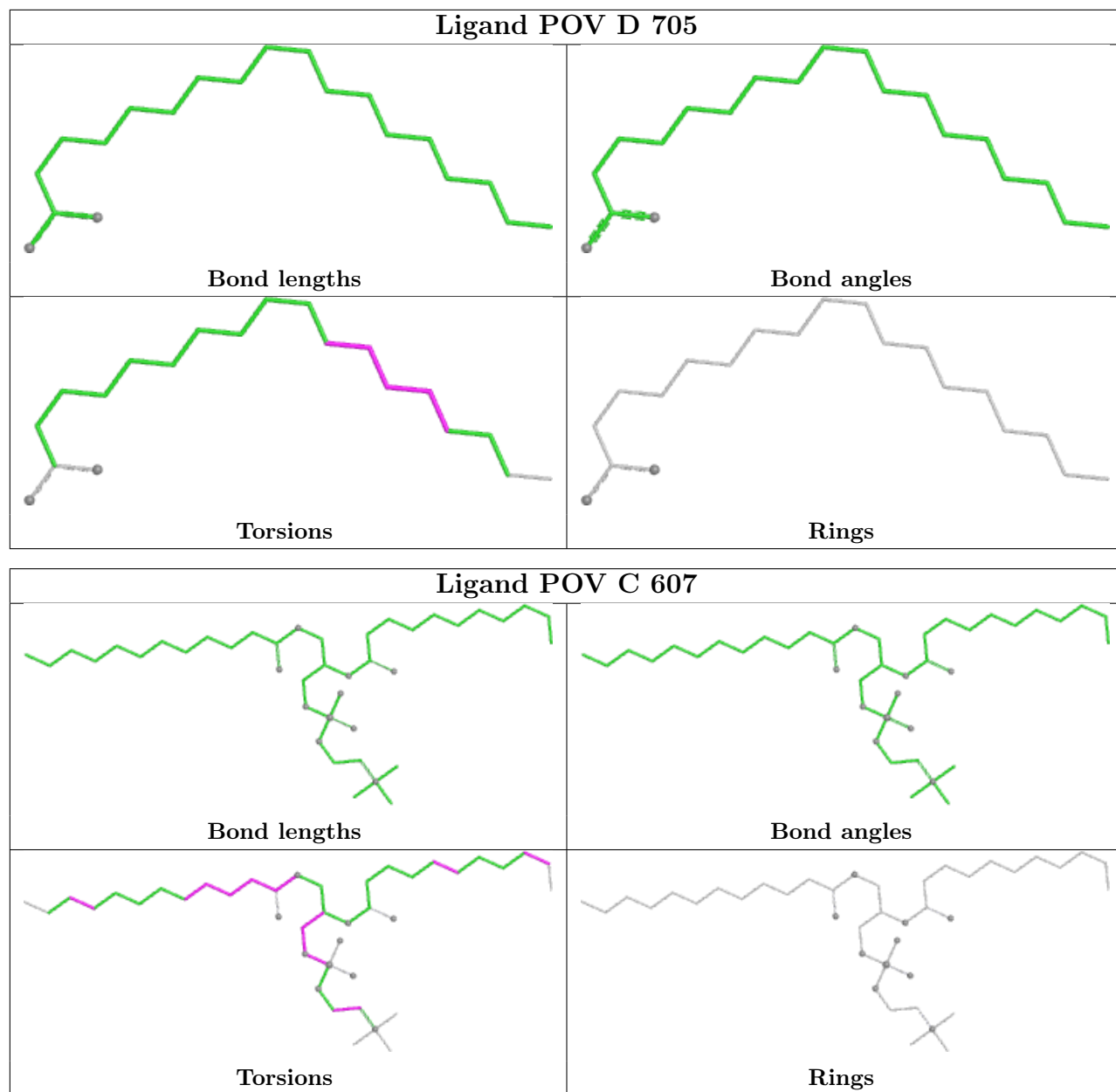


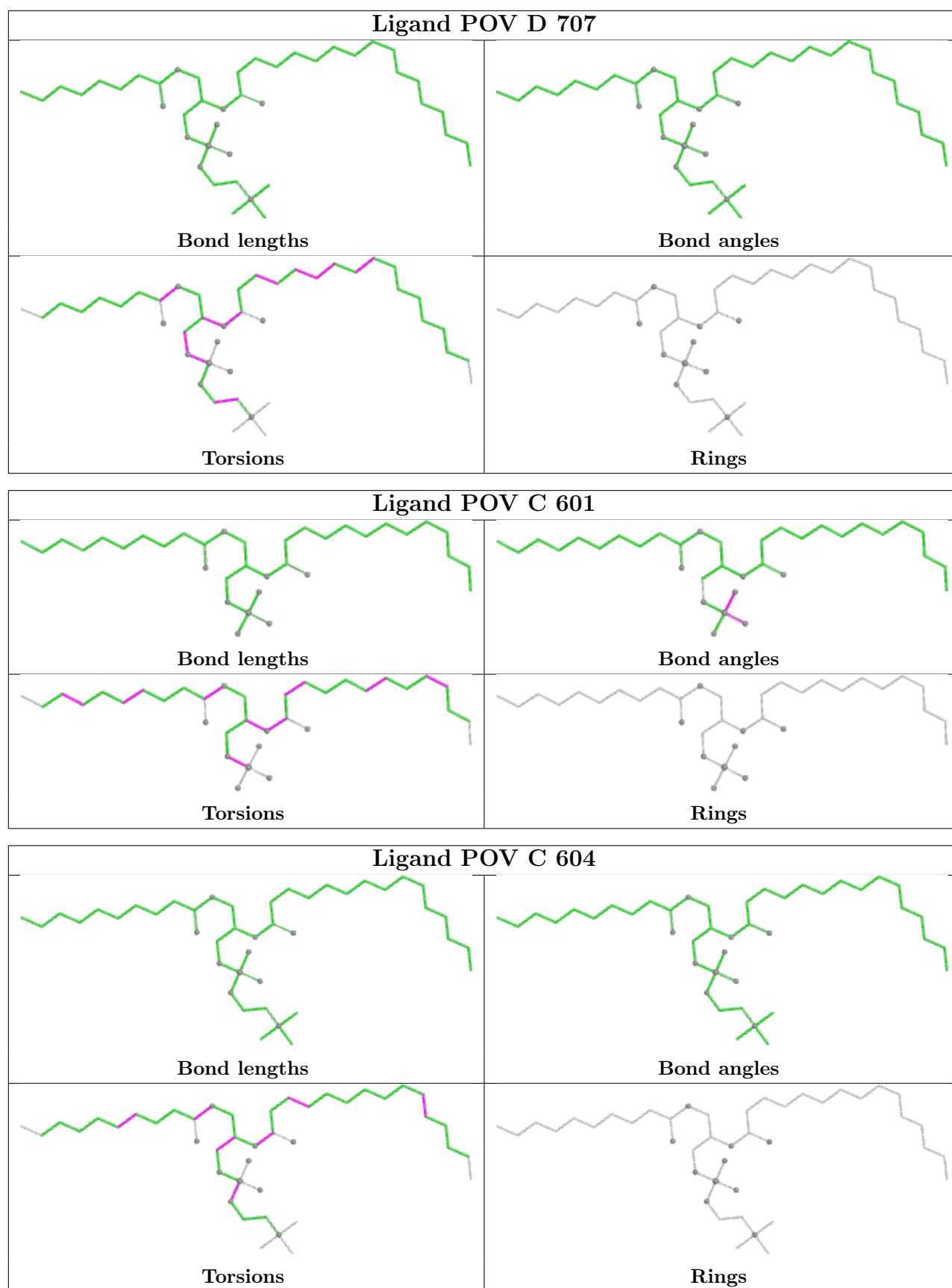


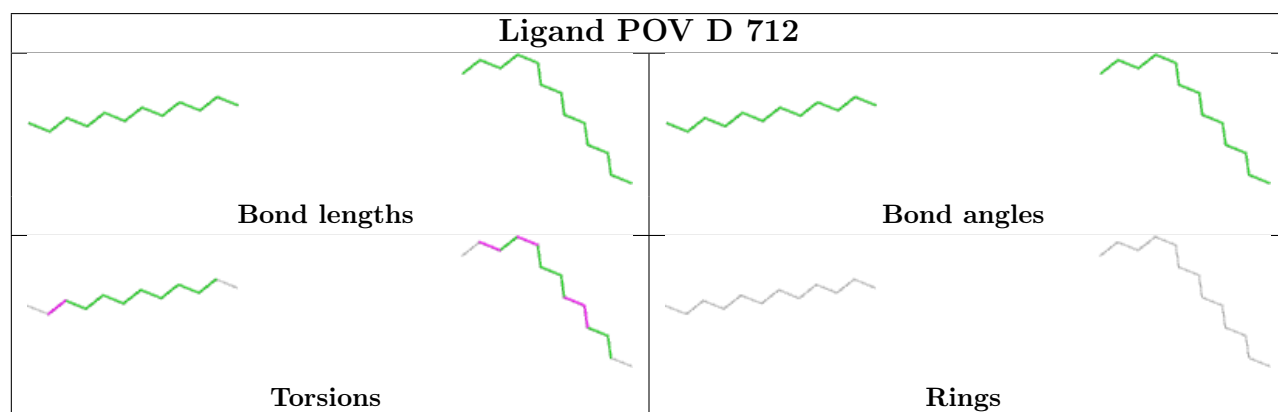
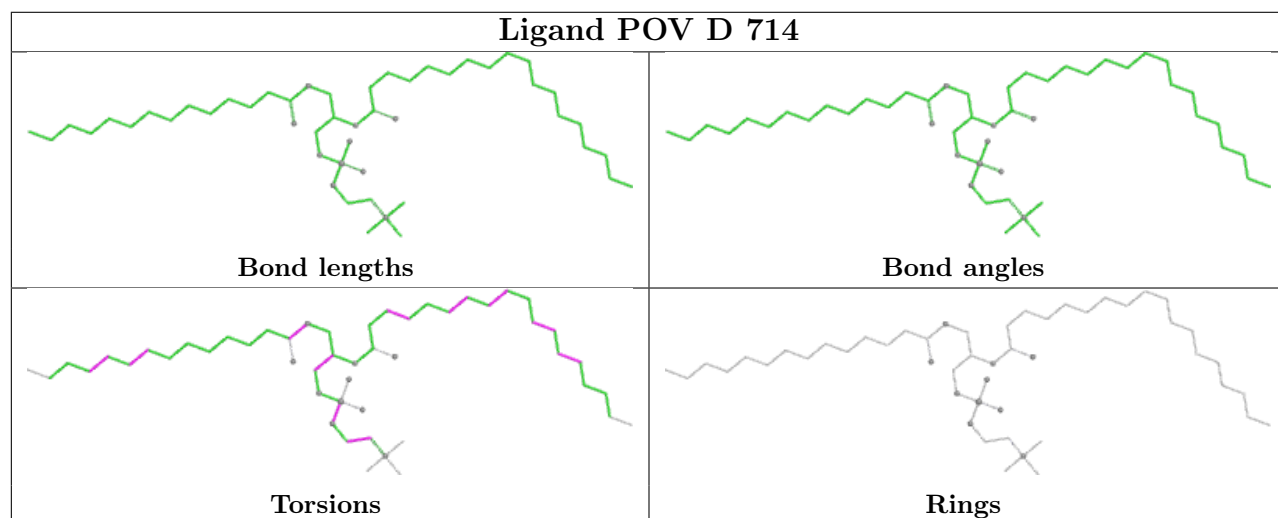
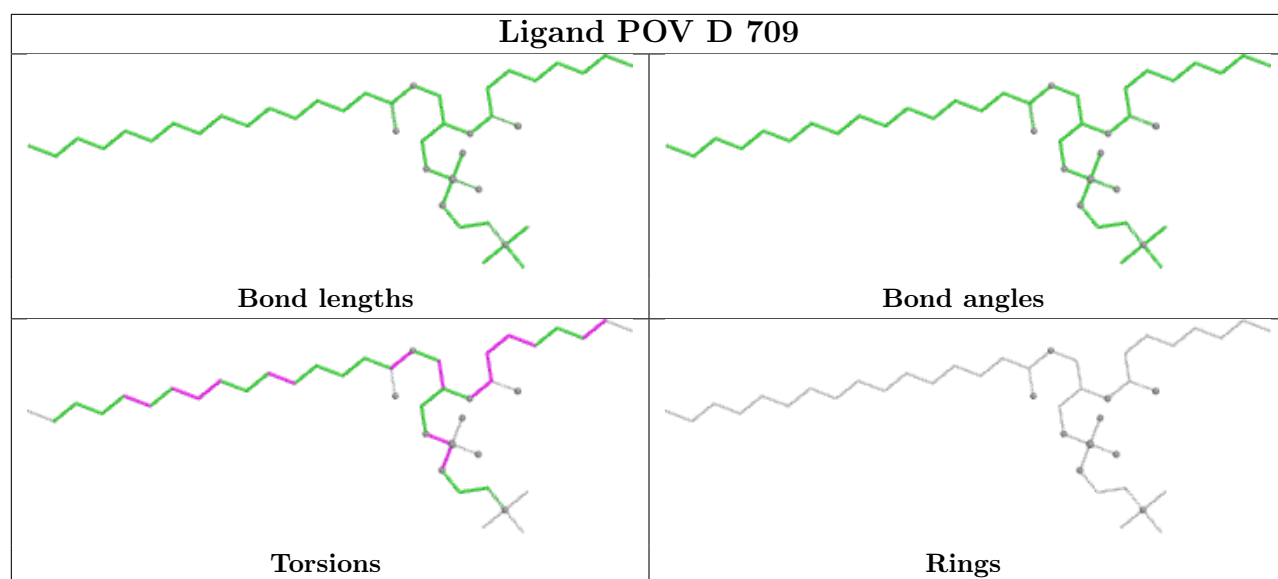


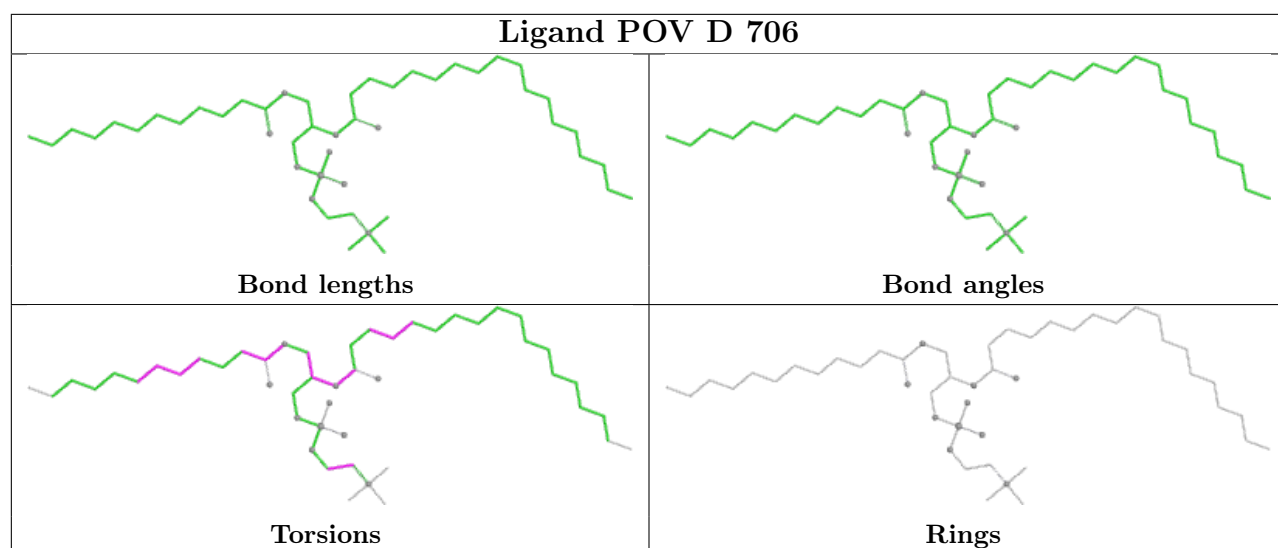












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

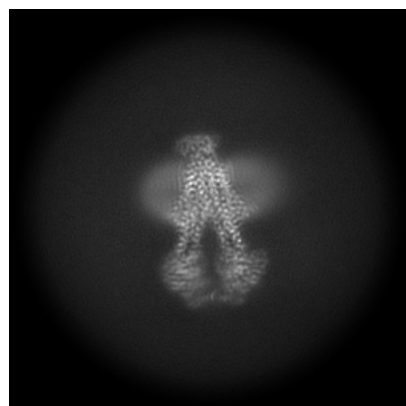
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29362. These allow visual inspection of the internal detail of the map and identification of artifacts.

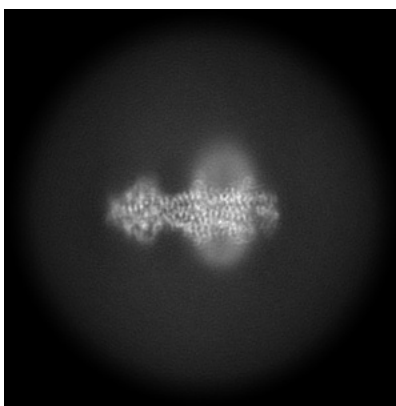
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

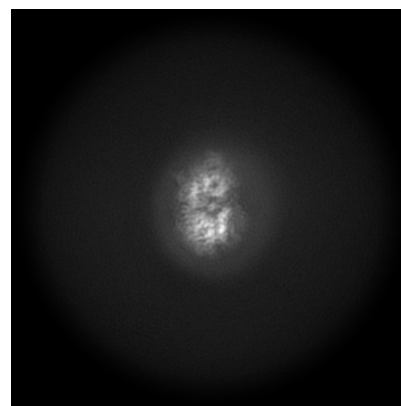
6.1.1 Primary map



X

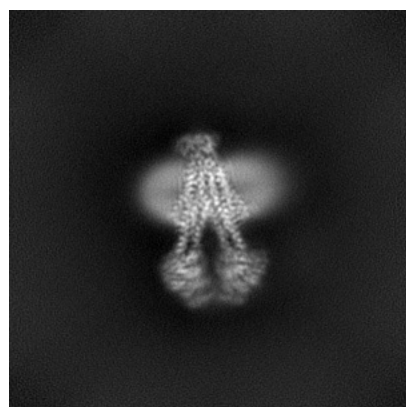


Y

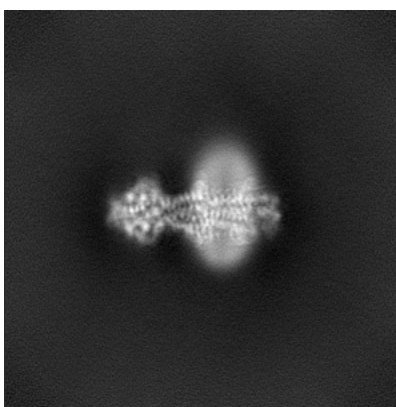


Z

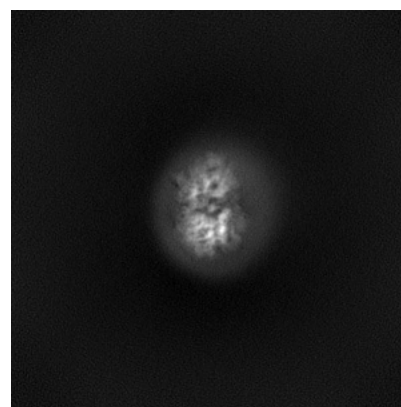
6.1.2 Raw map



X



Y

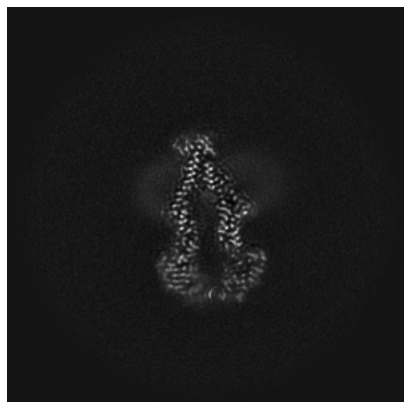


Z

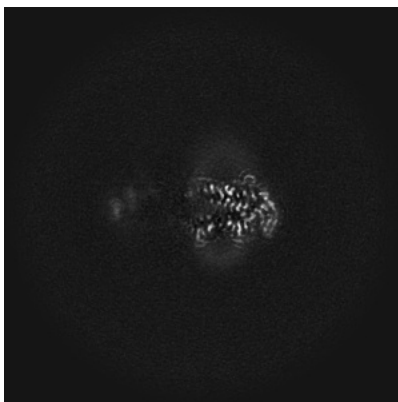
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

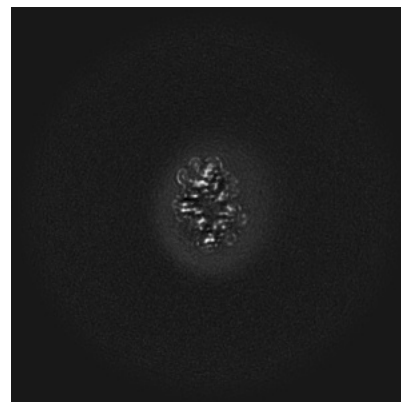
6.2.1 Primary map



X Index: 150

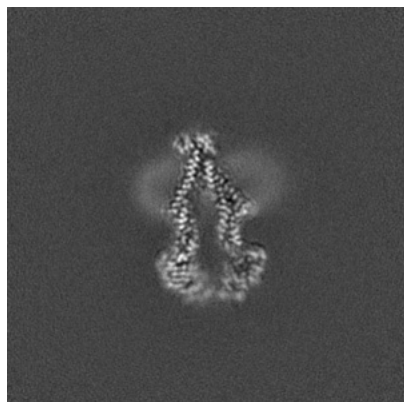


Y Index: 150

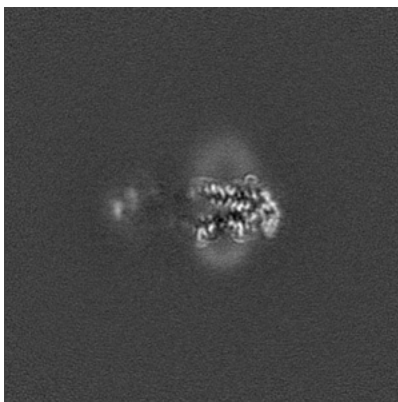


Z Index: 150

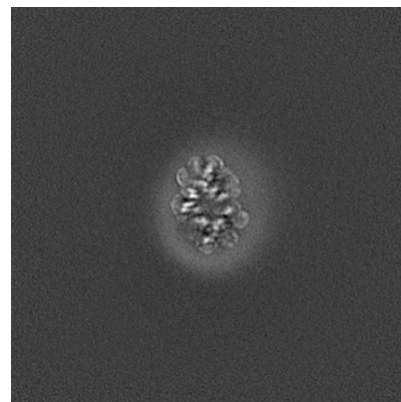
6.2.2 Raw map



X Index: 150



Y Index: 150

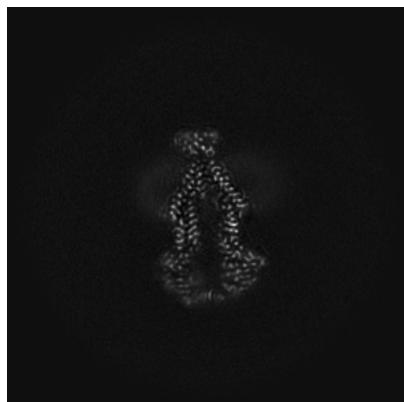


Z Index: 150

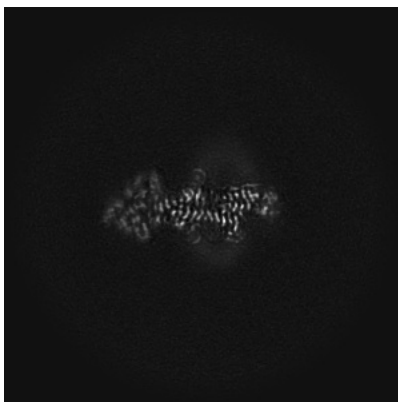
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

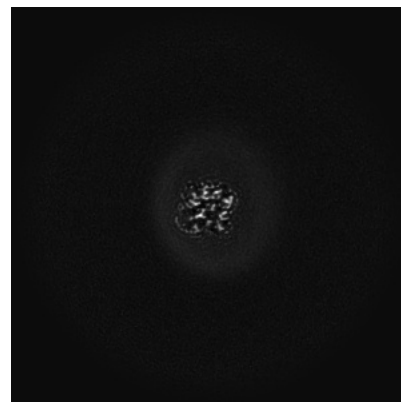
6.3.1 Primary map



X Index: 146

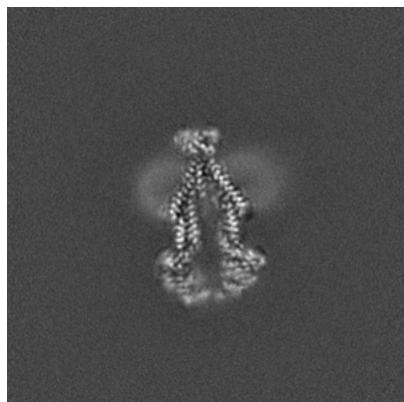


Y Index: 137

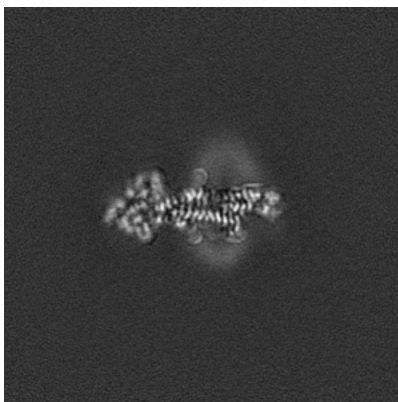


Z Index: 171

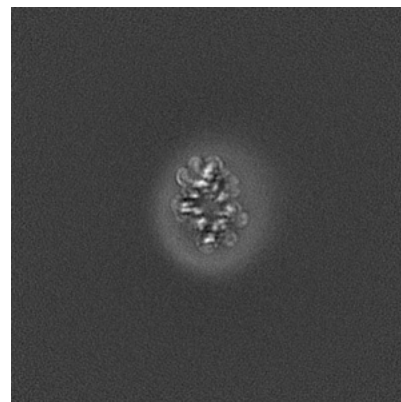
6.3.2 Raw map



X Index: 146



Y Index: 137

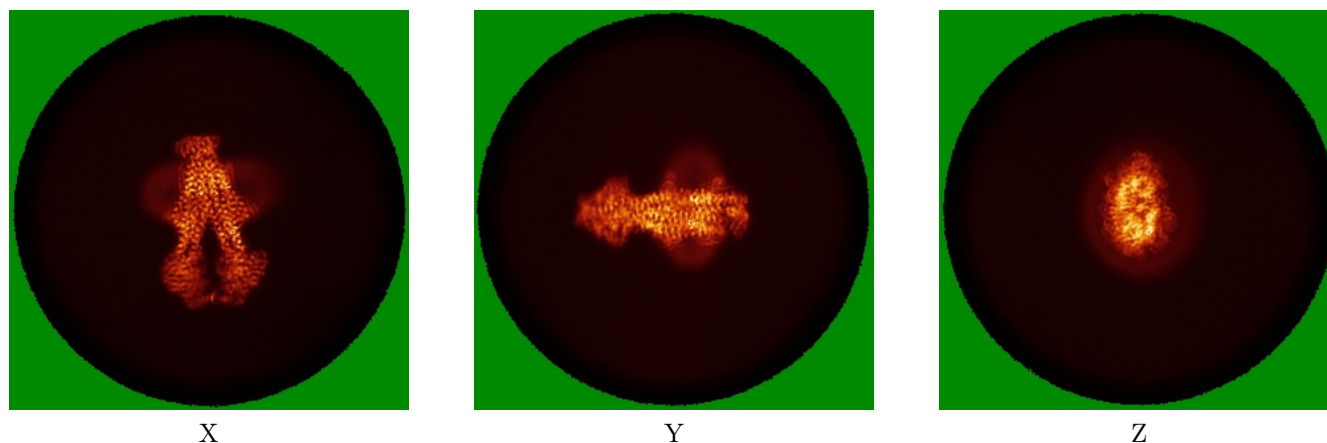


Z Index: 151

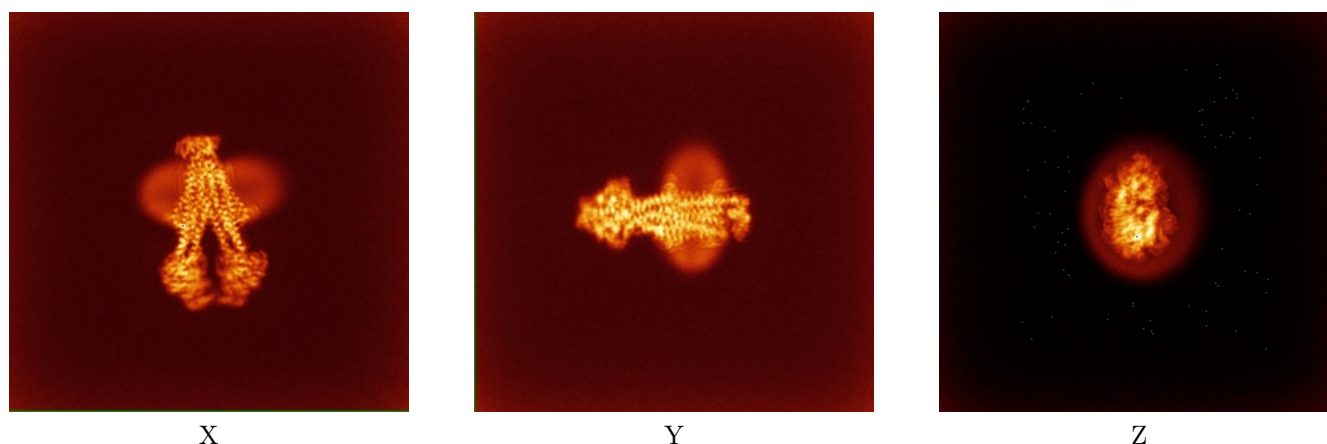
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

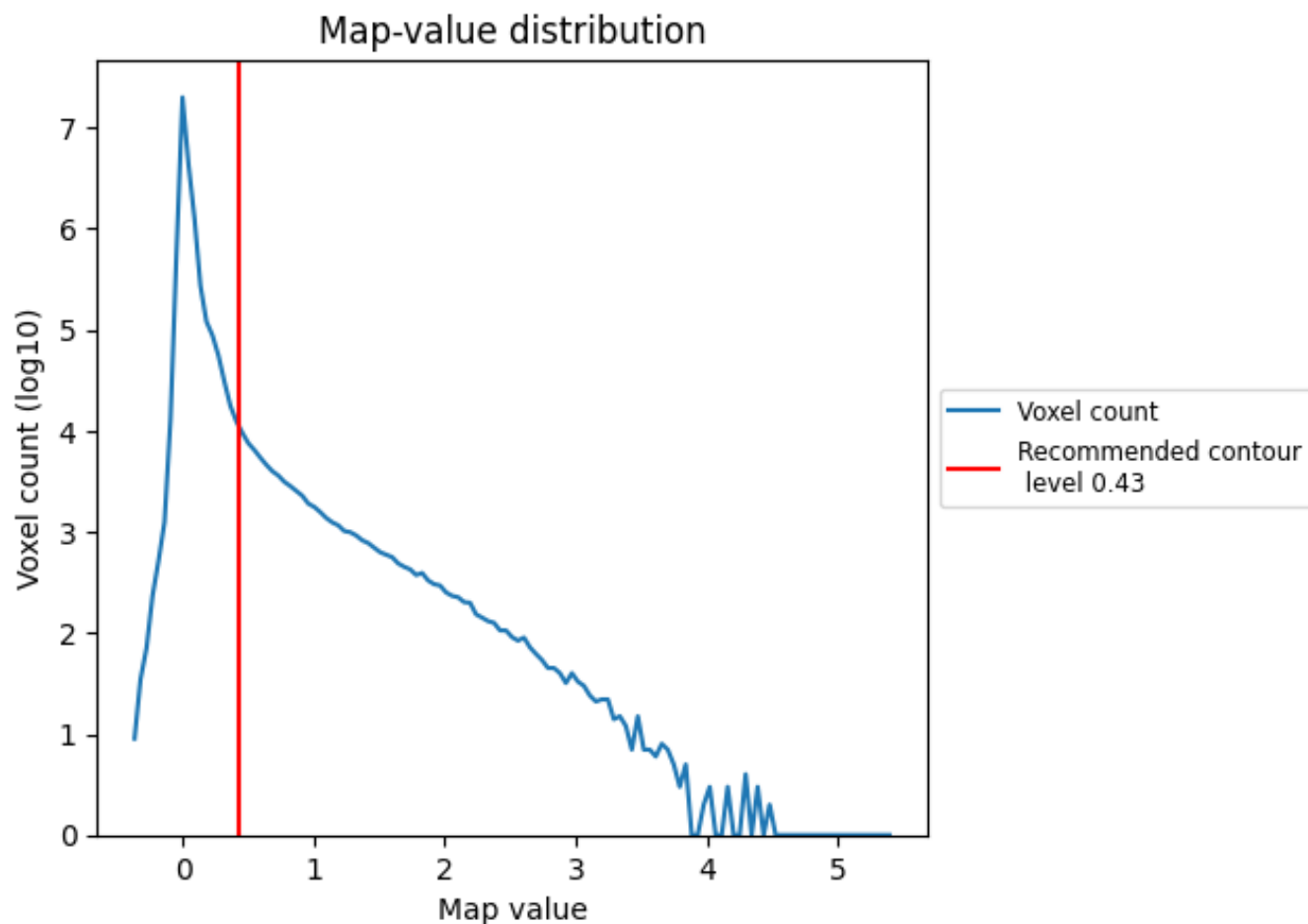
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

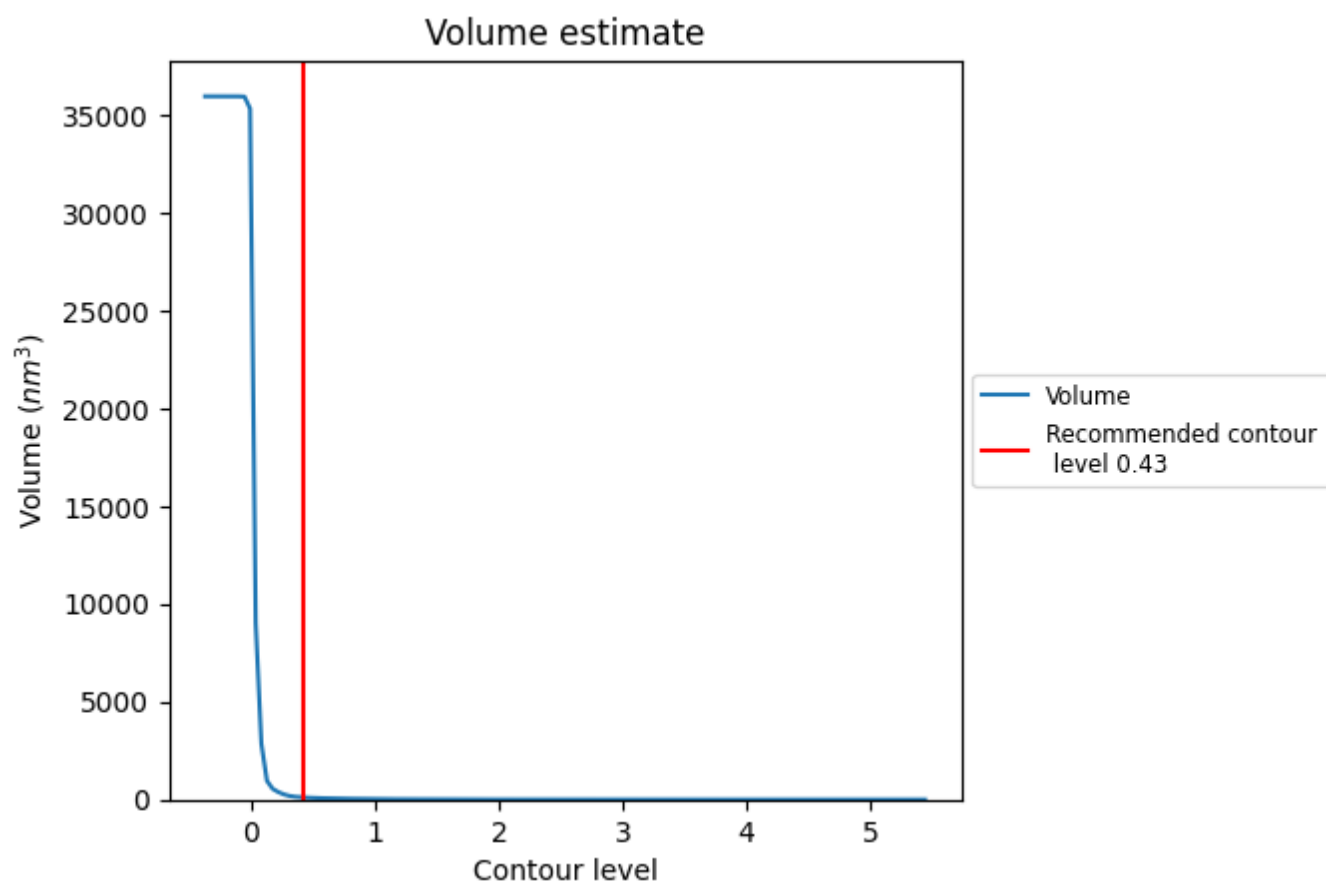
This section contains the results of statistical analysis of the map.

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.

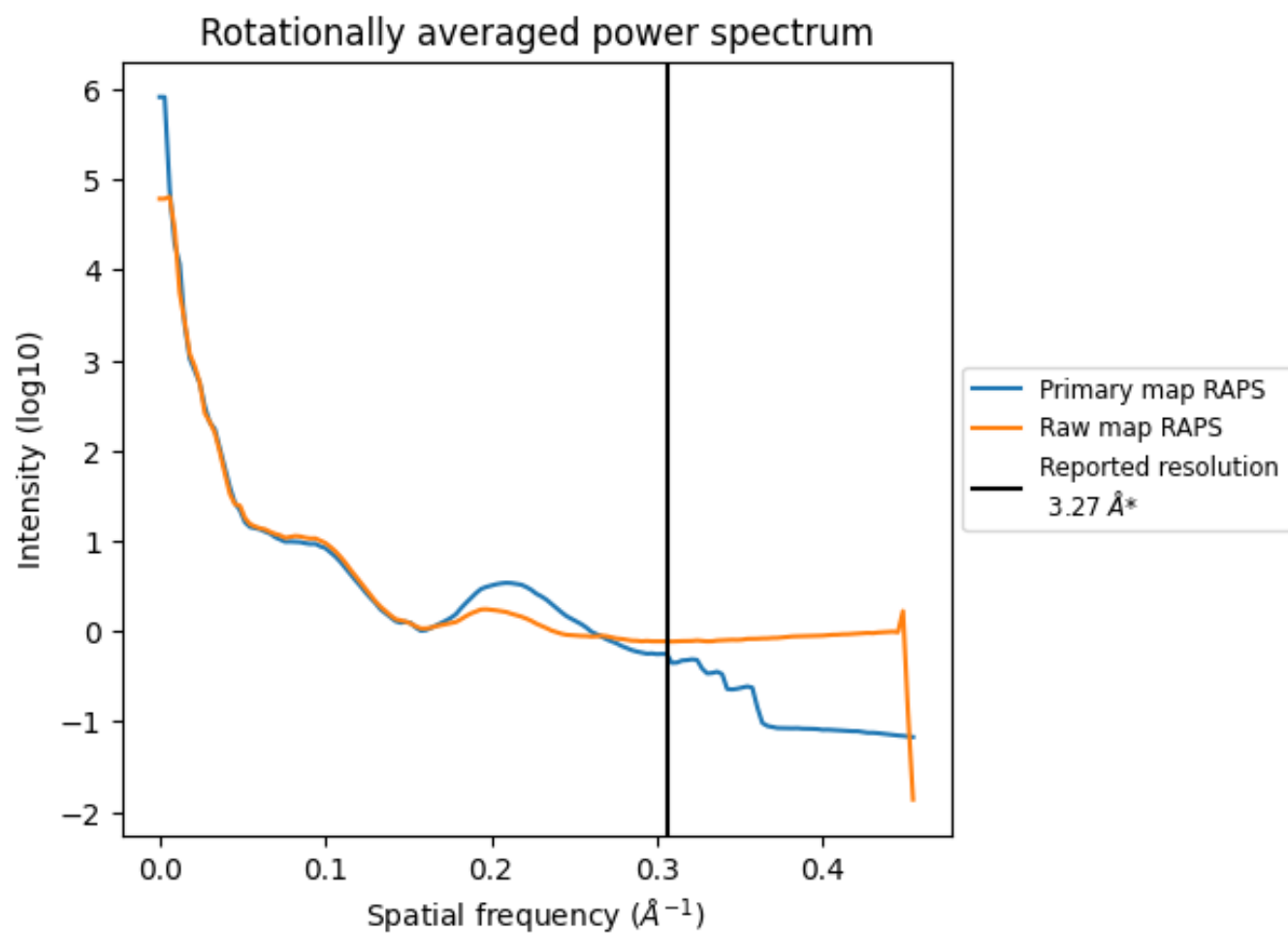
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 107 nm³; this corresponds to an approximate mass of 97 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 ⓘ

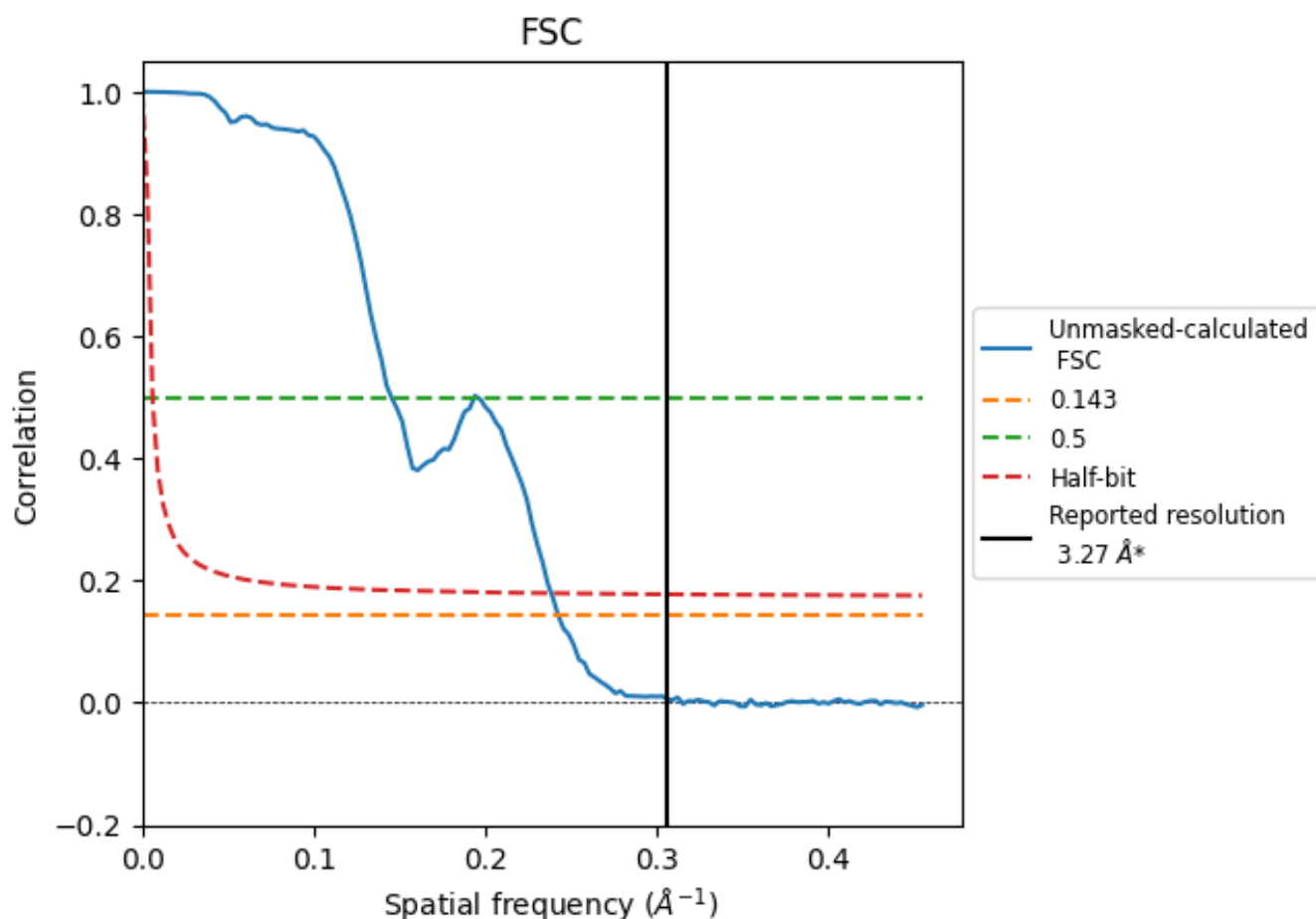


*Reported resolution corresponds to spatial frequency of $0.306 \text{ \AA}^{-1}$

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 $0.306 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

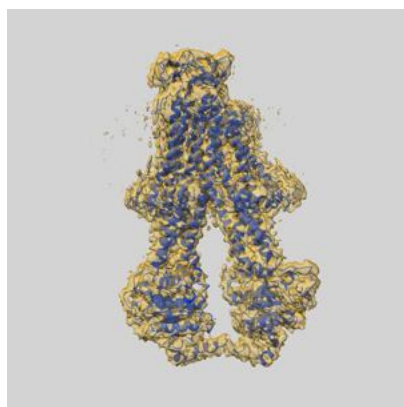
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.27	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.13	6.89	4.19

*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 4.13 differs from the reported value 3.27 by more than 10 %

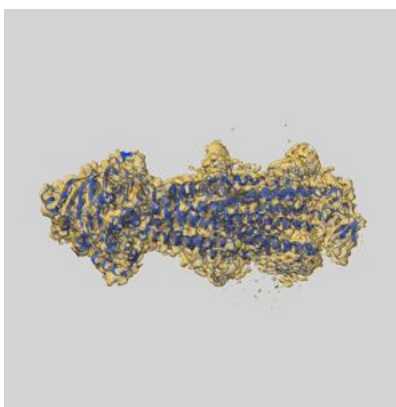
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29362 and PDB model 8FPF. Per-residue inclusion information can be found in section [3](#) on page [7](#).

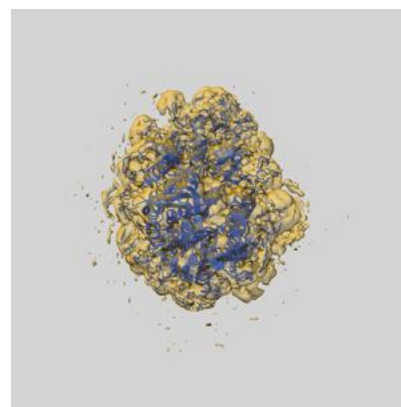
9.1 Map-model overlay [i](#)



X



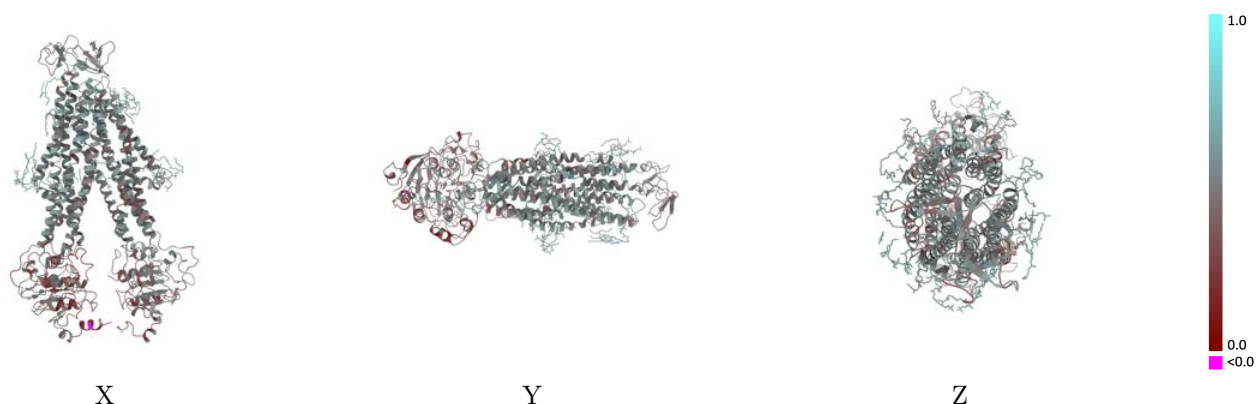
Y



Z

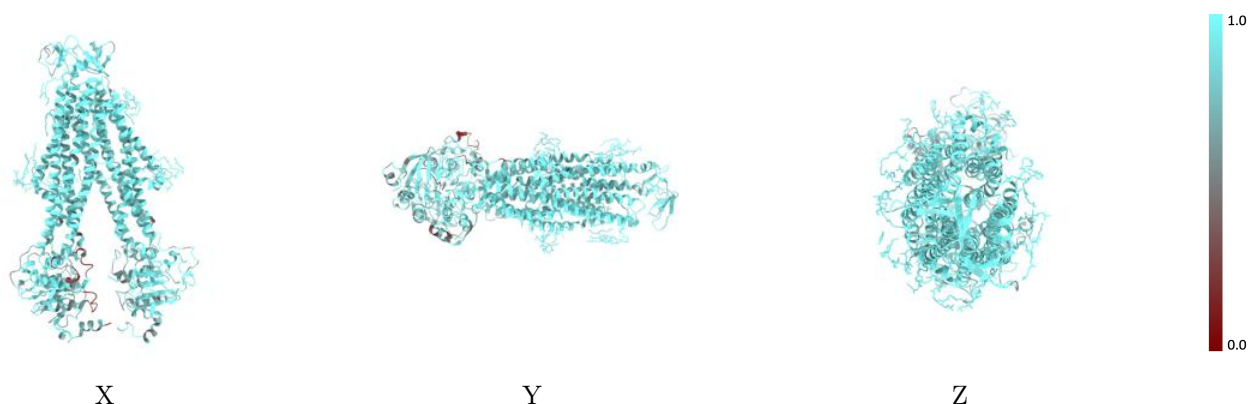
The images above show the 3D surface view of the map at the recommended contour level 0.43 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



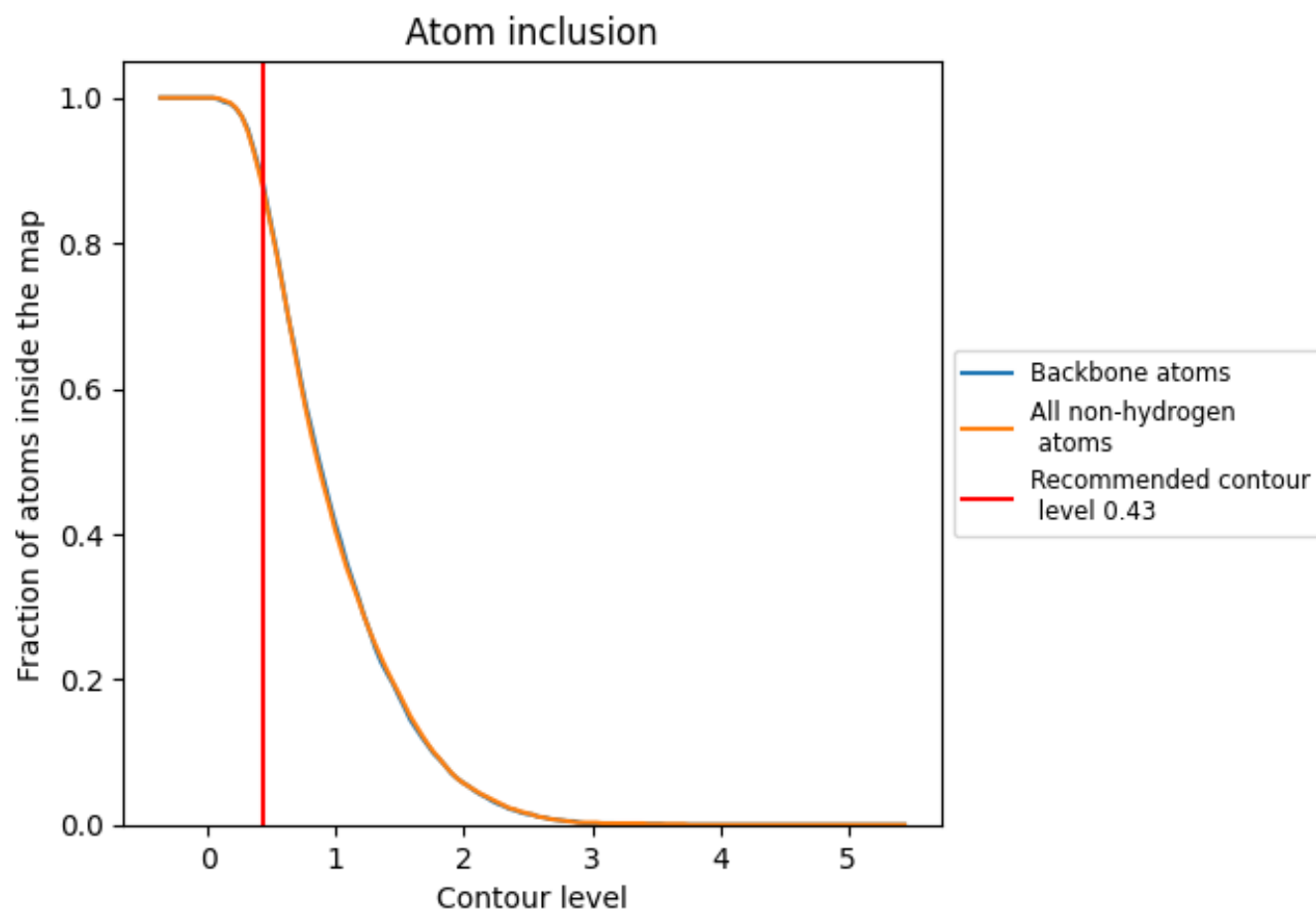
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.43).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 88% 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 (0.43) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8770	0.4590
C	0.8870	0.4660
D	0.8720	0.4540

