



Full wwPDB EM Validation Report ⓘ

Mar 7, 2026 – 11:13 PM UTC

PDB ID : 9UDF / pdb_00009udf
EMDB ID : EMD-64068
Title : Cryo-EM structure of Na⁺-translocating NADH-ubiquinone oxidoreductase NqrB-G141A mutant from *Vibrio cholerae* reduced by NADH, with bound korormicin A, shifted state
Authors : Ishikawa-Fukuda, M.; Kishikawa, J.; Kato, T.; Murai, M.
Deposited on : 2025-04-06
Resolution : 2.93 Å (reported)
Based on initial model : 9LRR

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:

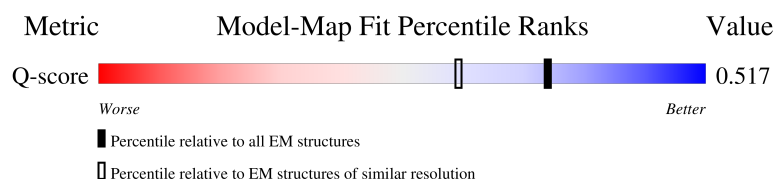
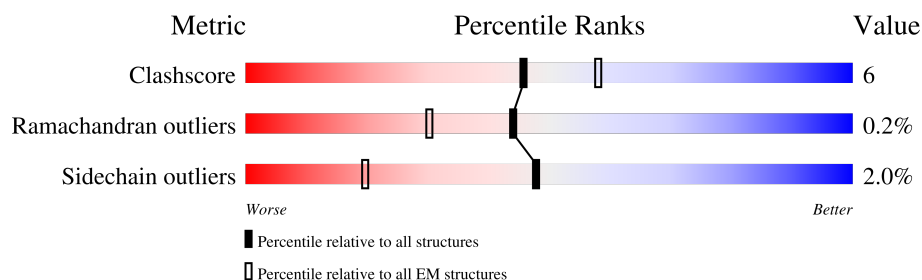
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 2.93 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	13037 (2.43 - 3.43)

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

Mol	Chain	Length	Quality of chain
1	A	446	
2	B	415	
3	C	257	

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Mol	Chain	Length	Quality of chain
4	D	210	86% 11%
5	E	198	82% 17%
6	F	414	35% 76% 22%

2 Entry composition

There are 15 unique types of molecules in this entry. The entry contains 15124 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Na(+)-translocating NADH-quinone reductase subunit A.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	446	Total	C	N	O	S	0	0
			3416	2165	584	650	17		

- Molecule 2 is a protein called Na(+)-translocating NADH-quinone reductase subunit B.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	412	Total	C	N	O	S	0	0
			3178	2103	517	536	22		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	141	ALA	GLY	engineered mutation	UNP A5F5X0

- Molecule 3 is a protein called Na(+)-translocating NADH-quinone reductase subunit C.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	252	Total	C	N	O	S	0	0
			1911	1208	329	370	4		

- Molecule 4 is a protein called Na(+)-translocating NADH-quinone reductase subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	206	Total	C	N	O	S	0	0
			1581	1048	250	273	10		

- Molecule 5 is a protein called Na(+)-translocating NADH-quinone reductase subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	198	Total	C	N	O	S	0	0
			1512	1013	230	258	11		

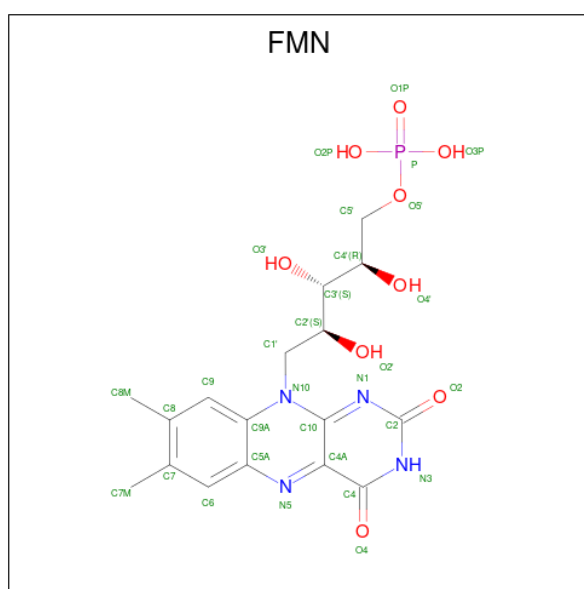
- Molecule 6 is a protein called Na(+)-translocating NADH-quinone reductase subunit F.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	408	Total	C	N	O	S	0	0
			3157	2021	517	595	24		

There are 6 discrepancies between the modelled and reference sequences:

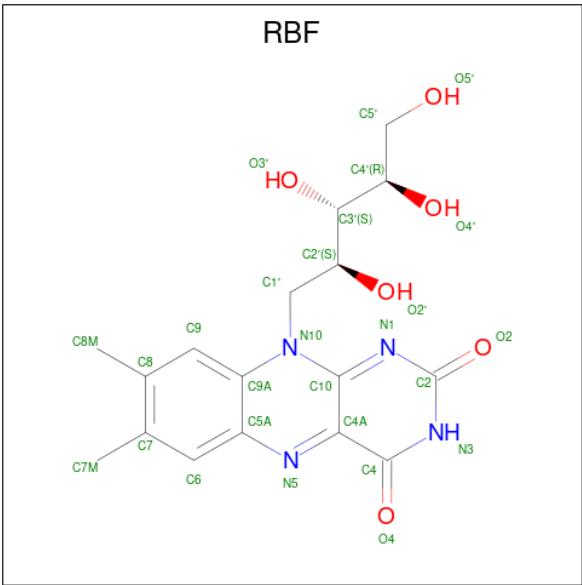
Chain	Residue	Modelled	Actual	Comment	Reference
F	409	HIS	-	expression tag	UNP A5F5Y4
F	410	HIS	-	expression tag	UNP A5F5Y4
F	411	HIS	-	expression tag	UNP A5F5Y4
F	412	HIS	-	expression tag	UNP A5F5Y4
F	413	HIS	-	expression tag	UNP A5F5Y4
F	414	HIS	-	expression tag	UNP A5F5Y4

- Molecule 7 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



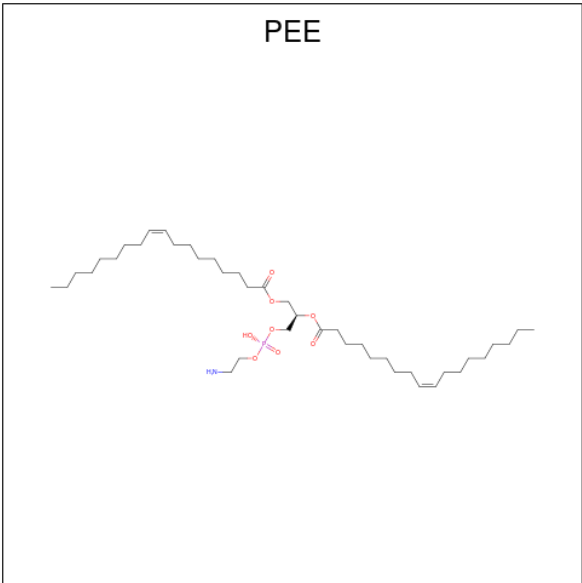
Mol	Chain	Residues	Atoms					AltConf
7	B	1	Total	C	N	O	P	0
			30	17	4	8	1	
7	C	1	Total	C	N	O	P	0
			30	17	4	8	1	

- Molecule 8 is RIBOFLAVIN (CCD ID: RBF) (formula: $C_{17}H_{20}N_4O_6$).



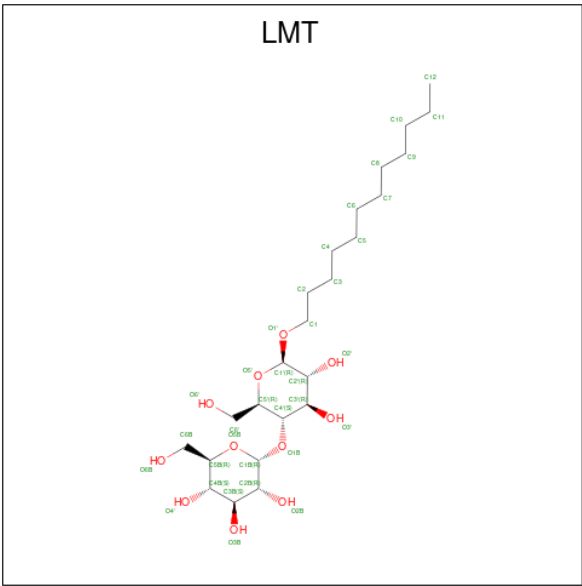
Mol	Chain	Residues	Atoms				AltConf
8	B	1	Total	C	N	O	0
			27	17	4	6	

- Molecule 9 is 1,2-dioleoyl-sn-glycero-3-phosphoethanolamine (CCD ID: PEE) (formula: $C_{41}H_{78}NO_8P$).



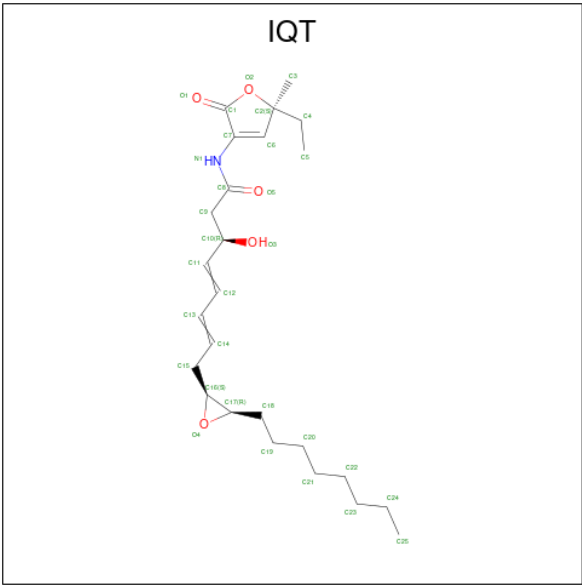
Mol	Chain	Residues	Atoms					AltConf
9	B	1	Total	C	N	O	P	0
			51	41	1	8	1	
9	B	1	Total	C	N	O	P	0
			51	41	1	8	1	

- Molecule 10 is DODECYL-BETA-D-MALTOSE (CCD ID: LMT) (formula: $C_{24}H_{46}O_{11}$).



Mol	Chain	Residues	Atoms			AltConf
10	B	1	Total	C	O	0
			35	24	11	
10	D	1	Total	C	O	0
			35	24	11	

- Molecule 11 is Korormicin (CCD ID: IQT) (formula: $C_{25}H_{39}NO_5$).

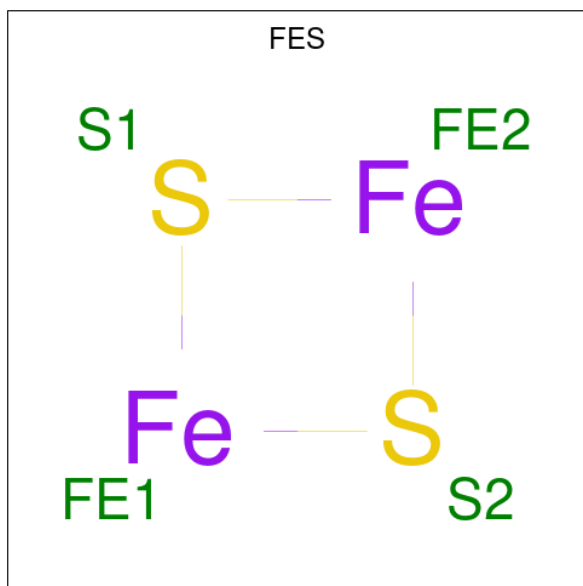


Mol	Chain	Residues	Atoms				AltConf
11	B	1	Total	C	N	O	0
			31	25	1	5	

- Molecule 12 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
12	C	1	Total	Ca	0
			1	1	

- Molecule 13 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



Mol	Chain	Residues	Atoms			AltConf
13	D	1	Total	Fe	S	0
			4	2	2	
13	F	1	Total	Fe	S	0
			4	2	2	

- Molecule 14 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).

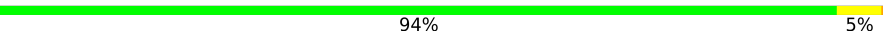
- Molecule 15 is water.

- | | | | | |
|----|---|---|---------|---|
| 15 | Λ | 7 | Total O | 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 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.

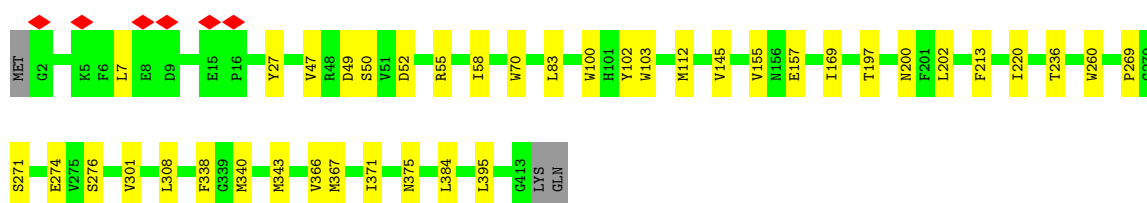
- Molecule 1: Na(+)-translocating NADH-quinone reductase subunit A

Chain A: 



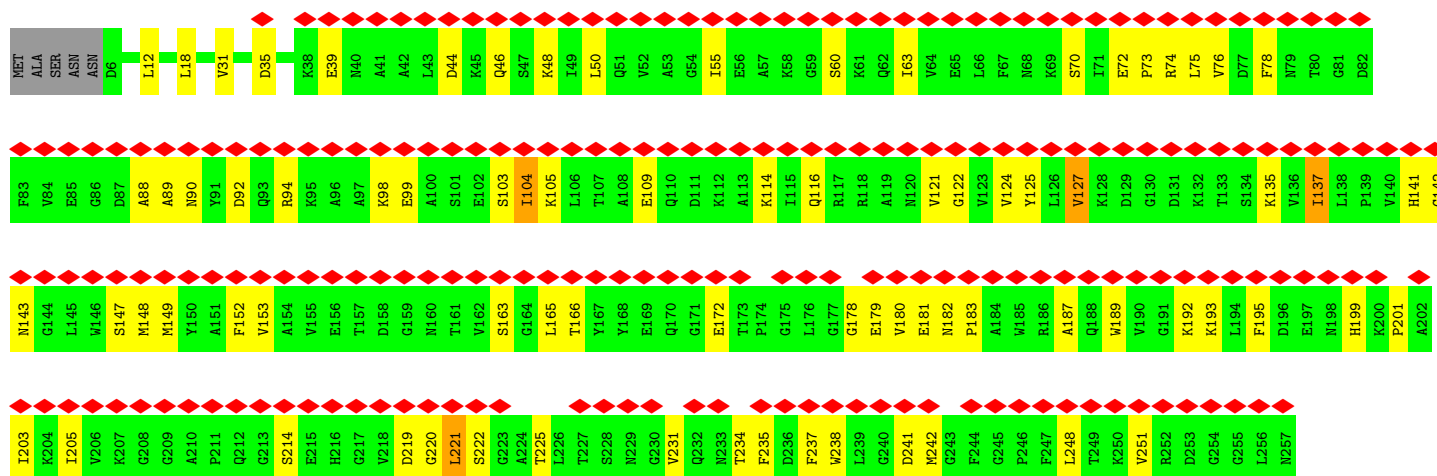
- Molecule 2: Na(+)-translocating NADH-quinone reductase subunit B

Chain B: 



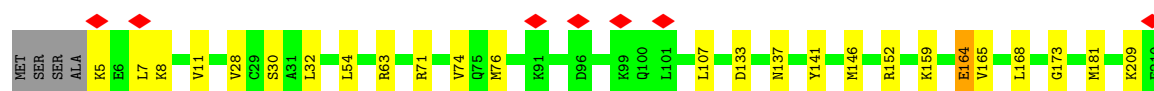
- Molecule 3: Na(+)-translocating NADH-quinone reductase subunit C

Chain C: 



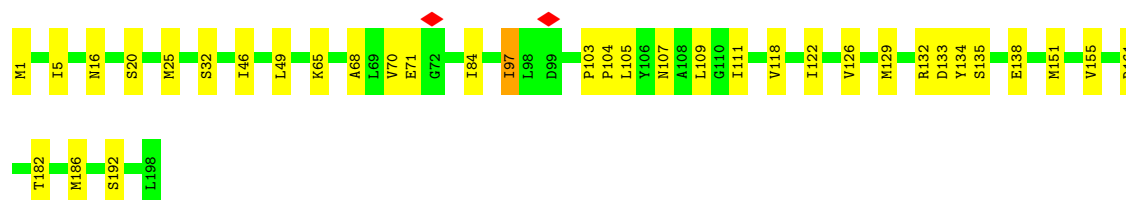
- Molecule 4: Na(+)-translocating NADH-quinone reductase subunit D

Chain D:  86% 11%



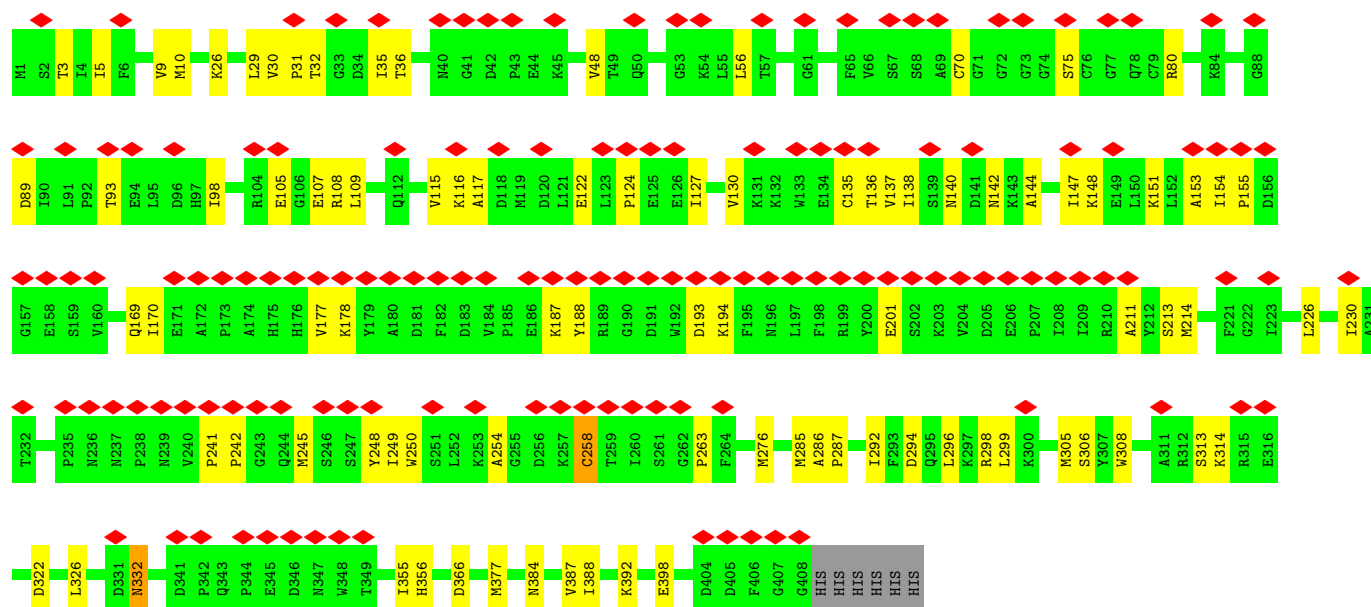
- Molecule 5: Na(+)-translocating NADH-quinone reductase subunit E

Chain E:  82% 17%



- Molecule 6: Na(+)-translocating NADH-quinone reductase subunit F

Chain F:  35% 76% 22%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	74083	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	65	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	4.093	Depositor
Minimum map value	-2.406	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.102	Depositor
Recommended contour level	0.488	Depositor
Map size ($\text{\AA}$)	281.6, 281.6, 281.6	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.88, 0.88, 0.88	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: FMN, LMT, FES, FAD, PEE, CA, IQT, RBF

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	0.15	0/3480	0.30	0/4715
2	B	0.15	0/3276	0.32	0/4459
3	C	0.18	0/1943	0.52	0/2621
4	D	0.15	0/1613	0.35	0/2187
5	E	0.17	0/1545	0.33	0/2094
6	F	0.14	0/3235	0.38	0/4381
All	All	0.15	0/15092	0.37	0/20457

There are no bond length outliers.

There are no bond angle outliers.

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	A	3416	0	3463	15	0
2	B	3178	0	3157	31	0
3	C	1911	0	1917	52	0
4	D	1581	0	1667	15	0
5	E	1512	0	1586	20	0
6	F	3157	0	3080	56	0
7	B	30	0	19	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	30	0	19	3	0
8	B	27	0	20	3	0
9	B	102	0	164	3	0
10	B	35	0	46	2	0
10	D	35	0	46	1	0
11	B	31	0	0	2	0
12	C	1	0	0	0	0
13	D	4	0	0	0	0
13	F	4	0	0	0	0
14	F	53	0	31	1	0
15	A	7	0	0	0	0
15	B	10	0	0	1	0
All	All	15124	0	15215	189	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 (189) 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)
3:C:237:PHE:HD2	3:C:242:MET:HE1	1.41	0.84
6:F:170:ILE:HG23	6:F:245:MET:HE3	1.61	0.80
3:C:109:GLU:HG3	3:C:242:MET:HE2	1.66	0.76
3:C:221:LEU:HD23	3:C:222:SER:H	1.54	0.73
4:D:76:MET:HE2	5:E:84:ILE:HG12	1.71	0.72
3:C:149:MET:HE2	3:C:181:GLU:HG2	1.74	0.70
5:E:71:GLU:N	5:E:71:GLU:OE2	2.23	0.70
6:F:142:ASN:OD1	6:F:148:LYS:NZ	2.25	0.70
3:C:178:GLY:HA3	3:C:221:LEU:HG	1.73	0.68
6:F:169:GLN:HA	6:F:211:ALA:HA	1.77	0.66
6:F:245:MET:O	6:F:249:ILE:HG12	1.97	0.65
6:F:70:CYS:HG	6:F:75:SER:HG	1.41	0.64
2:B:157:GLU:HG2	11:B:506:IQT:C5	2.28	0.64
6:F:142:ASN:ND2	6:F:144:ALA:O	2.32	0.62
4:D:30:SER:HG	4:D:141:TYR:HH	1.45	0.61
2:B:271:SER:HB2	2:B:274:GLU:HG2	1.84	0.60
1:A:277:THR:OG1	1:A:278:VAL:N	2.34	0.59
1:A:295:GLU:OE2	1:A:330:ARG:NH1	2.36	0.59
3:C:35:ASP:O	3:C:39:GLU:HG2	2.03	0.59
5:E:132:ARG:NH1	5:E:134:TYR:OH	2.35	0.59
2:B:55:ARG:HA	2:B:58:ILE:HG22	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:197:THR:HG23	8:B:502:RBF:HC52	1.84	0.59
3:C:127:VAL:HG21	3:C:135:LYS:HB2	1.83	0.59
3:C:165:LEU:HB2	3:C:182:ASN:ND2	2.19	0.58
6:F:98:ILE:HD11	6:F:109:LEU:HG	1.85	0.58
3:C:183:PRO:HD2	3:C:187:ALA:HA	1.86	0.58
6:F:187:LYS:HE2	6:F:188:TYR:CE1	2.39	0.57
1:A:97:ASP:N	1:A:97:ASP:OD1	2.36	0.57
3:C:125:TYR:HB3	3:C:137:ILE:HG22	1.85	0.56
3:C:180:VAL:HB	3:C:220:GLY:HA2	1.87	0.56
1:A:233:ASP:OD1	1:A:233:ASP:N	2.37	0.56
1:A:396:ASP:OD2	1:A:404:SER:OG	2.15	0.56
3:C:237:PHE:CD2	3:C:242:MET:HE1	2.32	0.54
5:E:1:MET:O	5:E:5:ILE:HD12	2.08	0.54
6:F:142:ASN:HA	6:F:148:LYS:HG3	1.90	0.54
2:B:83:LEU:HD13	2:B:112:MET:HE1	1.89	0.54
6:F:377:MET:HE1	6:F:388:ILE:HG13	1.91	0.53
1:A:260:ARG:HB2	1:A:277:THR:HG23	1.91	0.53
3:C:72:GLU:HB3	3:C:74:ARG:HH21	1.74	0.53
6:F:306:SER:OG	6:F:308:TRP:NE1	2.42	0.53
3:C:135:LYS:HB3	3:C:152:PHE:CE1	2.43	0.52
1:A:309:ALA:HB1	1:A:314:ALA:HA	1.90	0.52
5:E:135:SER:OG	5:E:138:GLU:OE2	2.28	0.52
6:F:10:MET:HE3	6:F:10:MET:HA	1.91	0.52
3:C:60:SER:HA	3:C:63:ILE:HG22	1.90	0.52
2:B:213:PHE:CZ	7:C:301:FMN:H6	2.45	0.51
3:C:165:LEU:HB2	3:C:182:ASN:HD21	1.74	0.51
6:F:356:HIS:HB3	6:F:387:VAL:HG22	1.92	0.51
3:C:201:PRO:HG2	3:C:238:TRP:HB3	1.93	0.51
3:C:75:LEU:HB2	3:C:88:ALA:HB2	1.92	0.51
6:F:308:TRP:HB3	6:F:355:ILE:HD11	1.92	0.51
5:E:107:ASN:C	5:E:109:LEU:H	2.19	0.50
6:F:154:ILE:HG13	6:F:155:PRO:HD2	1.93	0.50
6:F:213:SER:OG	14:F:502:FAD:H6	2.11	0.50
6:F:105:GLU:N	6:F:105:GLU:OE1	2.45	0.49
6:F:313:SER:OG	6:F:314:LYS:N	2.45	0.49
6:F:130:VAL:HA	6:F:263:PRO:HD2	1.94	0.49
5:E:49:LEU:HD13	5:E:122:ILE:HG23	1.94	0.49
6:F:136:THR:HB	6:F:153:ALA:HB3	1.93	0.49
5:E:16:ASN:H	5:E:20:SER:HB2	1.76	0.49
4:D:133:ASP:OD1	4:D:137:ASN:ND2	2.46	0.49
2:B:100:TRP:HA	2:B:103:TRP:NE1	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:115:VAL:O	6:F:116:LYS:HD3	2.13	0.48
3:C:18:LEU:HD22	4:D:74:VAL:HG13	1.94	0.48
3:C:78:PHE:HB2	3:C:122:GLY:HA3	1.94	0.48
6:F:193:ASP:OD1	6:F:194:LYS:N	2.46	0.48
2:B:7:LEU:O	2:B:27:TYR:OH	2.28	0.48
4:D:168:LEU:O	4:D:173:GLY:N	2.45	0.48
3:C:152:PHE:HB3	3:C:163:SER:HB2	1.95	0.48
3:C:73:PRO:HB3	3:C:89:ALA:N	2.28	0.48
5:E:129:MET:HE3	5:E:129:MET:HB3	1.73	0.48
4:D:5:LYS:HG3	4:D:7:LEU:H	1.77	0.48
3:C:75:LEU:HD21	3:C:121:VAL:HB	1.96	0.48
6:F:306:SER:HG	6:F:308:TRP:NE1	2.12	0.48
2:B:169:ILE:HD13	2:B:343:MET:HE1	1.96	0.48
2:B:338:PHE:HE1	2:B:343:MET:HE2	1.78	0.48
6:F:26:LYS:O	6:F:30:VAL:HG22	2.14	0.48
2:B:375:ASN:O	15:B:601:HOH:O	2.20	0.47
3:C:205:ILE:HG23	3:C:219:ASP:HA	1.96	0.47
6:F:31:PRO:O	6:F:32:THR:HG22	2.14	0.47
2:B:384:LEU:HB3	5:E:192:SER:HB3	1.95	0.47
6:F:135:CYS:N	6:F:258:CYS:O	2.45	0.47
9:B:503:PEE:H20	9:B:503:PEE:H13	1.61	0.47
1:A:258:THR:HA	1:A:279:MET:HE3	1.95	0.47
1:A:345:PHE:O	2:B:50:SER:OG	2.16	0.47
2:B:200:ASN:ND2	2:B:202:LEU:O	2.47	0.47
4:D:146:MET:HE3	4:D:146:MET:HB3	1.76	0.47
5:E:182:THR:O	5:E:186:MET:HG3	2.15	0.47
6:F:138:ILE:HB	6:F:151:LYS:HB3	1.96	0.47
6:F:89:ASP:O	6:F:108:ARG:NH2	2.48	0.46
6:F:286:ALA:HB3	6:F:287:PRO:HD3	1.98	0.46
3:C:44:ASP:O	3:C:48:LYS:HB2	2.15	0.46
3:C:72:GLU:HB3	3:C:74:ARG:NH2	2.30	0.46
4:D:8:LYS:HA	4:D:11:VAL:HG22	1.97	0.46
3:C:152:PHE:H	3:C:163:SER:HB2	1.81	0.46
6:F:296:LEU:HD13	6:F:326:LEU:HD22	1.97	0.46
2:B:157:GLU:CD	2:B:157:GLU:H	2.23	0.46
6:F:285:MET:HG3	6:F:285:MET:O	2.15	0.46
3:C:193:LYS:HB2	3:C:214:SER:O	2.16	0.46
5:E:97:ILE:HG12	6:F:29:LEU:HD11	1.97	0.46
6:F:294:ASP:OD1	6:F:298:ARG:NE	2.31	0.46
6:F:178:LYS:HD3	6:F:201:GLU:HB3	1.97	0.46
3:C:103:SER:O	3:C:104:ILE:HG22	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:159:LYS:HG2	4:D:164:GLU:HA	1.99	0.45
3:C:248:LEU:HA	3:C:251:VAL:HG13	1.98	0.45
2:B:236:THR:HG22	2:B:269:PRO:HG2	1.98	0.45
6:F:276:MET:HB2	6:F:305:MET:HG2	1.97	0.45
3:C:199:HIS:HA	3:C:248:LEU:HD11	1.99	0.45
2:B:308:LEU:HD22	2:B:366:VAL:HG13	1.98	0.45
6:F:5:ILE:O	6:F:9:VAL:HG12	2.17	0.45
3:C:92:ASP:OD2	3:C:94:ARG:N	2.47	0.45
3:C:203:ILE:HG21	3:C:235:PHE:HB2	1.99	0.45
6:F:241:PRO:HA	6:F:242:PRO:HD3	1.87	0.45
3:C:114:LYS:HB3	3:C:143:ASN:HB3	1.99	0.45
3:C:195:PHE:HE1	3:C:238:TRP:NE1	2.15	0.45
3:C:55:ILE:HD13	3:C:70:SER:HB3	1.99	0.44
5:E:103:PRO:N	5:E:104:PRO:HD2	2.33	0.44
3:C:141:HIS:HA	3:C:148:MET:HA	2.00	0.44
2:B:213:PHE:HA	2:B:220:ILE:HG21	2.00	0.44
3:C:50:LEU:HD12	3:C:55:ILE:HB	1.99	0.44
4:D:152:ARG:HA	4:D:181:MET:HE3	2.00	0.44
6:F:366:ASP:OD1	6:F:366:ASP:N	2.50	0.44
9:B:505:PEE:H20	9:B:505:PEE:H53	2.00	0.44
4:D:28:VAL:O	4:D:32:LEU:HG	2.17	0.44
2:B:112:MET:HE3	2:B:112:MET:HB3	1.77	0.44
3:C:31:VAL:HG21	6:F:3:THR:HG21	1.99	0.44
3:C:172:GLU:HB2	3:C:179:GLU:OE1	2.17	0.44
6:F:35:ILE:HG21	6:F:117:ALA:HA	2.00	0.44
3:C:99:GLU:OE1	3:C:99:GLU:N	2.50	0.43
3:C:105:LYS:HB2	3:C:116:GLN:HA	1.99	0.43
4:D:209:LYS:HE2	4:D:209:LYS:HB2	1.81	0.43
6:F:392:LYS:NZ	6:F:398:GLU:OE1	2.48	0.43
1:A:10:LEU:HD12	1:A:11:PRO:HD2	1.99	0.43
2:B:70:TRP:CD1	2:B:276:SER:HB2	2.53	0.43
4:D:54:LEU:HD23	4:D:54:LEU:HA	1.83	0.43
6:F:124:PRO:O	6:F:127:ILE:HG23	2.18	0.43
1:A:363:MET:SD	2:B:47:VAL:HG11	2.57	0.43
10:B:504:LMT:H42	10:B:504:LMT:H11	1.82	0.43
2:B:395:LEU:HD21	5:E:155:VAL:HG12	2.00	0.43
2:B:367:MET:HE3	2:B:371:ILE:HD11	2.00	0.43
5:E:25:MET:HA	5:E:25:MET:HE2	2.00	0.43
10:B:504:LMT:H121	10:B:504:LMT:H92	1.87	0.43
3:C:46:GLN:CD	3:C:90:ASN:HD21	2.26	0.43
3:C:98:LYS:HG2	3:C:99:GLU:OE1	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:151:MET:O	5:E:155:VAL:HG23	2.19	0.43
6:F:214:MET:HA	6:F:226:LEU:HD23	2.00	0.43
2:B:301:VAL:HG11	2:B:340:MET:HB2	2.01	0.42
10:D:301:LMT:H5'	10:D:301:LMT:H1B	1.77	0.42
6:F:36:THR:HG22	6:F:48:VAL:HG22	2.02	0.42
6:F:80:ARG:HB2	6:F:107:GLU:HG2	2.01	0.42
6:F:201:GLU:OE1	6:F:201:GLU:N	2.52	0.42
6:F:306:SER:HG	6:F:308:TRP:CD1	2.38	0.42
3:C:189:TRP:CE3	3:C:192:LYS:HB3	2.54	0.42
4:D:107:LEU:HD12	4:D:107:LEU:H	1.83	0.42
1:A:154:MET:SD	1:A:218:GLY:HA2	2.60	0.42
6:F:144:ALA:HB3	6:F:147:ILE:HB	2.01	0.42
6:F:140:ASN:ND2	6:F:148:LYS:HD2	2.34	0.42
4:D:63:ARG:HG2	4:D:63:ARG:HH11	1.85	0.41
6:F:127:ILE:C	6:F:127:ILE:HD12	2.45	0.41
3:C:125:TYR:HB3	3:C:137:ILE:CG2	2.49	0.41
3:C:152:PHE:CB	3:C:163:SER:HB2	2.50	0.41
7:C:301:FMN:HM73	7:C:301:FMN:HM81	1.79	0.41
2:B:260:TRP:CH2	9:B:503:PEE:H8	2.55	0.41
8:B:502:RBF:HC3'	8:B:502:RBF:N1	2.35	0.41
5:E:46:ILE:HD11	5:E:151:MET:HE2	2.02	0.41
1:A:39:MET:HE1	1:A:58:PHE:CD1	2.56	0.41
3:C:225:THR:OG1	7:C:301:FMN:H5'2	2.19	0.41
6:F:377:MET:HG2	6:F:384:ASN:OD1	2.20	0.41
1:A:381:GLU:N	1:A:381:GLU:OE1	2.53	0.41
2:B:83:LEU:HD11	2:B:102:TYR:CZ	2.56	0.41
7:B:501:FMN:H9	7:B:501:FMN:H1'1	1.81	0.41
3:C:142:GLY:N	3:C:147:SER:O	2.38	0.41
3:C:221:LEU:HD23	3:C:222:SER:N	2.29	0.41
3:C:76:VAL:HG13	3:C:124:VAL:HG11	2.01	0.41
5:E:65:LYS:HB3	5:E:68:ALA:HB2	2.03	0.41
5:E:122:ILE:O	5:E:126:VAL:HG23	2.21	0.40
6:F:230:ILE:HB	6:F:250:TRP:CE2	2.57	0.40
2:B:200:ASN:OD1	8:B:502:RBF:O5'	2.27	0.40
6:F:332:ASN:OD1	6:F:332:ASN:N	2.53	0.40
2:B:100:TRP:HA	2:B:103:TRP:CD1	2.56	0.40
3:C:231:VAL:O	3:C:234:THR:HG22	2.22	0.40
6:F:137:VAL:HG12	6:F:254:ALA:HA	2.04	0.40
1:A:345:PHE:HE2	1:A:423:PHE:HE2	1.70	0.40
2:B:52:ASP:OD1	2:B:52:ASP:N	2.52	0.40
2:B:145:VAL:HG21	11:B:506:IQT:O4	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:395:LEU:HA	2:B:395:LEU:HD23	1.87	0.40
5:E:32:SER:O	5:E:161:ARG:HD3	2.22	0.40
6:F:248:TYR:CD2	6:F:248:TYR:C	2.98	0.40
6:F:299:LEU:HD23	6:F:299:LEU:HA	1.95	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	444/446 (100%)	430 (97%)	14 (3%)	0	100	100
2	B	410/415 (99%)	406 (99%)	4 (1%)	0	100	100
3	C	250/257 (97%)	204 (82%)	43 (17%)	3 (1%)	10	27
4	D	204/210 (97%)	200 (98%)	4 (2%)	0	100	100
5	E	196/198 (99%)	186 (95%)	10 (5%)	0	100	100
6	F	406/414 (98%)	391 (96%)	15 (4%)	0	100	100
All	All	1910/1940 (98%)	1817 (95%)	90 (5%)	3 (0%)	44	65

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	104	ILE
3	C	153	VAL
3	C	241	ASP

5.3.2 Protein sidechains [i](#)

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

entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	376/376 (100%)	369 (98%)	7 (2%)	50	71
2	B	317/320 (99%)	315 (99%)	2 (1%)	78	89
3	C	201/205 (98%)	196 (98%)	5 (2%)	42	66
4	D	173/176 (98%)	170 (98%)	3 (2%)	53	73
5	E	165/165 (100%)	159 (96%)	6 (4%)	31	56
6	F	335/343 (98%)	327 (98%)	8 (2%)	43	67
All	All	1567/1585 (99%)	1536 (98%)	31 (2%)	48	70

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	MET
1	A	97	ASP
1	A	162	GLU
1	A	277	THR
1	A	284	GLU
1	A	350	SER
1	A	442	ILE
2	B	49	ASP
2	B	155	VAL
3	C	12	LEU
3	C	127	VAL
3	C	137	ILE
3	C	166	THR
3	C	221	LEU
4	D	71	ARG
4	D	164	GLU
4	D	165	VAL
5	E	70	VAL
5	E	97	ILE
5	E	105	LEU
5	E	111	ILE
5	E	118	VAL
5	E	133	ASP
6	F	56	LEU
6	F	93	THR
6	F	122	GLU

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Mol	Chain	Res	Type
6	F	177	VAL
6	F	258	CYS
6	F	292	ILE
6	F	322	ASP
6	F	332	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	223	HIS
2	B	267	ASN
3	C	188	GLN
4	D	24	GLN
5	E	56	ASN
6	F	140	ASN
6	F	176	HIS
6	F	196	ASN
6	F	362	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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
10	LMT	D	301	-	36,36,36	0.53	0	47,47,47	0.76	0
11	IQT	B	506	-	31,32,32	2.00	5 (16%)	33,42,42	0.74	0
10	LMT	B	504	-	36,36,36	0.56	0	47,47,47	0.66	0
7	FMN	B	501	-	29,32,33	1.13	2 (6%)	41,47,50	1.30	8 (19%)
14	FAD	F	502	-	58,58,58	1.26	3 (5%)	85,89,89	0.73	1 (1%)
13	FES	F	501	6	0,4,4	-	-	-	-	-
7	FMN	C	301	-	29,32,33	1.11	2 (6%)	41,47,50	1.32	7 (17%)
9	PEE	B	503	-	50,50,50	0.89	2 (4%)	53,55,55	0.72	1 (1%)
13	FES	D	302	4,5	0,4,4	-	-	-	-	-
9	PEE	B	505	-	50,50,50	0.90	2 (4%)	53,55,55	0.62	0
8	RBF	B	502	-	29,29,29	1.01	1 (3%)	42,43,43	0.84	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
10	LMT	D	301	-	-	9/21/61/61	0/2/2/2
11	IQT	B	506	-	-	10/27/46/46	0/2/2/2
10	LMT	B	504	-	-	10/21/61/61	0/2/2/2
7	FMN	B	501	-	-	7/15/17/18	0/3/3/3
14	FAD	F	502	-	-	10/34/50/50	0/6/6/6
13	FES	F	501	6	-	-	0/1/1/1
7	FMN	C	301	-	-	11/15/17/18	0/3/3/3
9	PEE	B	503	-	-	17/54/54/54	-
13	FES	D	302	4,5	-	-	0/1/1/1
9	PEE	B	505	-	-	19/54/54/54	-
8	RBF	B	502	-	-	6/14/14/14	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	B	506	IQT	C6-C7	6.95	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	F	502	FAD	P-O3P	6.17	1.66	1.59
11	B	506	IQT	C13-C12	-4.98	1.28	1.44
11	B	506	IQT	C7-C1	-4.61	1.38	1.50
9	B	505	PEE	C39-C38	3.73	1.52	1.31
9	B	503	PEE	C18-C19	3.72	1.52	1.31
9	B	503	PEE	C39-C38	3.72	1.52	1.31
9	B	505	PEE	C18-C19	3.69	1.52	1.31
14	F	502	FAD	PA-O3P	3.63	1.63	1.59
7	B	501	FMN	C4A-N5	3.42	1.38	1.30
7	C	301	FMN	C4A-N5	3.39	1.38	1.30
11	B	506	IQT	C7-N1	-3.31	1.32	1.41
7	C	301	FMN	C10-N1	2.69	1.38	1.33
7	B	501	FMN	C10-N1	2.52	1.38	1.33
8	B	502	RBF	C5A-N5	-2.38	1.35	1.39
11	B	506	IQT	C8-N1	-2.30	1.32	1.37
14	F	502	FAD	C5X-N5	-2.28	1.35	1.39

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	301	FMN	C4-N3-C2	-3.32	119.74	125.64
7	B	501	FMN	C4-N3-C2	-3.19	119.98	125.64
7	C	301	FMN	O4-C4-C4A	-2.75	119.26	126.53
7	C	301	FMN	C4A-C4-N3	2.73	120.20	113.25
7	B	501	FMN	C4A-C4-N3	2.69	120.10	113.25
7	B	501	FMN	C4A-C10-N10	2.56	120.14	116.48
7	B	501	FMN	O4-C4-C4A	-2.44	120.10	126.53
7	B	501	FMN	C10-C4A-N5	-2.35	120.00	124.81
7	C	301	FMN	C4A-C10-N10	2.35	119.85	116.48
7	B	501	FMN	C5A-C9A-N10	2.20	119.96	117.97
7	B	501	FMN	C9A-C5A-N5	-2.19	120.13	122.45
9	B	503	PEE	C2-O2-C10	2.19	123.04	117.80
14	F	502	FAD	O2P-P-O1P	2.19	122.64	112.44
7	C	301	FMN	C10-C4A-N5	-2.15	120.42	124.81
7	C	301	FMN	C4A-C10-N1	-2.15	119.33	124.59
7	C	301	FMN	C4'-C3'-C2'	-2.08	110.11	113.57
7	B	501	FMN	C4A-C10-N1	-2.02	119.63	124.59

There are no chirality outliers.

All (99) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	501	FMN	C3'-C4'-C5'-O5'
7	B	501	FMN	O4'-C4'-C5'-O5'
7	C	301	FMN	C1'-C2'-C3'-O3'
7	C	301	FMN	C1'-C2'-C3'-C4'
7	C	301	FMN	O2'-C2'-C3'-O3'
7	C	301	FMN	O2'-C2'-C3'-C4'
7	C	301	FMN	O4'-C4'-C5'-O5'
8	B	502	RBF	C2'-C3'-C4'-O4'
8	B	502	RBF	C2'-C3'-C4'-C5'
8	B	502	RBF	O3'-C3'-C4'-C5'
9	B	503	PEE	O4P-C4-C5-N
9	B	505	PEE	C1-O3P-P-O2P
9	B	505	PEE	C1-O3P-P-O1P
9	B	505	PEE	C1-O3P-P-O4P
10	D	301	LMT	O5'-C1'-O1'-C1
11	B	506	IQT	C11-C12-C13-C14
11	B	506	IQT	C10-C11-C12-C13
11	B	506	IQT	C12-C13-C14-C15
14	F	502	FAD	C5B-O5B-PA-O1A
14	F	502	FAD	C5B-O5B-PA-O2A
14	F	502	FAD	C5B-O5B-PA-O3P
14	F	502	FAD	C3'-C4'-C5'-O5'
9	B	503	PEE	O5-C30-O3-C3
9	B	505	PEE	O5-C30-O3-C3
9	B	503	PEE	C31-C30-O3-C3
10	D	301	LMT	C5'-C4'-O1B-C1B
9	B	505	PEE	C31-C30-O3-C3
8	B	502	RBF	O3'-C3'-C4'-O4'
9	B	503	PEE	C37-C38-C39-C40
10	B	504	LMT	O5'-C5'-C6'-O6'
9	B	505	PEE	C37-C38-C39-C40
7	B	501	FMN	C4'-C5'-O5'-P
10	B	504	LMT	C4'-C5'-C6'-O6'
10	D	301	LMT	C2'-C1'-O1'-C1
9	B	505	PEE	C17-C18-C19-C20
9	B	503	PEE	C11-C12-C13-C14
10	B	504	LMT	C5-C6-C7-C8
10	D	301	LMT	C2-C3-C4-C5
11	B	506	IQT	C21-C22-C23-C24
10	B	504	LMT	C2-C1-O1'-C1'
9	B	503	PEE	C10-C11-C12-C13
14	F	502	FAD	O4'-C4'-C5'-O5'
14	F	502	FAD	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
9	B	503	PEE	C12-C13-C14-C15
7	B	501	FMN	O3'-C3'-C4'-C5'
7	B	501	FMN	C2'-C3'-C4'-O4'
10	B	504	LMT	O1'-C1-C2-C3
9	B	505	PEE	C40-C41-C42-C43
10	D	301	LMT	O1'-C1-C2-C3
9	B	505	PEE	C11-C10-O2-C2
14	F	502	FAD	C3B-C4B-C5B-O5B
9	B	505	PEE	O2-C2-C3-O3
7	C	301	FMN	C3'-C4'-C5'-O5'
10	D	301	LMT	C4-C5-C6-C7
9	B	505	PEE	C1-C2-C3-O3
10	B	504	LMT	C1-C2-C3-C4
11	B	506	IQT	C13-C14-C15-C16
10	B	504	LMT	C2-C3-C4-C5
11	B	506	IQT	C14-C15-C16-O4
10	D	301	LMT	C3'-C4'-O1B-C1B
10	B	504	LMT	C7-C8-C9-C10
7	C	301	FMN	O3'-C3'-C4'-C5'
7	C	301	FMN	C4'-C5'-O5'-P
10	D	301	LMT	C2-C1-O1'-C1'
7	B	501	FMN	C2'-C3'-C4'-C5'
7	C	301	FMN	C2'-C3'-C4'-C5'
9	B	505	PEE	O4-C10-O2-C2
9	B	505	PEE	C13-C14-C15-C16
7	B	501	FMN	O3'-C3'-C4'-O4'
7	C	301	FMN	O3'-C3'-C4'-O4'
11	B	506	IQT	C1-C7-N1-C8
9	B	503	PEE	C18-C19-C20-C21
9	B	503	PEE	O3P-C1-C2-O2
9	B	503	PEE	C1-C2-C3-O3
7	C	301	FMN	C2'-C3'-C4'-O4'
9	B	503	PEE	O2-C2-C3-O3
14	F	502	FAD	C2B-C1B-N9A-C8A
9	B	503	PEE	O3P-C1-C2-C3
8	B	502	RBF	N10-C1'-C2'-C3'
11	B	506	IQT	C6-C2-C4-C5
9	B	505	PEE	C36-C37-C38-C39
9	B	505	PEE	C4-O4P-P-O1P
9	B	505	PEE	C38-C39-C40-C41
9	B	503	PEE	C42-C43-C44-C45
11	B	506	IQT	C20-C21-C22-C23

Continued on next page...

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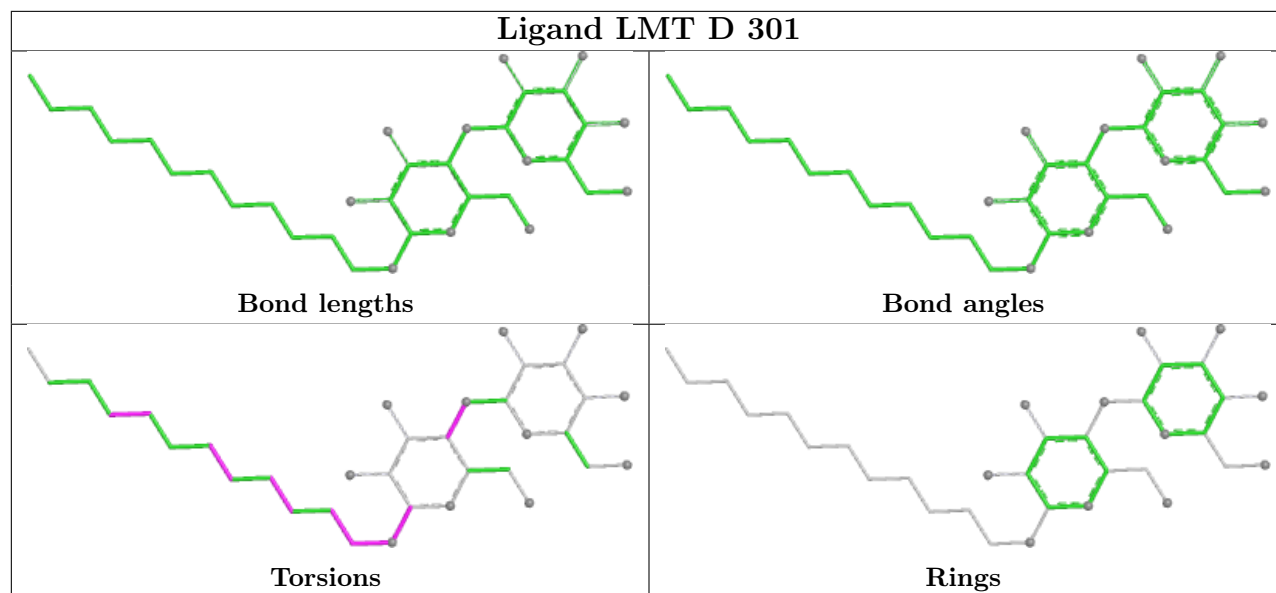
Mol	Chain	Res	Type	Atoms
9	B	503	PEE	C15-C16-C17-C18
9	B	503	PEE	C2-C1-O3P-P
10	B	504	LMT	C11-C10-C9-C8
14	F	502	FAD	C2B-C1B-N9A-C4A
9	B	505	PEE	C18-C19-C20-C21
9	B	505	PEE	C2-C1-O3P-P
14	F	502	FAD	O4B-C1B-N9A-C8A
9	B	503	PEE	C38-C39-C40-C41
8	B	502	RBF	O2'-C2'-C3'-C4'
9	B	503	PEE	C36-C37-C38-C39
9	B	505	PEE	C16-C17-C18-C19
10	D	301	LMT	C7-C8-C9-C10
10	B	504	LMT	C9-C10-C11-C12
11	B	506	IQT	C19-C20-C21-C22

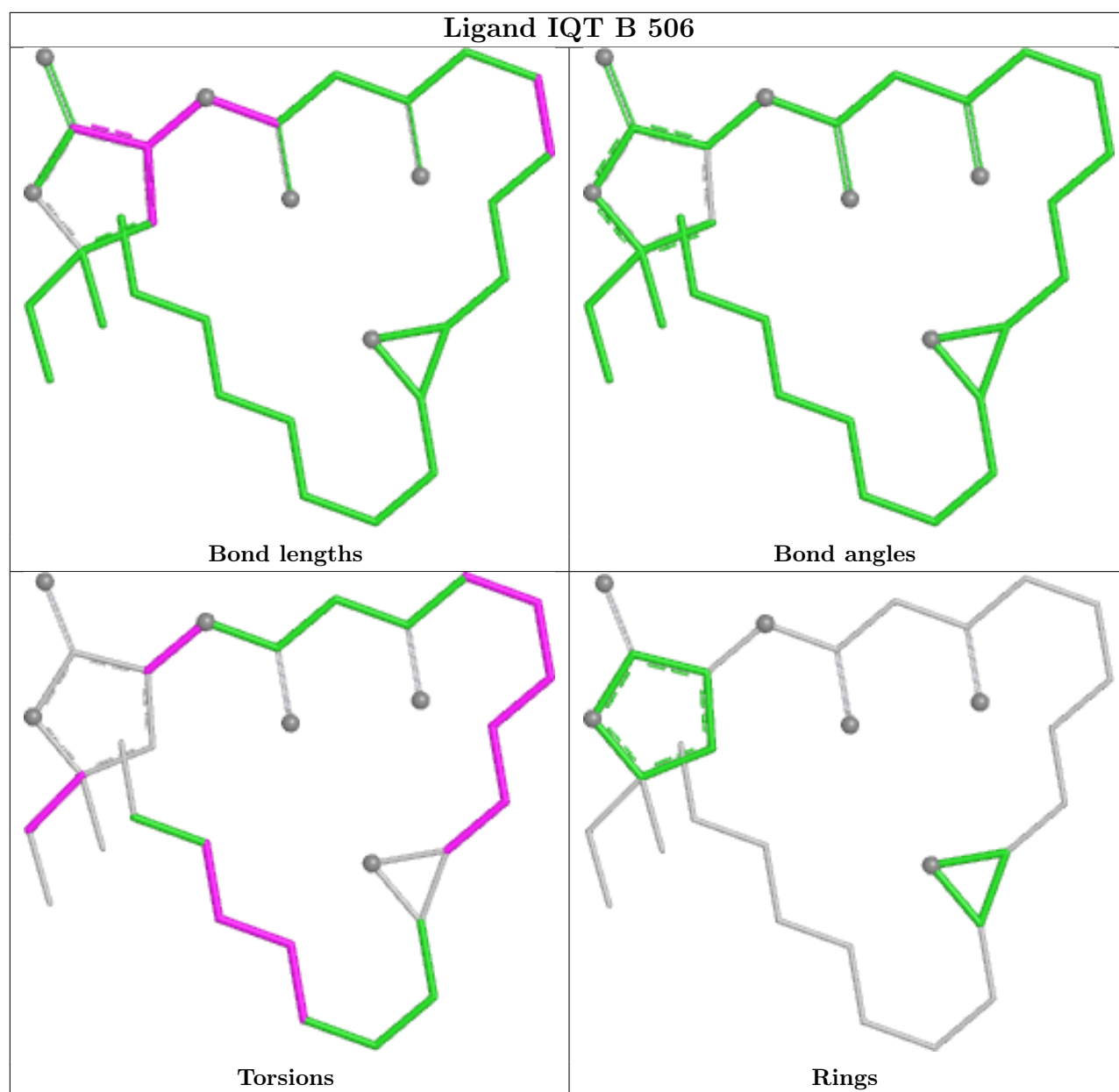
There are no ring outliers.

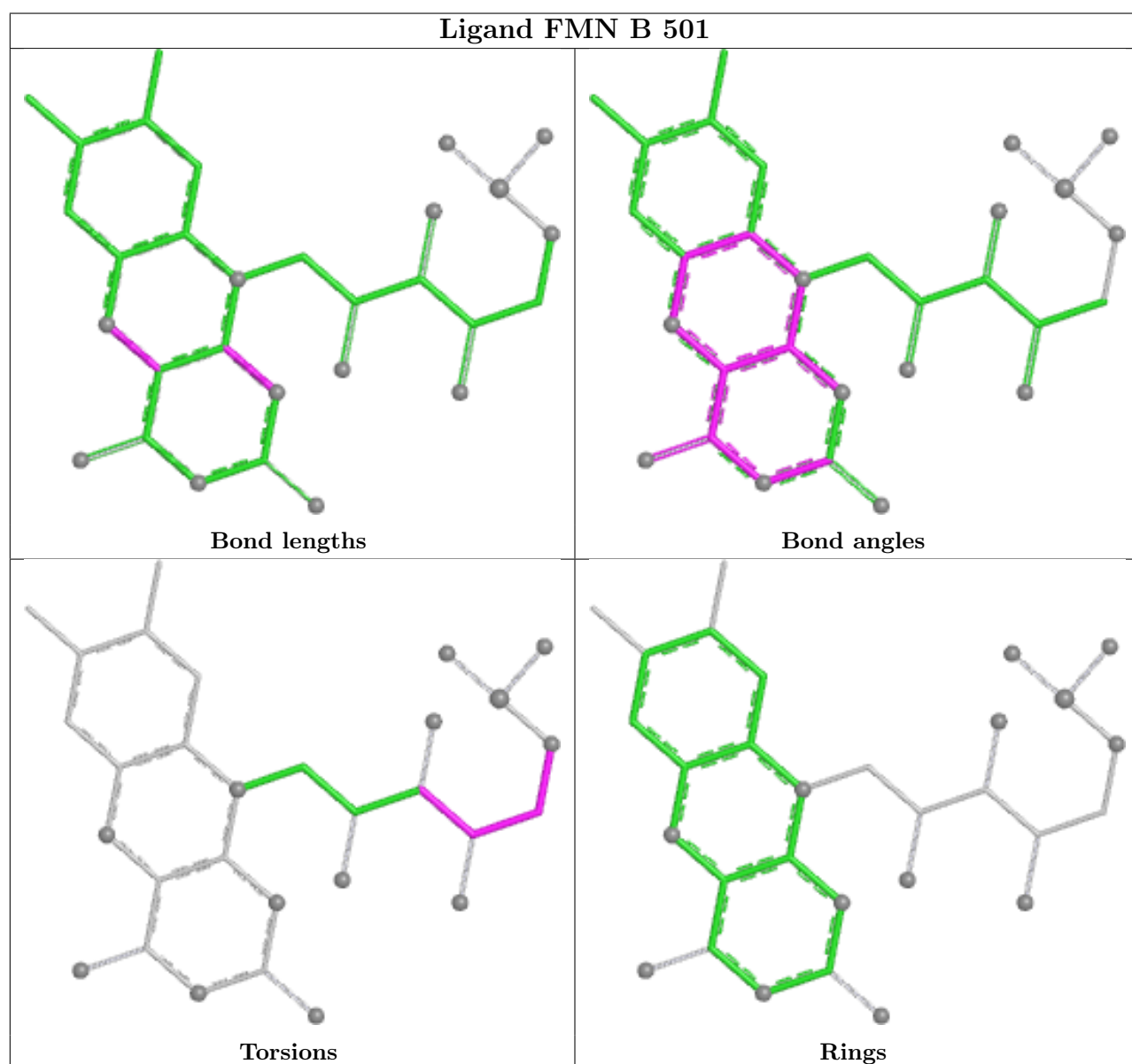
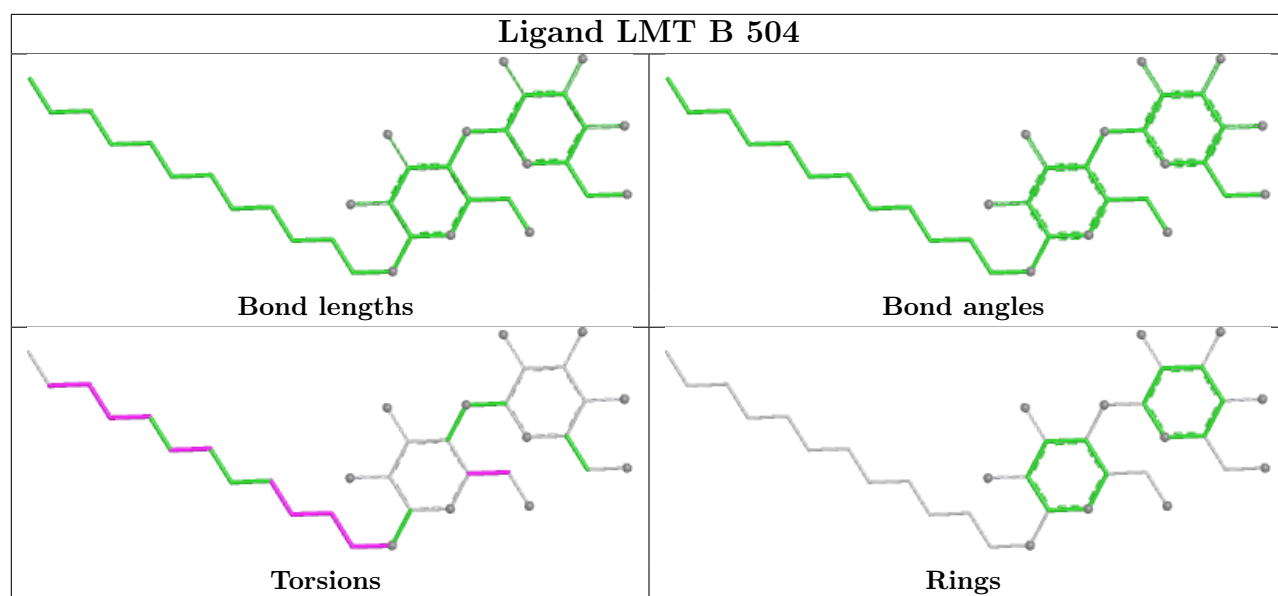
9 monomers are involved in 16 short contacts:

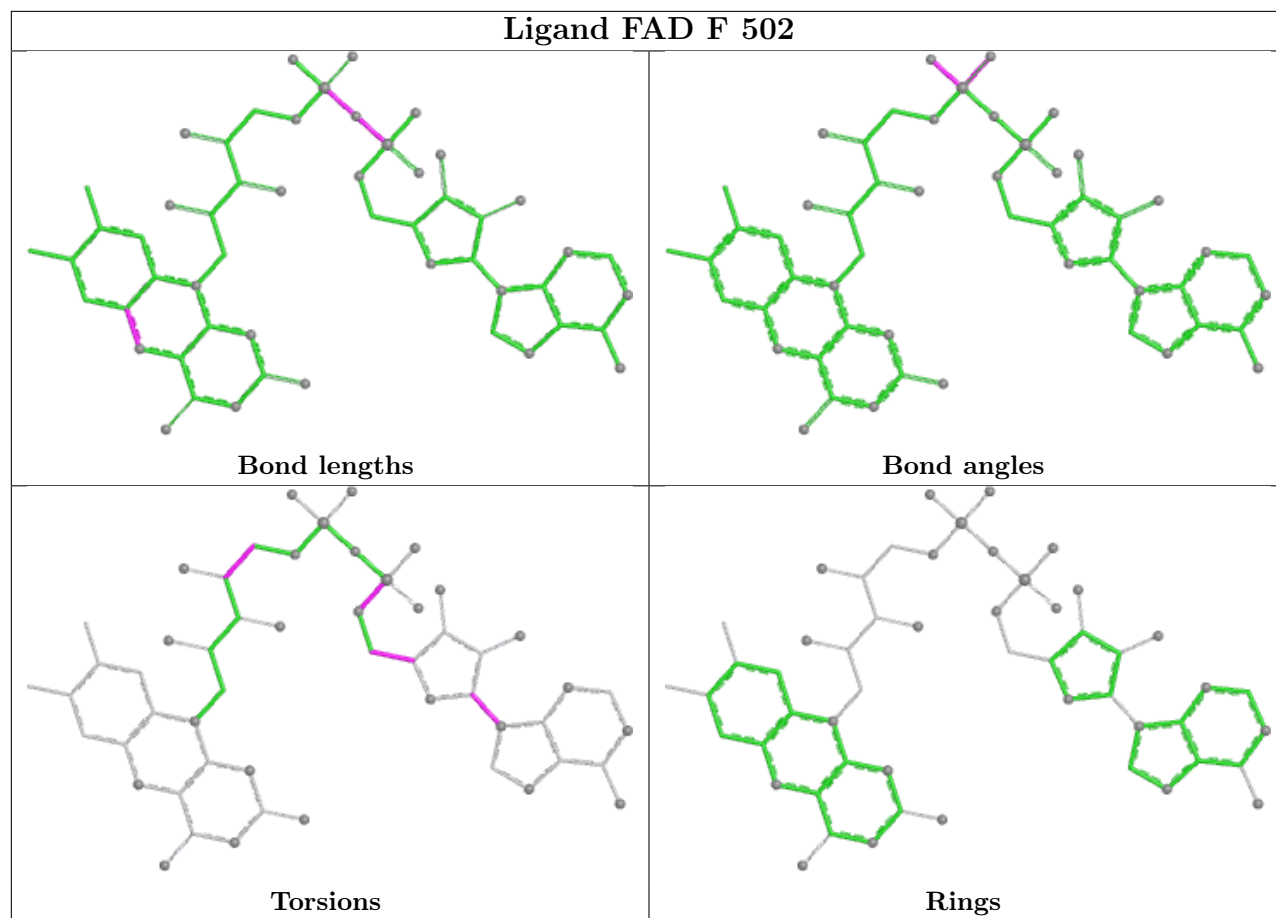
Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	D	301	LMT	1	0
11	B	506	IQT	2	0
10	B	504	LMT	2	0
7	B	501	FMN	1	0
14	F	502	FAD	1	0
7	C	301	FMN	3	0
9	B	503	PEE	2	0
9	B	505	PEE	1	0
8	B	502	RBF	3	0

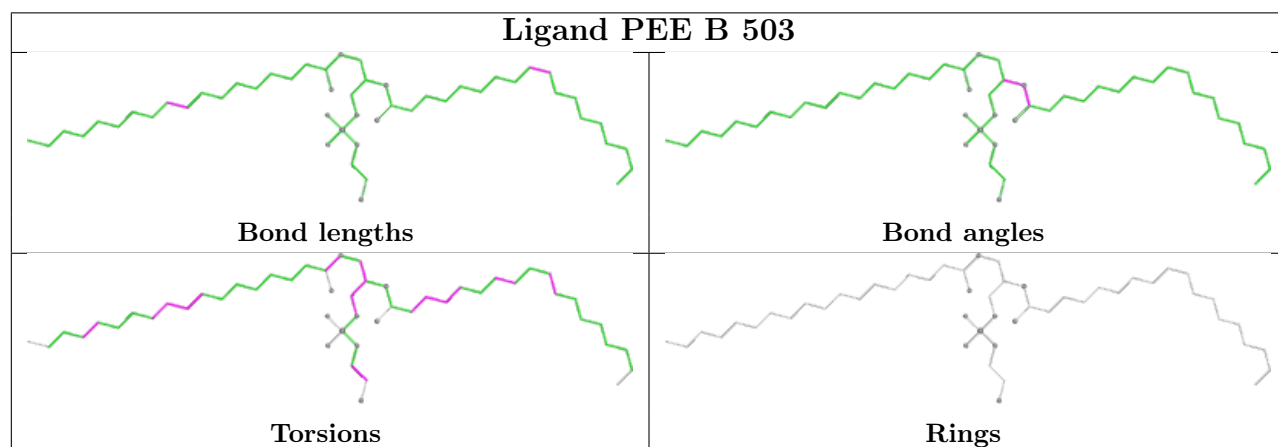
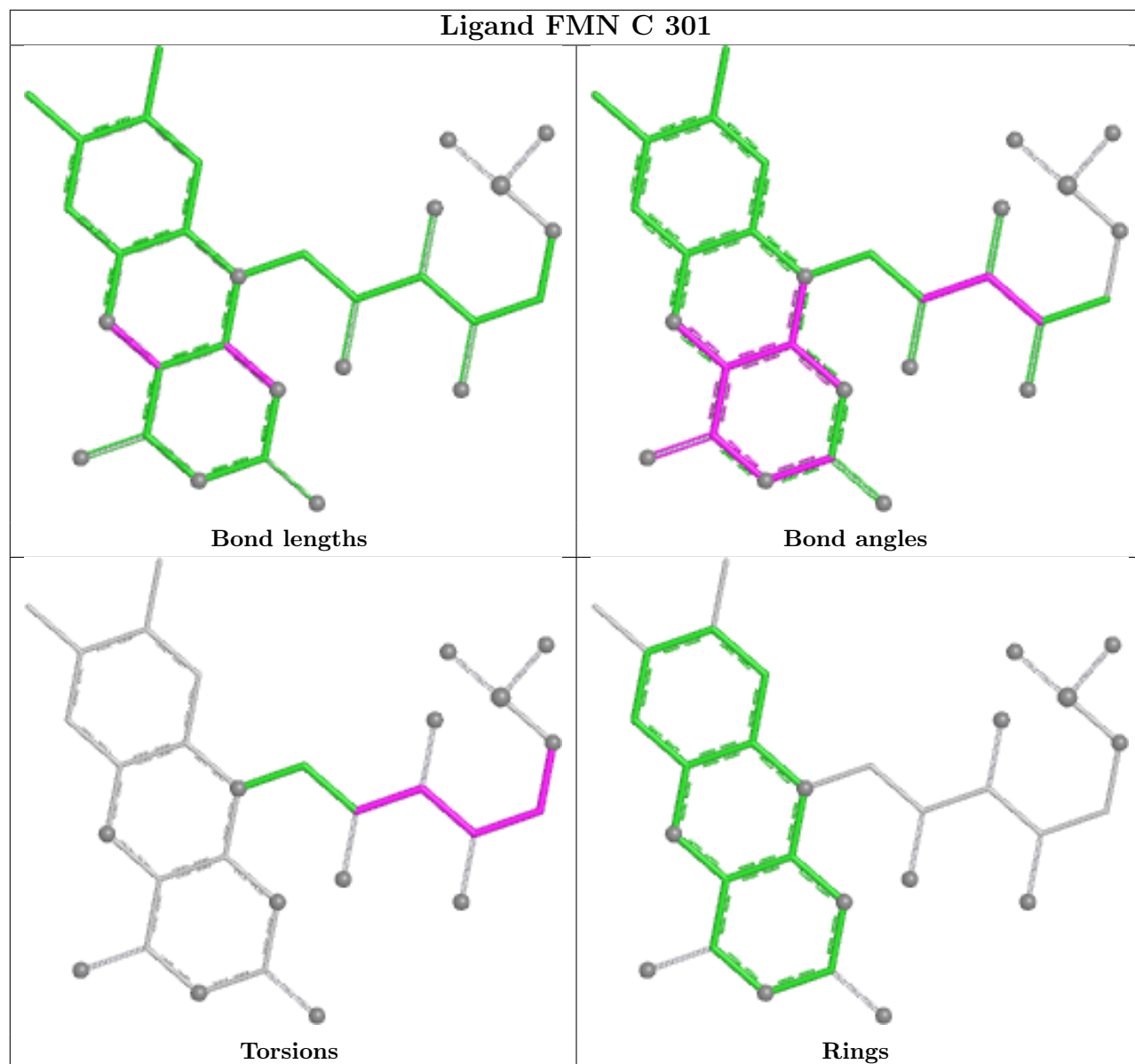
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

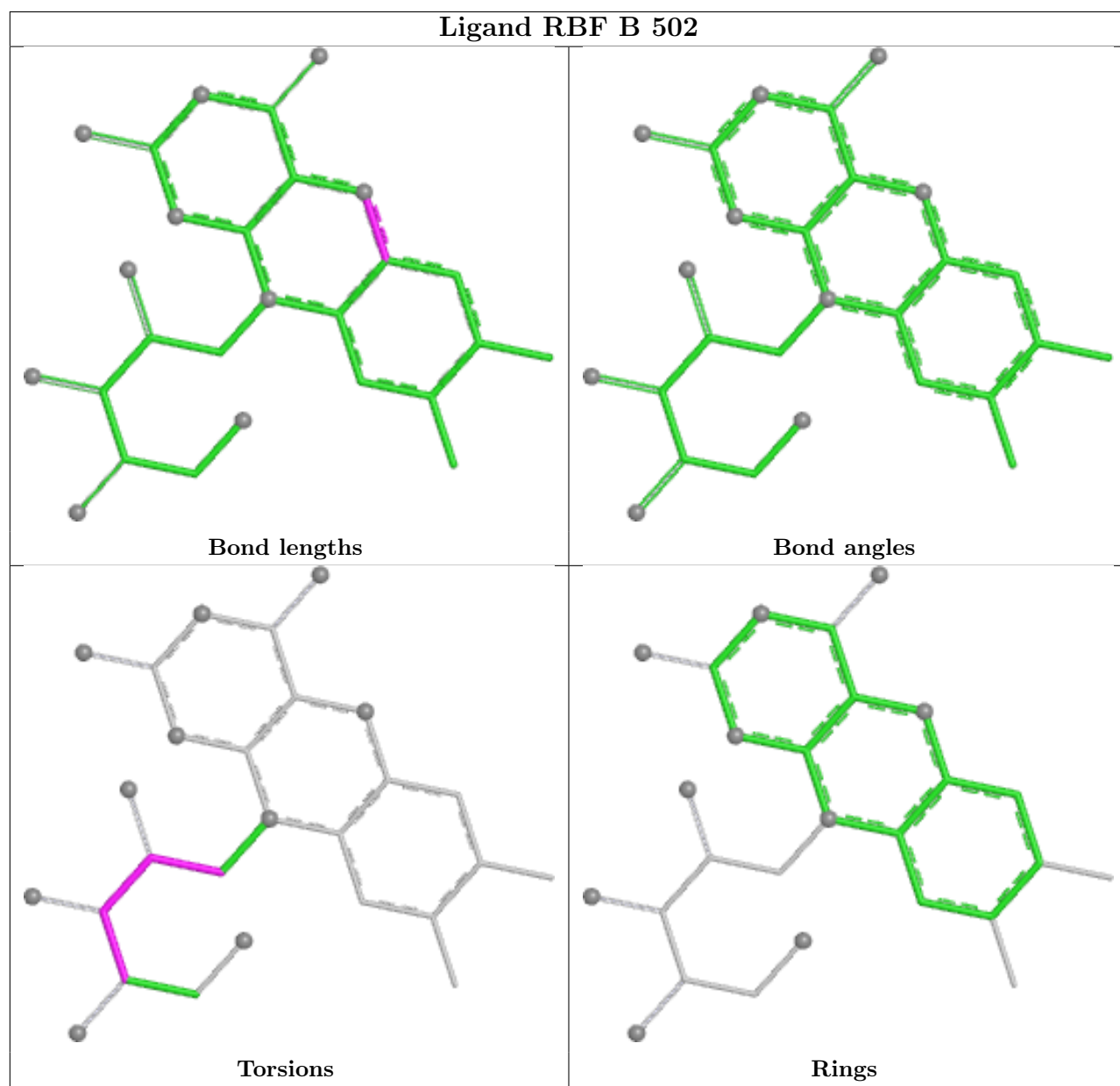
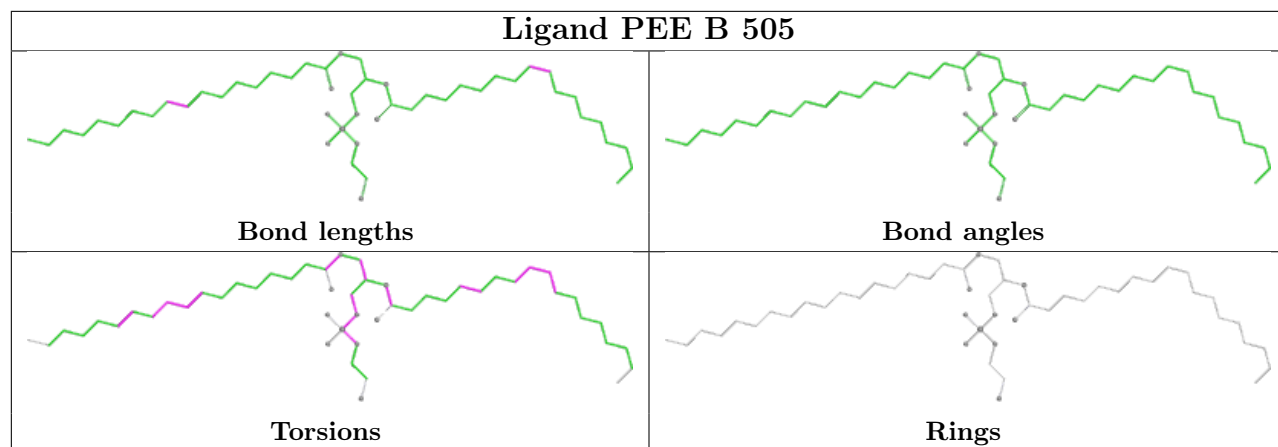












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

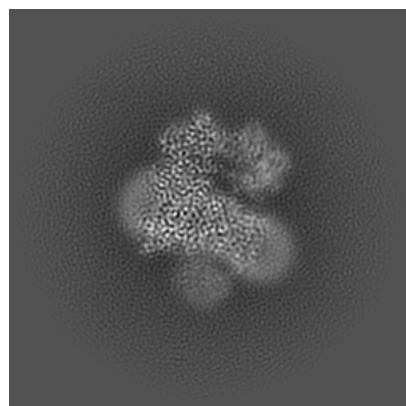
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-64068. These allow visual inspection of the internal detail of the map and identification of artifacts.

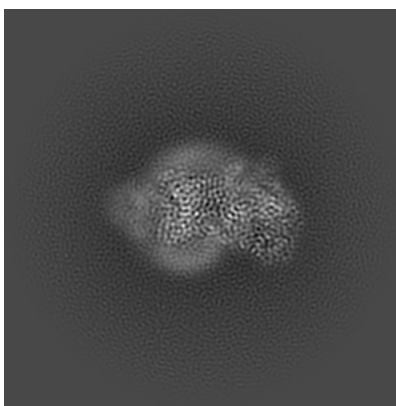
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

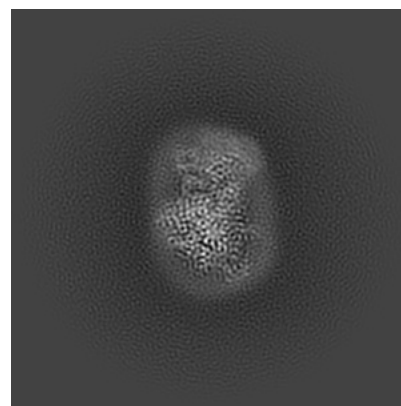
6.1.1 Primary map



X

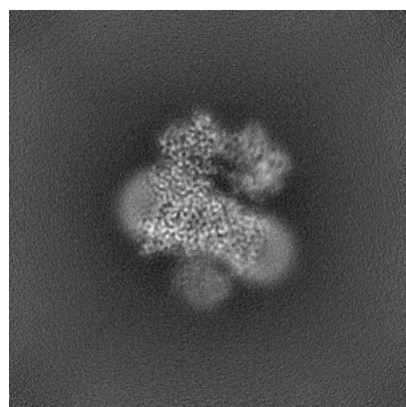


Y

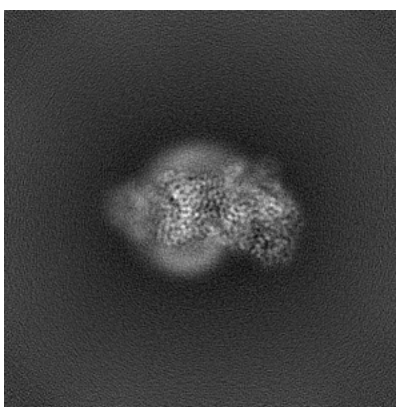


Z

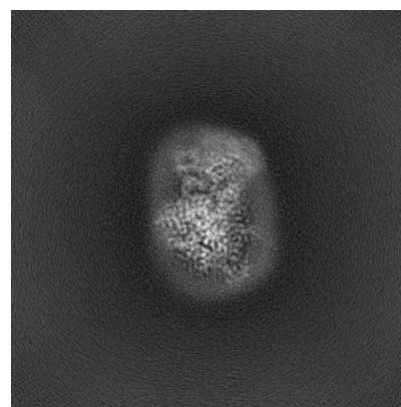
6.1.2 Raw map



X



Y

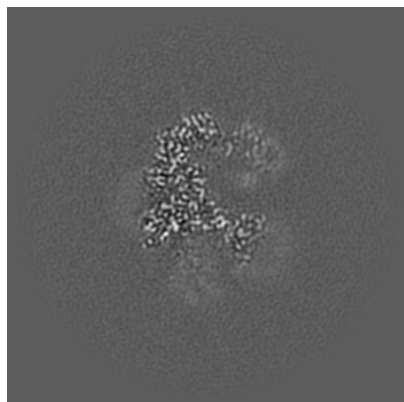


Z

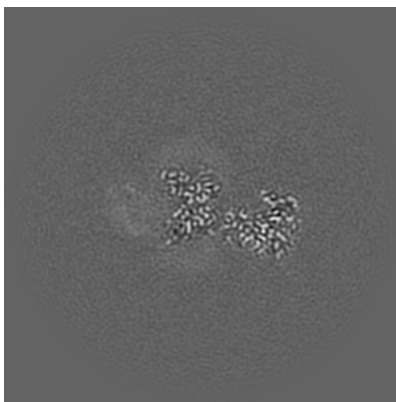
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

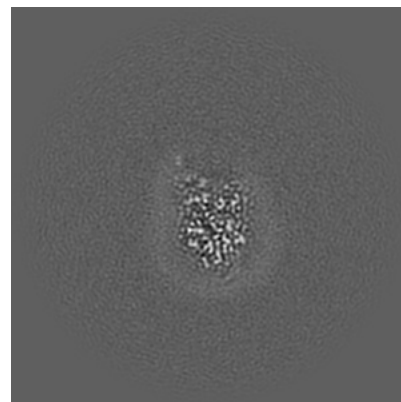
6.2.1 Primary map



X Index: 160

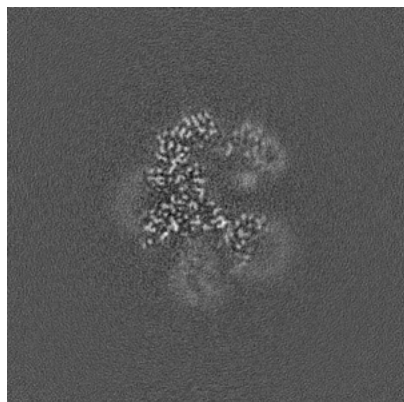


Y Index: 160

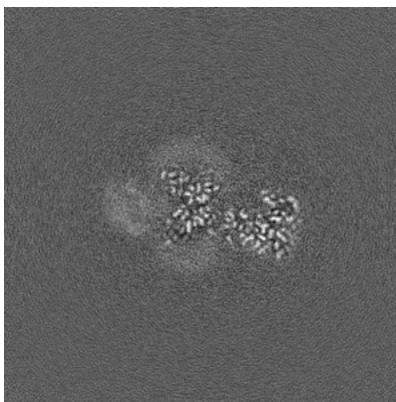


Z Index: 160

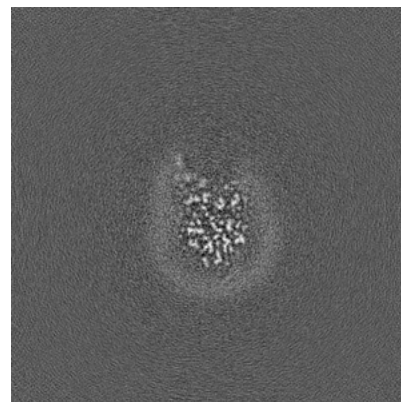
6.2.2 Raw map



X Index: 160



Y Index: 160

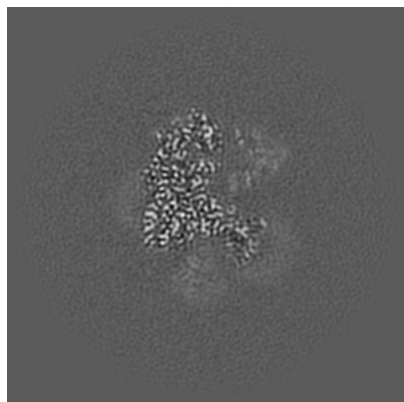


Z Index: 160

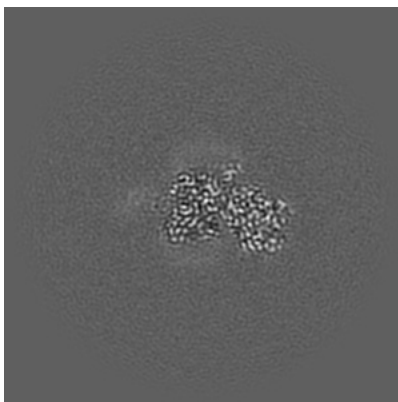
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

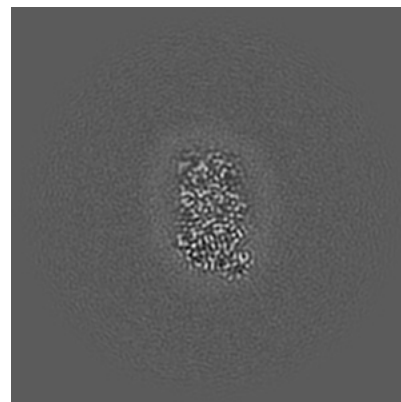
6.3.1 Primary map



X Index: 155

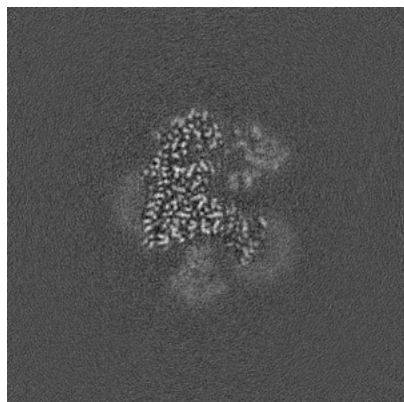


Y Index: 136

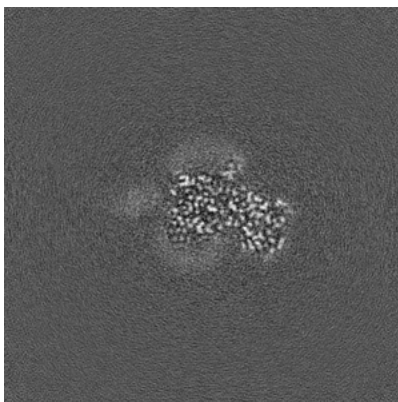


Z Index: 142

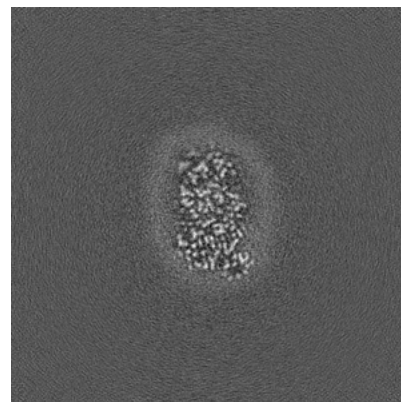
6.3.2 Raw map



X Index: 156



Y Index: 137

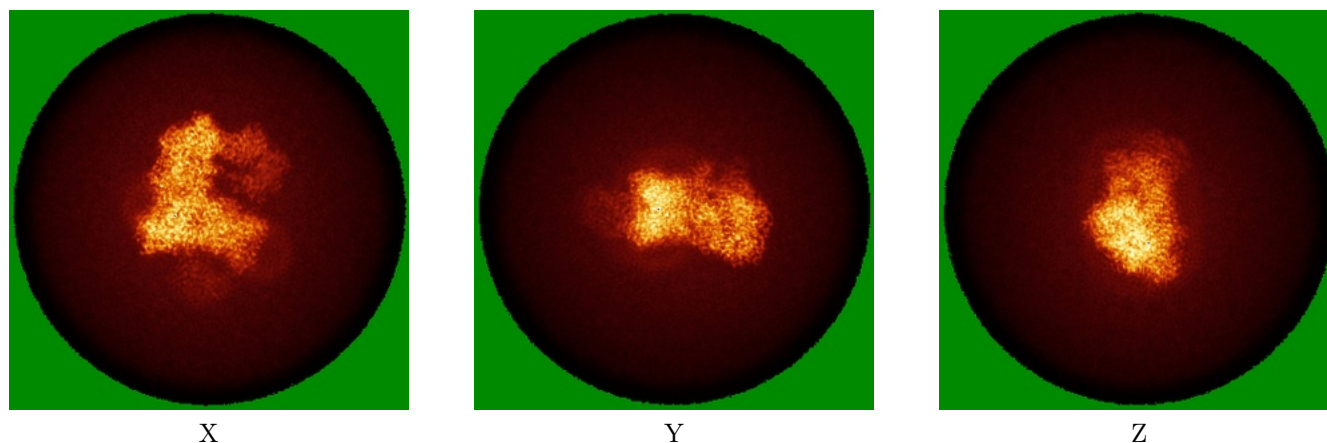


Z Index: 142

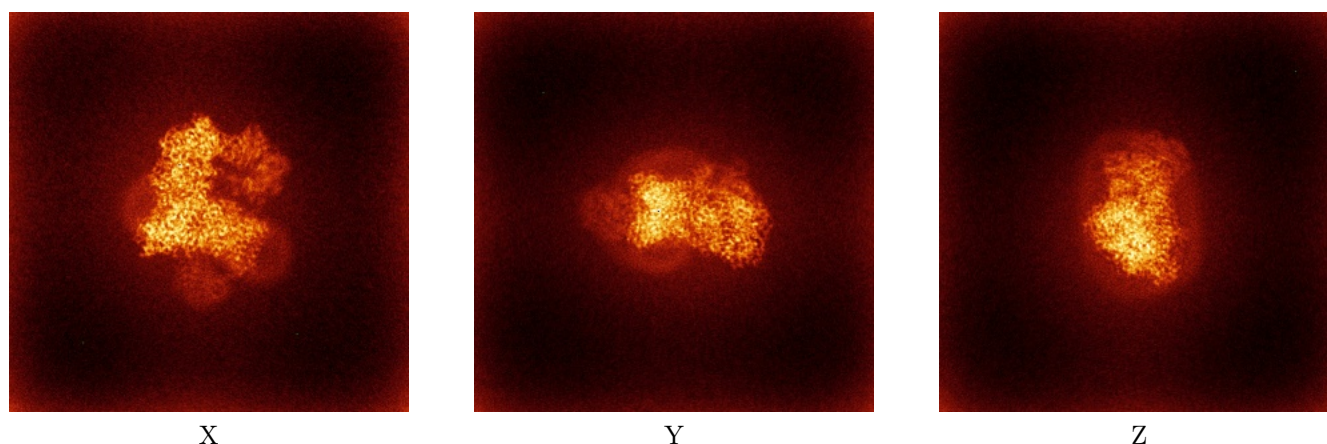
The images above show the largest variance slices of the map in three orthogonal directions.

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.

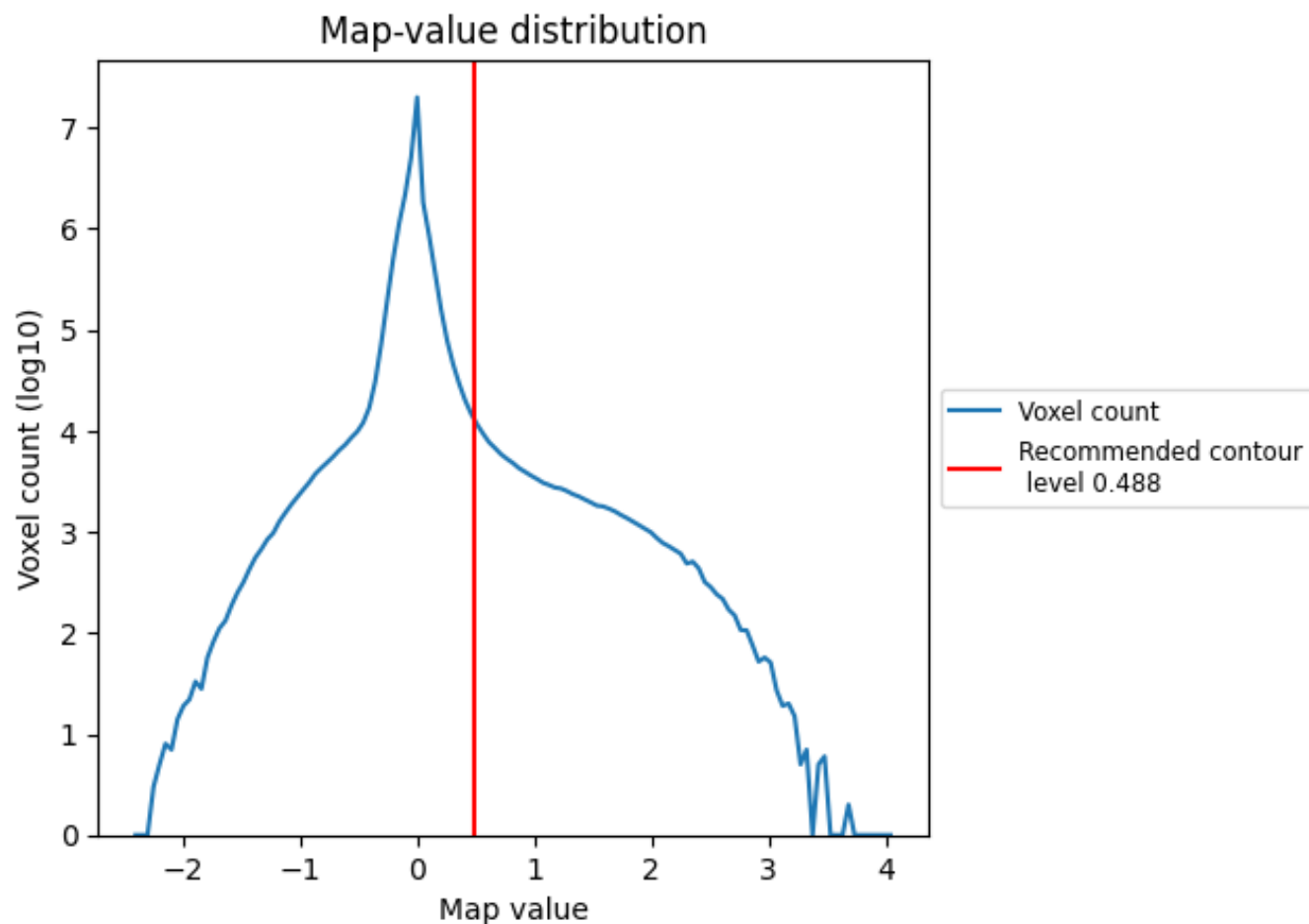
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

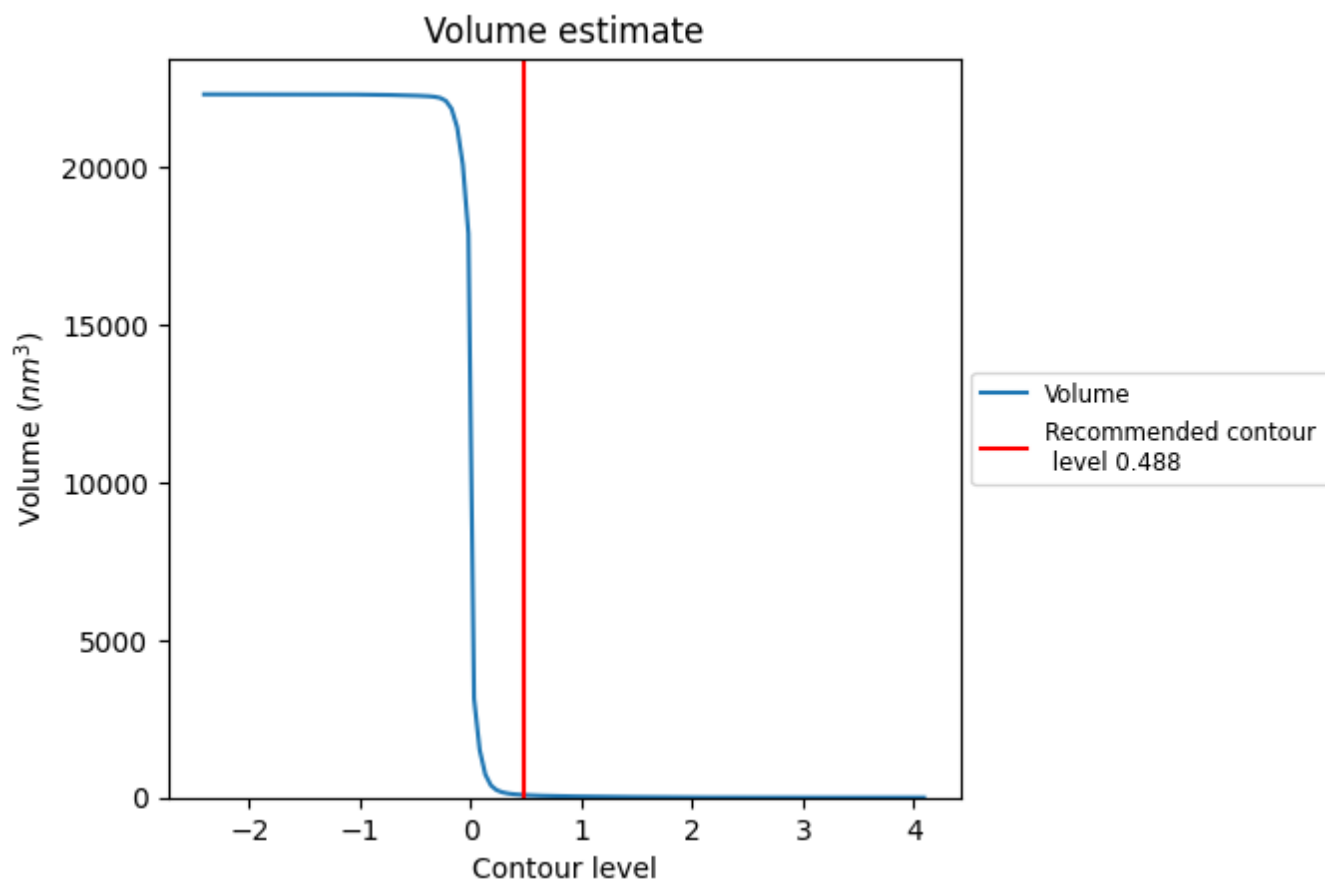
This section contains the results of statistical analysis of the map.

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.

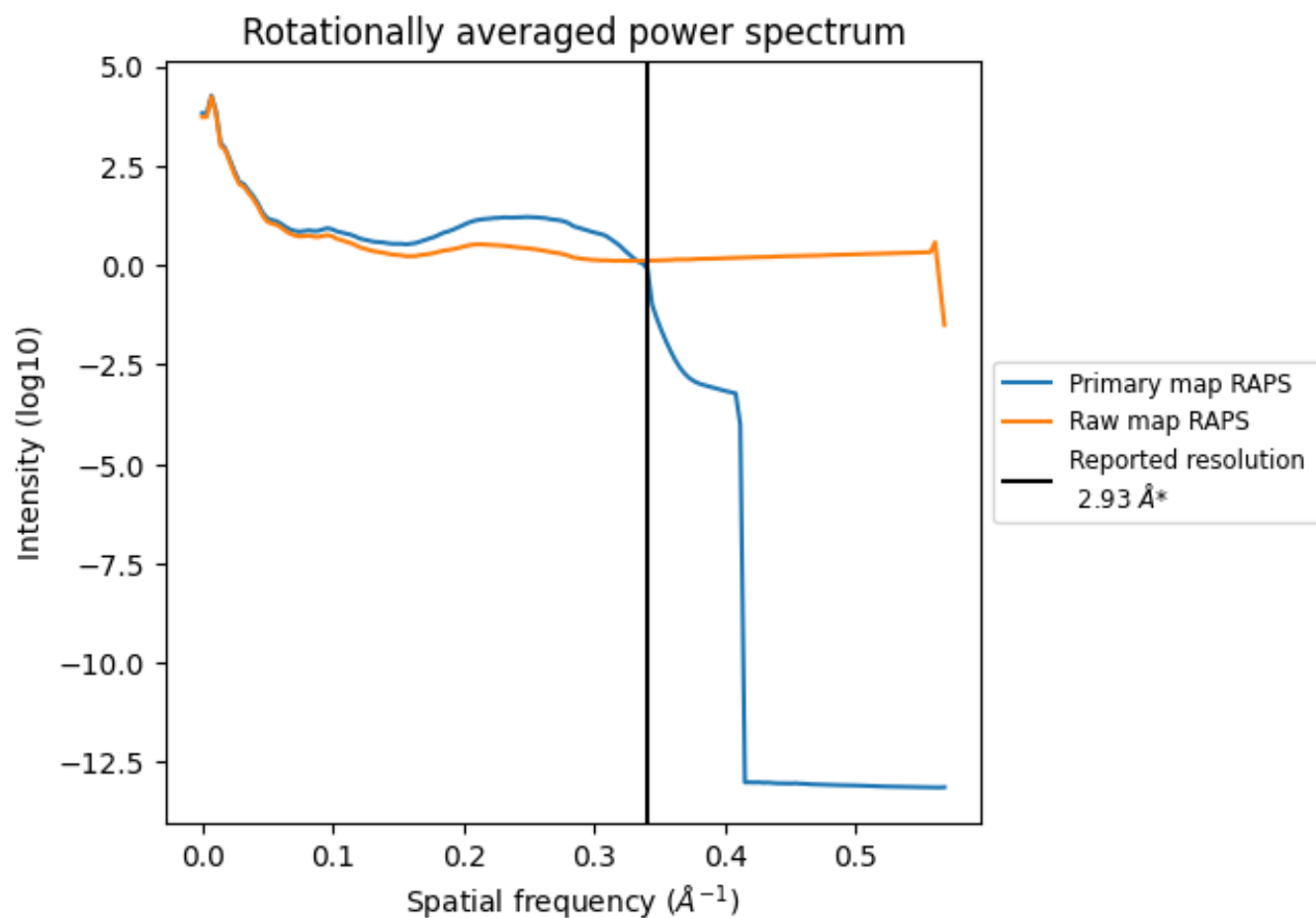
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 79 nm^3 ; this corresponds to an approximate mass of 72 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 ⓘ

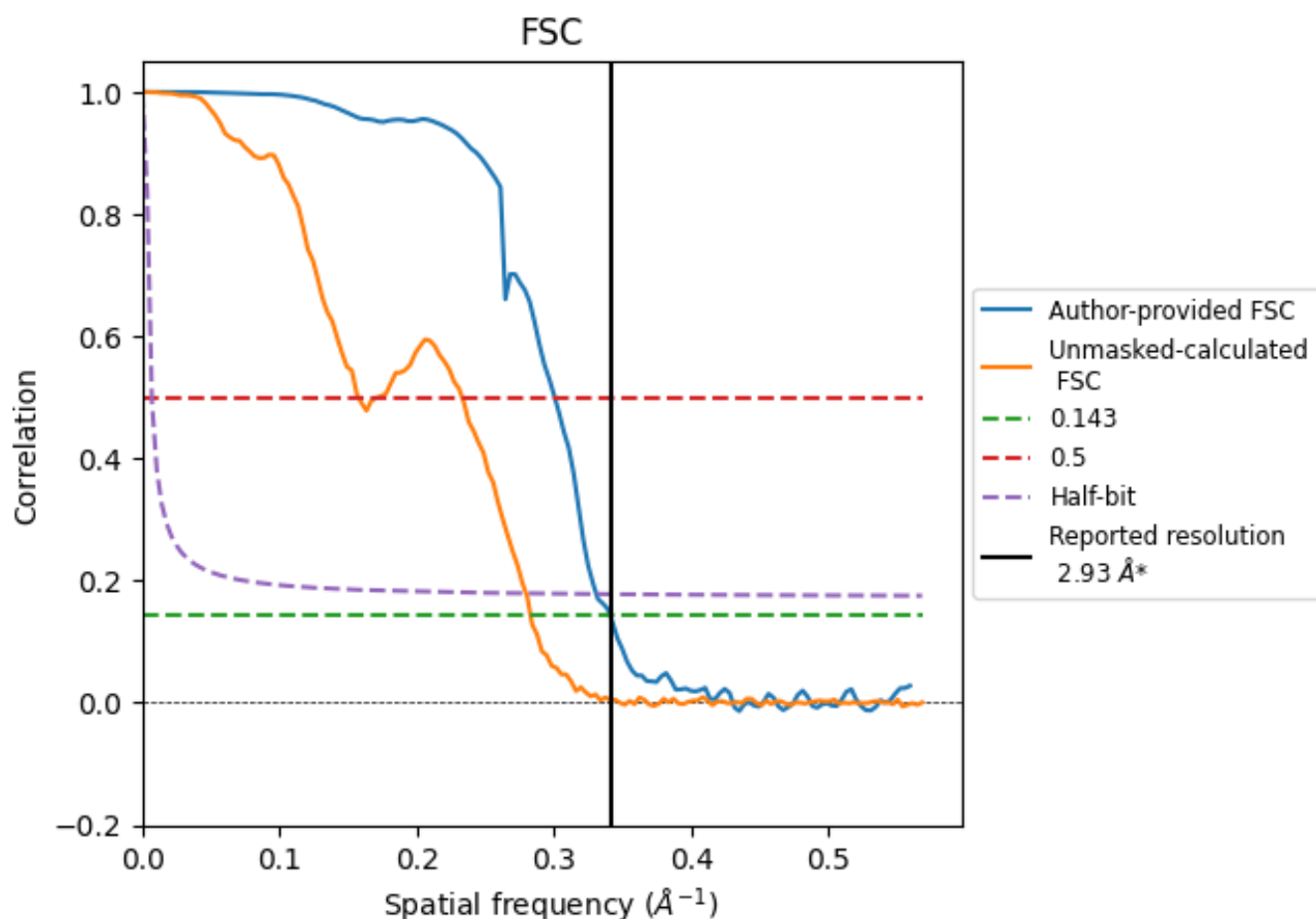


*Reported resolution corresponds to spatial frequency of 0.341 $\text{\AA}^{-1}$

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 0.341 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

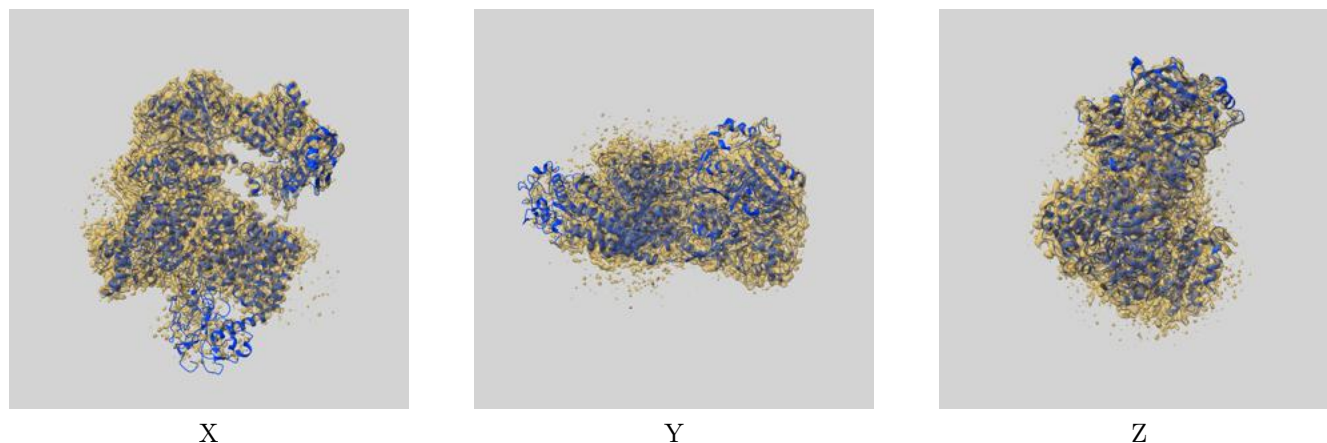
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.93	-	-
Author-provided FSC curve	2.93	3.33	3.02
Unmasked-calculated*	3.53	6.35	3.58

*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 3.53 differs from the reported value 2.93 by more than 10 %

9 Map-model fit [i](#)

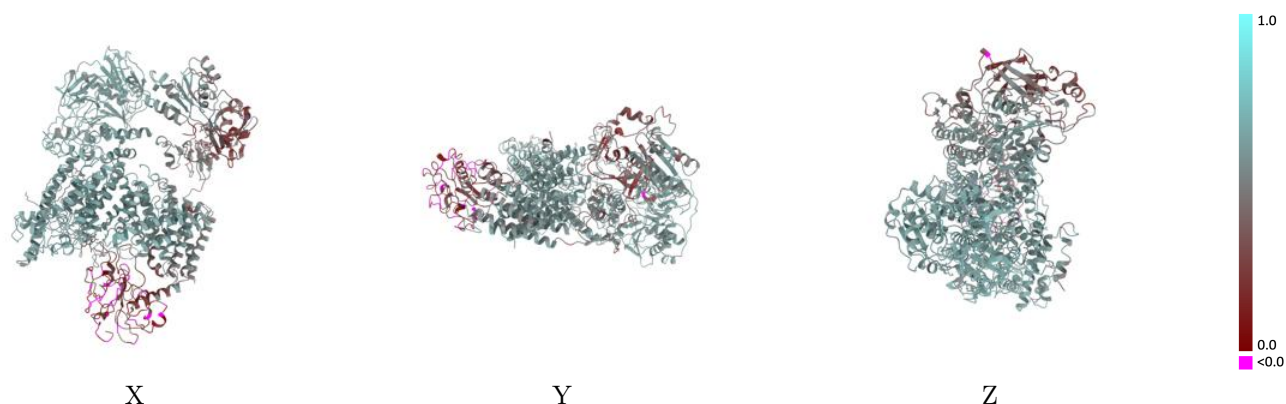
This section contains information regarding the fit between EMDB map EMD-64068 and PDB model 9UDF. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.488 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](#)

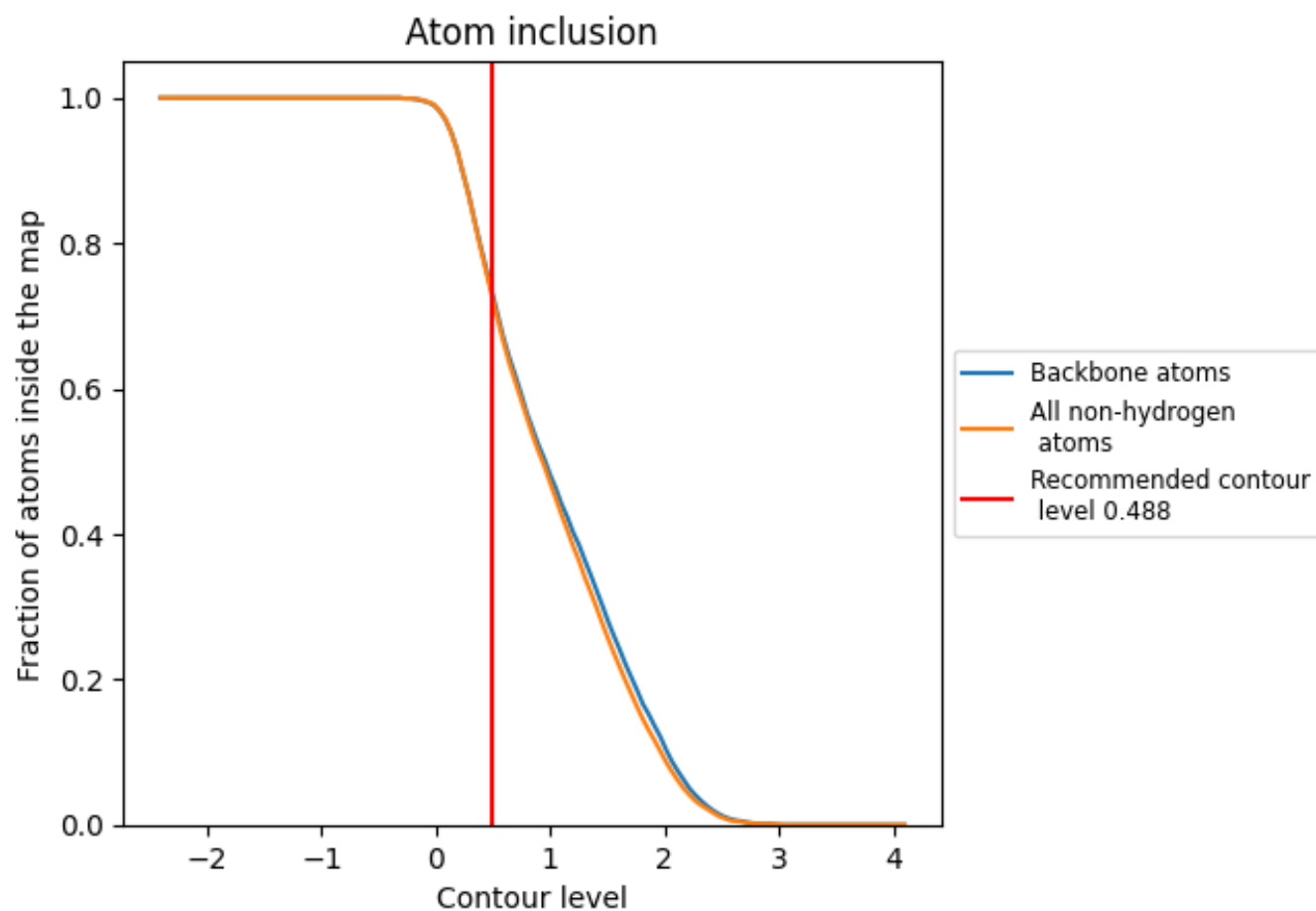


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 73% 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 (0.488) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7280	0.5170
A	0.9260	0.6150
B	0.9150	0.6090
C	0.1520	0.1920
D	0.8410	0.5650
E	0.8940	0.5950
F	0.5400	0.4500

1.0

0.0

<0.0