



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 02:47 PM UTC

PDB ID : 9GY7 / pdb_00009gy7
EMDB ID : EMD-51691
Title : C. thermocellum UvrA-UvrB in complex with DNA with a fluorescein modification and AMPPNP
Authors : Nirwal, S.; Czarnocki-Cieciura, M.; Nowotny, M.
Deposited on : 2024-10-01
Resolution : 3.18 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:

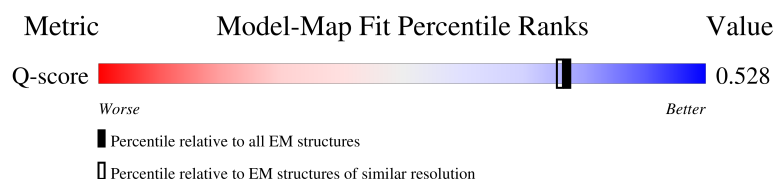
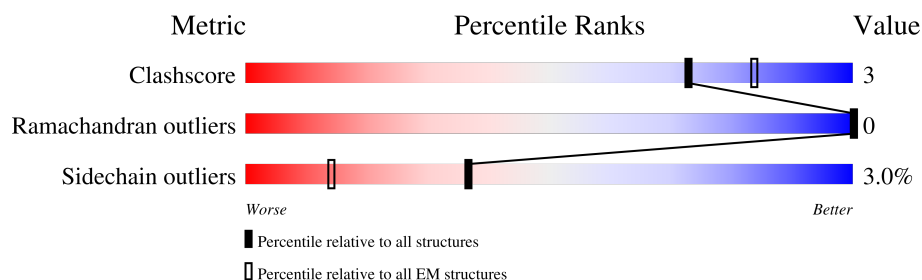
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.18 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	14470 (2.68 - 3.68)

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	962	<div> <div>11%</div> <div>70%</div> <div>7%</div> <div>23%</div> </div>
1	B	962	<div> <div>11%</div> <div>70%</div> <div>7%</div> <div>23%</div> </div>
2	C	50	<div> <div>42%</div> <div>58%</div> <div>30%</div> <div>12%</div> </div>
2	D	50	<div> <div>42%</div> <div>56%</div> <div>32%</div> <div>12%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	E	661	<div><div>53%</div><div><div></div><div></div><div></div><div></div></div><div>80%</div><div>7%</div><div>13%</div></div>
3	F	661	<div><div>52%</div><div><div></div><div></div><div></div><div></div></div><div>80%</div><div>7%</div><div>13%</div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 21176 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 UvrABC system protein A.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	743	Total	C	N	O	S	0	0
			5547	3518	983	1030	16		
1	B	743	Total	C	N	O	S	0	0
			5547	3518	983	1030	16		

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A3DC70
A	-18	GLY	-	expression tag	UNP A3DC70
A	-17	SER	-	expression tag	UNP A3DC70
A	-16	SER	-	expression tag	UNP A3DC70
A	-15	HIS	-	expression tag	UNP A3DC70
A	-14	HIS	-	expression tag	UNP A3DC70
A	-13	HIS	-	expression tag	UNP A3DC70
A	-12	HIS	-	expression tag	UNP A3DC70
A	-11	HIS	-	expression tag	UNP A3DC70
A	-10	HIS	-	expression tag	UNP A3DC70
A	-9	SER	-	expression tag	UNP A3DC70
A	-8	SER	-	expression tag	UNP A3DC70
A	-7	GLY	-	expression tag	UNP A3DC70
A	-6	LEU	-	expression tag	UNP A3DC70
A	-5	VAL	-	expression tag	UNP A3DC70
A	-4	PRO	-	expression tag	UNP A3DC70
A	-3	ARG	-	expression tag	UNP A3DC70
A	-2	GLY	-	expression tag	UNP A3DC70
A	-1	SER	-	expression tag	UNP A3DC70
A	0	HIS	-	expression tag	UNP A3DC70
B	-19	MET	-	initiating methionine	UNP A3DC70
B	-18	GLY	-	expression tag	UNP A3DC70
B	-17	SER	-	expression tag	UNP A3DC70
B	-16	SER	-	expression tag	UNP A3DC70
B	-15	HIS	-	expression tag	UNP A3DC70
B	-14	HIS	-	expression tag	UNP A3DC70

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	HIS	-	expression tag	UNP A3DC70
B	-12	HIS	-	expression tag	UNP A3DC70
B	-11	HIS	-	expression tag	UNP A3DC70
B	-10	HIS	-	expression tag	UNP A3DC70
B	-9	SER	-	expression tag	UNP A3DC70
B	-8	SER	-	expression tag	UNP A3DC70
B	-7	GLY	-	expression tag	UNP A3DC70
B	-6	LEU	-	expression tag	UNP A3DC70
B	-5	VAL	-	expression tag	UNP A3DC70
B	-4	PRO	-	expression tag	UNP A3DC70
B	-3	ARG	-	expression tag	UNP A3DC70
B	-2	GLY	-	expression tag	UNP A3DC70
B	-1	SER	-	expression tag	UNP A3DC70
B	0	HIS	-	expression tag	UNP A3DC70

- Molecule 2 is a DNA chain called DNA (50-MER) with a fluorescein modification - (random sequence built in model).

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	44	Total	C	N	O	P	0	0
			884	423	162	258	41		
2	D	44	Total	C	N	O	P	0	0
			884	423	162	258	41		

- Molecule 3 is a protein called UvrABC system protein B.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	576	Total	C	N	O	S	0	0
			4093	2614	713	749	17		
3	F	576	Total	C	N	O	S	0	0
			4093	2614	713	749	17		

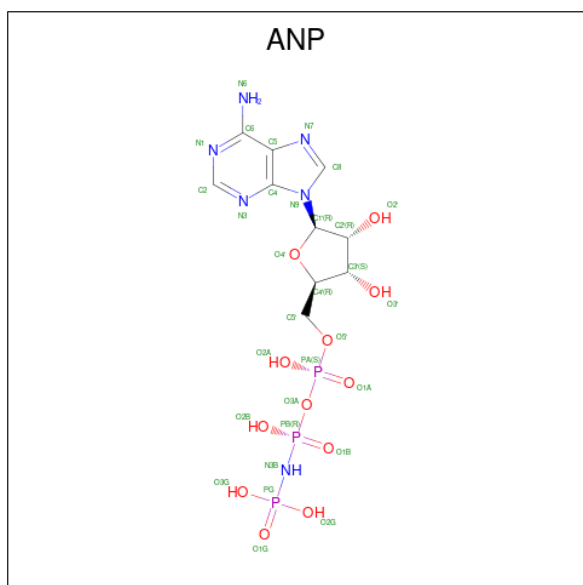
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	0	SER	-	expression tag	UNP A3DC68
F	0	SER	-	expression tag	UNP A3DC68

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Zn	0
			1	1	
4	B	1	Total	Zn	0
			1	1	

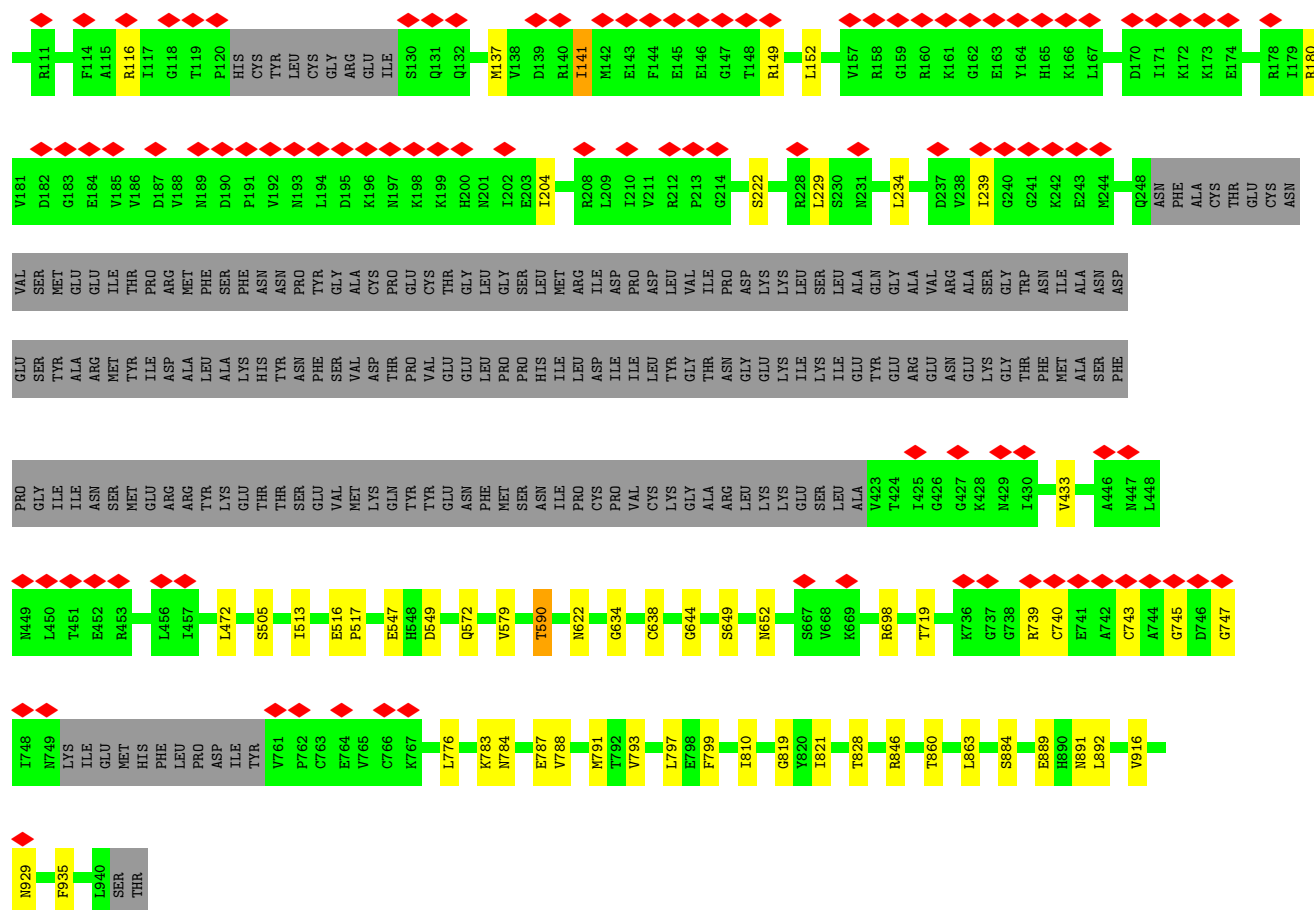
- Molecule 5 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$) (labeled as "Ligand of Interest" by depositor).



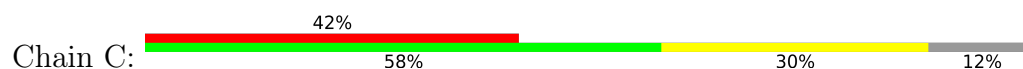
Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total	C	N	O	P	0
			31	10	6	12	3	
5	A	1	Total	C	N	O	P	0
			31	10	6	12	3	
5	B	1	Total	C	N	O	P	0
			31	10	6	12	3	
5	B	1	Total	C	N	O	P	0
			31	10	6	12	3	

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

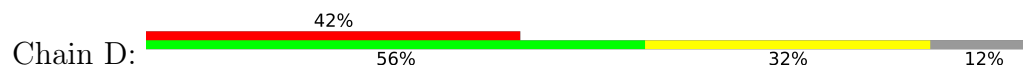
Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total	Mg	0
			1	1	
6	B	1	Total	Mg	0
			1	1	



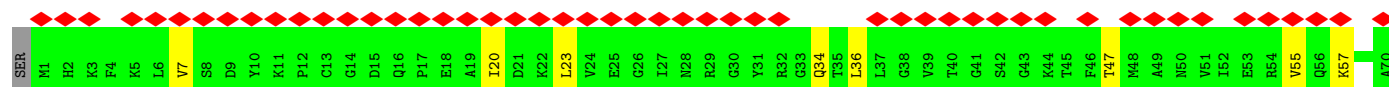
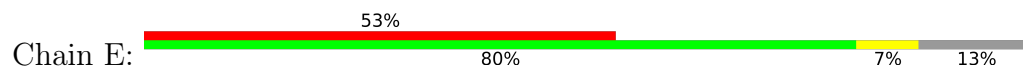
- Molecule 2: DNA (50-MER) with a fluorescein modification - (random sequence built in model)

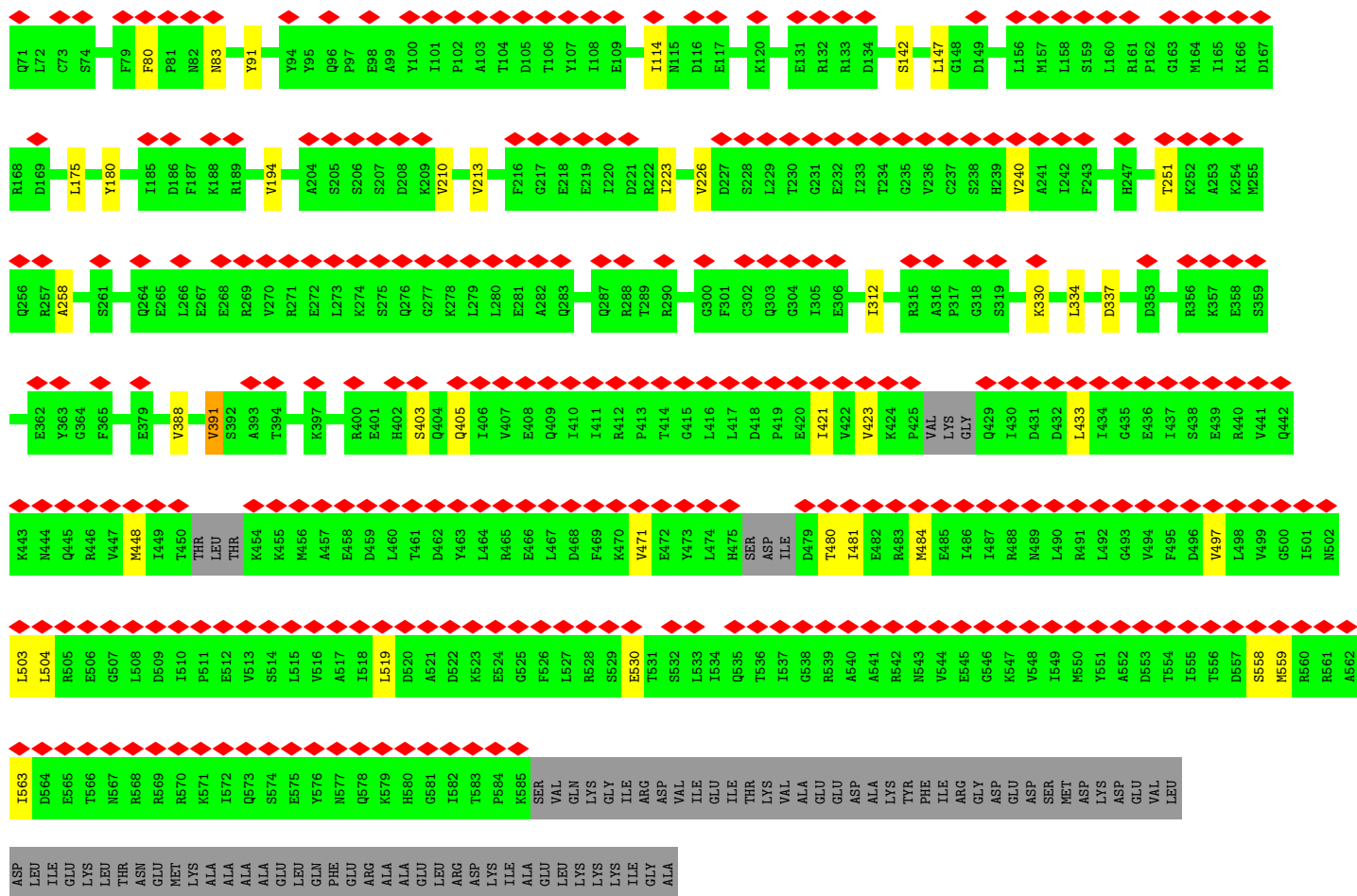


- Molecule 2: DNA (50-MER) with a fluorescein modification - (random sequence built in model)

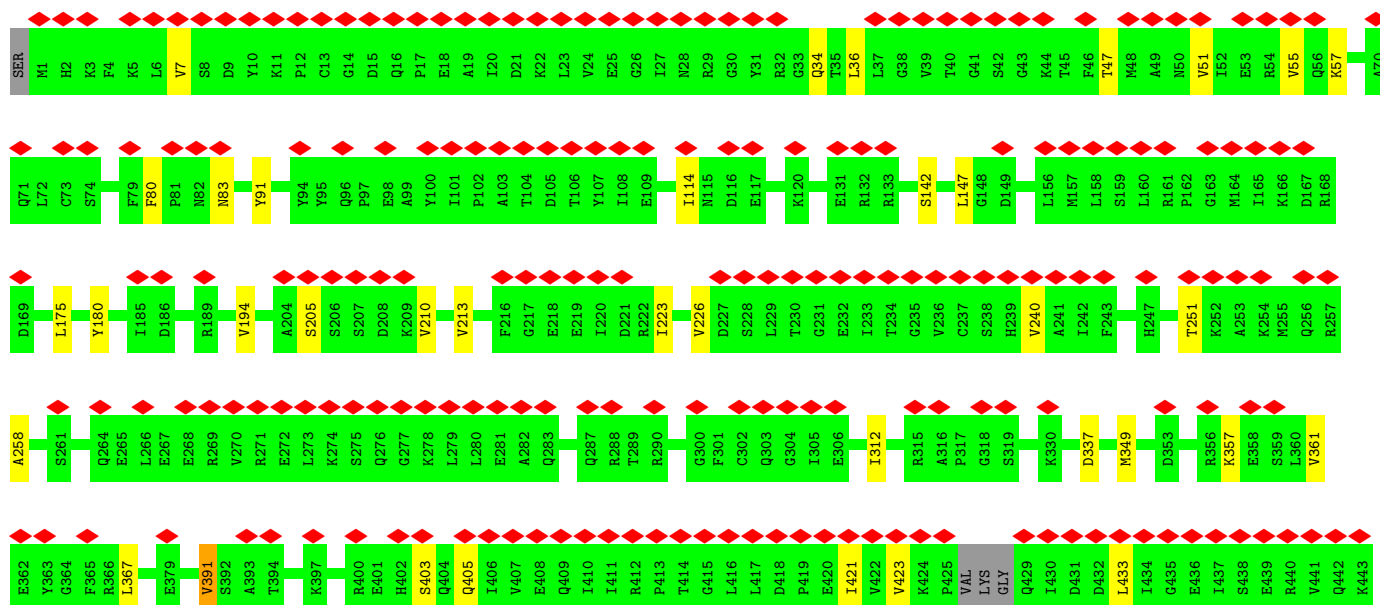
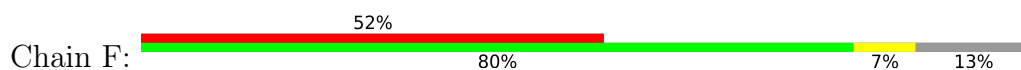


- Molecule 3: UvrABC system protein B





• Molecule 3: UvrABC system protein B





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	136850	Depositor
Resolution determination method	OTHER	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40.73	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.252	Depositor
Minimum map value	0.000	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.037	Depositor
Recommended contour level	0.35	Depositor
Map size (Å)	322.56, 322.56, 322.56	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.84, 0.84, 0.84	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ANP, ZN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.10	0/5620	0.23	0/7600
1	B	0.10	0/5620	0.23	0/7600
2	C	0.17	0/989	0.32	0/1521
2	D	0.17	0/989	0.33	0/1521
3	E	0.08	0/4159	0.21	0/5656
3	F	0.08	0/4159	0.22	0/5656
All	All	0.10	0/21536	0.24	0/29554

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

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	5547	0	5571	36	0
1	B	5547	0	5571	36	0
2	C	884	0	496	10	0
2	D	884	0	496	11	0
3	E	4093	0	3740	24	0
3	F	4093	0	3740	24	0
4	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	A	62	0	26	4	0
5	B	62	0	26	4	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
All	All	21176	0	19666	135	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 3.

All (135) 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:644:GLY:H	5:B:1003:ANP:HNB1	1.32	0.78
1:A:644:GLY:H	5:A:1003:ANP:HNB1	1.32	0.78
1:B:634:GLY:H	1:B:884:SER:HB3	1.56	0.71
1:A:634:GLY:H	1:A:884:SER:HB3	1.56	0.70
1:B:547:GLU:HG3	1:B:549:ASP:H	1.57	0.69
3:F:80:PHE:HB3	3:F:83:ASN:HB2	1.75	0.69
1:A:547:GLU:HG3	1:A:549:ASP:H	1.57	0.67
3:E:80:PHE:HB3	3:E:83:ASN:HB2	1.76	0.67
3:E:421:ILE:HG21	3:E:563:ILE:HD12	1.80	0.64
3:F:421:ILE:HG21	3:F:563:ILE:HD12	1.80	0.64
3:F:433:LEU:HD11	3:F:519:LEU:HD21	1.79	0.64
3:E:433:LEU:HD11	3:E:519:LEU:HD21	1.81	0.63
3:E:481:ILE:HA	3:E:484:MET:HE2	1.81	0.61
1:A:889:GLU:HG3	1:A:891:ASN:H	1.66	0.61
2:C:3:DG:H2"	2:C:4:DT:H5"	1.83	0.61
1:B:889:GLU:HG3	1:B:891:ASN:H	1.65	0.60
1:A:67:LEU:HD11	1:B:67:LEU:HD11	1.84	0.59
2:D:-3:DC:H5"	3:F:142:SER:HB2	1.85	0.59
1:A:36:GLY:H	5:A:1002:ANP:HNB1	1.51	0.59
3:F:481:ILE:HA	3:F:484:MET:HE2	1.86	0.58
1:A:40:SER:OG	5:A:1002:ANP:O1B	2.22	0.58
1:B:36:GLY:H	5:B:1002:ANP:HNB1	1.51	0.58
1:A:152:LEU:HD12	1:A:222:SER:HB3	1.86	0.58
1:B:40:SER:OG	5:B:1002:ANP:O1B	2.22	0.57
3:E:337:ASP:HA	3:E:391:VAL:HG23	1.86	0.57
3:F:337:ASP:HA	3:F:391:VAL:HG23	1.86	0.57
1:B:783:LYS:HG3	1:B:787:GLU:HB2	1.87	0.56
1:B:152:LEU:HD12	1:B:222:SER:HB3	1.86	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:3:DG:H2''	2:D:4:DT:H5''	1.87	0.56
1:A:41:SER:HA	1:A:45:ASP:HB2	1.88	0.54
2:C:-3:DC:H5''	3:E:142:SER:HB2	1.90	0.54
1:A:698:ARG:HD2	2:C:11:DT:H5'	1.89	0.54
1:B:41:SER:HA	1:B:45:ASP:HB2	1.89	0.53
3:F:357:LYS:O	3:F:361:VAL:HG23	2.08	0.53
1:B:698:ARG:HD2	2:D:11:DT:H5'	1.91	0.53
1:A:32:THR:HG22	1:A:590:THR:HG21	1.90	0.53
3:F:47:THR:O	3:F:51:VAL:HG12	2.09	0.53
1:B:32:THR:HG22	1:B:590:THR:HG21	1.90	0.52
1:A:776:LEU:HB3	1:A:784:ASN:HD22	1.75	0.52
1:A:48:TYR:HB2	1:A:513:ILE:HD13	1.92	0.52
1:B:776:LEU:HB3	1:B:784:ASN:HD22	1.75	0.52
1:B:48:TYR:HB2	1:B:513:ILE:HD13	1.92	0.51
1:A:783:LYS:HG3	1:A:787:GLU:HB2	1.91	0.51
3:E:55:VAL:HG12	3:E:57:LYS:HB2	1.93	0.51
2:C:39:DT:H2'	2:C:40:DG:C8	2.46	0.50
1:A:149:ARG:HB2	1:A:239:ILE:HD12	1.93	0.50
1:A:797:LEU:HG	1:A:810:ILE:HG22	1.94	0.50
3:F:480:THR:O	3:F:484:MET:HG2	2.11	0.50
2:D:39:DT:H2'	2:D:40:DG:C8	2.47	0.49
1:B:149:ARG:HB2	1:B:239:ILE:HD12	1.93	0.49
1:B:797:LEU:HG	1:B:810:ILE:HG22	1.94	0.49
3:E:480:THR:O	3:E:484:MET:HG2	2.13	0.49
3:F:55:VAL:HG12	3:F:57:LYS:HB2	1.94	0.49
3:F:559:MET:O	3:F:563:ILE:HG12	2.13	0.49
3:E:559:MET:O	3:E:563:ILE:HG12	2.13	0.48
1:A:788:VAL:HA	1:A:791:MET:HE3	1.94	0.48
2:C:5:DT:H2''	2:C:6:DA:C8	2.49	0.48
1:B:788:VAL:HA	1:B:791:MET:HE3	1.96	0.47
2:D:5:DT:H2''	2:D:6:DA:C8	2.50	0.47
1:A:204:ILE:HD12	1:A:229:LEU:HD21	1.97	0.46
1:B:793:VAL:HB	1:B:819:GLY:HA2	1.97	0.46
1:B:517:PRO:HD2	1:B:547:GLU:OE1	2.16	0.46
1:B:7:VAL:HG22	1:B:22:LYS:HG2	1.98	0.46
1:A:793:VAL:HB	1:A:819:GLY:HA2	1.97	0.46
1:A:791:MET:HE1	1:A:799:PHE:CD2	2.51	0.46
3:F:91:TYR:HE2	3:F:114:ILE:HD12	1.81	0.46
3:F:530:GLU:HB2	3:F:558:SER:HB2	1.97	0.46
1:A:784:ASN:OD1	1:A:787:GLU:HG3	2.15	0.46
1:A:517:PRO:HD2	1:A:547:GLU:OE1	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:791:MET:HE1	1:B:799:PHE:CD2	2.51	0.45
1:A:783:LYS:HB2	1:A:783:LYS:HE3	1.61	0.45
1:A:19:ILE:HG21	1:A:572:GLN:HA	1.99	0.45
1:A:7:VAL:HG22	1:A:22:LYS:HG2	1.98	0.45
2:D:2:DT:H2''	2:D:3:DG:H5''	1.98	0.45
3:E:448:MET:SD	3:E:503:LEU:HD21	2.56	0.45
1:B:204:ILE:HD12	1:B:229:LEU:HD21	1.97	0.45
3:E:530:GLU:HB2	3:E:558:SER:HB2	1.98	0.45
2:C:2:DT:H2''	2:C:3:DG:H5''	1.99	0.45
3:F:448:MET:SD	3:F:503:LEU:HD21	2.57	0.45
1:A:644:GLY:N	5:A:1003:ANP:HNB1	2.09	0.45
1:A:743:CYS:HB2	1:A:747:GLY:H	1.82	0.45
1:B:19:ILE:HG21	1:B:572:GLN:HA	1.99	0.45
3:E:213:VAL:HG12	3:E:223:ILE:HG12	1.99	0.45
1:B:743:CYS:HB2	1:B:747:GLY:H	1.82	0.44
3:E:91:TYR:HE2	3:E:114:ILE:HD12	1.81	0.44
1:A:740:CYS:HB3	1:A:743:CYS:SG	2.57	0.44
1:B:644:GLY:N	5:B:1003:ANP:HNB1	2.09	0.44
3:F:213:VAL:HG12	3:F:223:ILE:HG12	1.99	0.44
1:B:740:CYS:HB3	1:B:743:CYS:SG	2.57	0.44
2:D:4:DT:C6	2:D:5:DT:H72	2.52	0.44
1:B:791:MET:HE1	1:B:799:PHE:CG	2.53	0.43
2:D:-4:DT:O2	3:F:349:MET:HE1	2.18	0.43
1:A:152:LEU:HD22	1:A:234:LEU:HD11	2.01	0.43
1:A:739:ARG:HD2	1:A:745:GLY:HA3	2.01	0.43
1:A:81:LEU:HD23	1:A:81:LEU:HA	1.86	0.42
2:C:20:DT:H2''	2:C:21:DG:C8	2.54	0.42
3:E:20:ILE:HG13	3:E:47:THR:HG23	2.01	0.42
3:E:175:LEU:O	3:E:180:TYR:HB2	2.18	0.42
2:D:20:DT:H2''	2:D:21:DG:C8	2.54	0.42
3:E:147:LEU:HD23	3:E:147:LEU:HA	1.87	0.42
3:F:175:LEU:O	3:F:180:TYR:HB2	2.18	0.42
1:B:137:MET:O	1:B:141:ILE:HD12	2.20	0.42
2:C:36:DC:H2''	2:C:37:DG:C8	2.54	0.42
3:E:36:LEU:HB3	3:E:391:VAL:HG12	2.01	0.42
1:A:110:LEU:HD23	1:A:110:LEU:HA	1.84	0.42
2:D:36:DC:H2''	2:D:37:DG:C8	2.54	0.42
3:E:504:LEU:HD23	3:E:504:LEU:HA	1.83	0.42
2:C:4:DT:C6	2:C:5:DT:H72	2.55	0.42
1:B:9:LYS:HB3	1:B:9:LYS:HE3	1.73	0.42
1:B:739:ARG:HD2	1:B:745:GLY:HA3	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:36:LEU:HB3	3:F:391:VAL:HG12	2.02	0.41
1:A:791:MET:HE1	1:A:799:PHE:CG	2.56	0.41
3:E:330:LYS:HA	3:E:330:LYS:HD2	1.81	0.41
3:E:258:ALA:HB2	3:E:312:ILE:HA	2.02	0.41
3:E:334:LEU:HD23	3:E:388:VAL:HG23	2.02	0.41
1:B:152:LEU:HD22	1:B:234:LEU:HD11	2.02	0.41
1:A:614:LEU:HD12	1:A:633:LEU:HD21	2.03	0.41
3:F:34:GLN:HG2	3:F:405:GLN:HB3	2.02	0.41
3:F:367:LEU:HD23	3:F:367:LEU:HA	1.96	0.41
1:A:622:ASN:HD22	1:A:649:SER:HB2	1.85	0.41
3:E:23:LEU:HD11	3:E:36:LEU:HD13	2.03	0.41
3:E:34:GLN:HG2	3:E:405:GLN:HB3	2.03	0.41
3:F:147:LEU:HD23	3:F:147:LEU:HA	1.86	0.41
1:A:805:ARG:H	1:A:805:ARG:HG3	1.73	0.41
3:E:210:VAL:HG22	3:E:226:VAL:O	2.21	0.41
1:B:622:ASN:HD22	1:B:649:SER:HB2	1.86	0.40
2:D:8:DG:H2"	2:D:9:DG:H8	1.86	0.40
2:C:13:DC:H2"	2:C:14:DA:C8	2.56	0.40
1:B:104:THR:HG21	1:B:472:LEU:HD11	2.03	0.40
1:B:892:LEU:HD11	1:B:935:PHE:HB3	2.04	0.40
3:F:258:ALA:HB2	3:F:312:ILE:HA	2.02	0.40
3:F:504:LEU:HA	3:F:539:ARG:HH21	1.86	0.40
1:B:783:LYS:HE3	1:B:783:LYS:HB2	1.61	0.40
1:B:860:THR:O	1:B:863:LEU:HB2	2.21	0.40
3:F:210:VAL:HG22	3:F:226:VAL:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	735/962 (76%)	714 (97%)	21 (3%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	735/962 (76%)	715 (97%)	20 (3%)	0	100	100
3	E	568/661 (86%)	558 (98%)	10 (2%)	0	100	100
3	F	568/661 (86%)	558 (98%)	10 (2%)	0	100	100
All	All	2606/3246 (80%)	2545 (98%)	61 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	564/823 (68%)	546 (97%)	18 (3%)	34	62
1	B	564/823 (68%)	545 (97%)	19 (3%)	32	61
3	E	356/584 (61%)	347 (98%)	9 (2%)	42	66
3	F	356/584 (61%)	346 (97%)	10 (3%)	38	64
All	All	1840/2814 (65%)	1784 (97%)	56 (3%)	37	63

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	GLN
1	A	82	SER
1	A	92	THR
1	A	116	ARG
1	A	180	ARG
1	A	433	VAL
1	A	505	SER
1	A	508	MET
1	A	516	GLU
1	A	579	VAL
1	A	590	THR
1	A	638	CYS
1	A	652	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	719	THR
1	A	821	ILE
1	A	828	THR
1	A	846	ARG
1	A	916	VAL
1	B	52	GLN
1	B	82	SER
1	B	92	THR
1	B	116	ARG
1	B	141	ILE
1	B	180	ARG
1	B	433	VAL
1	B	505	SER
1	B	516	GLU
1	B	579	VAL
1	B	590	THR
1	B	638	CYS
1	B	652	ASN
1	B	719	THR
1	B	821	ILE
1	B	828	THR
1	B	846	ARG
1	B	916	VAL
1	B	929	ASN
3	E	7	VAL
3	E	194	VAL
3	E	240	VAL
3	E	251	THR
3	E	391	VAL
3	E	403	SER
3	E	423	VAL
3	E	471	VAL
3	E	497	VAL
3	F	7	VAL
3	F	194	VAL
3	F	205	SER
3	F	240	VAL
3	F	251	THR
3	F	391	VAL
3	F	403	SER
3	F	423	VAL
3	F	471	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	F	497	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	GLN
1	A	201	ASN
1	A	216	GLN
1	A	466	ASN
1	A	622	ASN
1	A	628	ASN
1	A	876	HIS
1	B	151	GLN
1	B	201	ASN
1	B	216	GLN
1	B	466	ASN
1	B	622	ASN
1	B	628	ASN
1	B	876	HIS
3	E	2	HIS
3	E	28	ASN
3	E	34	GLN
3	E	115	ASN
3	E	239	HIS
3	E	247	HIS
3	E	307	ASN
3	E	345	GLN
3	E	405	GLN
3	F	2	HIS
3	F	28	ASN
3	F	64	HIS
3	F	115	ASN
3	F	239	HIS
3	F	247	HIS
3	F	307	ASN
3	F	345	GLN
3	F	405	GLN

5.3.3 RNA ⓘ

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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
5	ANP	B	1003	-	33,33,33	1.19	4 (12%)	45,52,52	0.56	0
5	ANP	A	1002	6	33,33,33	1.20	4 (12%)	45,52,52	0.63	1 (2%)
5	ANP	A	1003	-	33,33,33	1.19	4 (12%)	45,52,52	0.56	0
5	ANP	B	1002	6	33,33,33	1.20	4 (12%)	45,52,52	0.63	1 (2%)

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
5	ANP	B	1003	-	-	3/18/38/38	0/3/3/3
5	ANP	A	1002	6	-	4/18/38/38	0/3/3/3
5	ANP	A	1003	-	-	3/18/38/38	0/3/3/3
5	ANP	B	1002	6	-	4/18/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1003	ANP	PB-O2B	-3.42	1.47	1.56

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1003	ANP	PB-O2B	-3.42	1.47	1.56
5	B	1002	ANP	PB-O2B	-3.38	1.47	1.56
5	A	1002	ANP	PB-O2B	-3.38	1.47	1.56
5	B	1002	ANP	PG-O2G	-3.30	1.48	1.56
5	A	1002	ANP	PG-O2G	-3.30	1.48	1.56
5	B	1002	ANP	PG-O3G	-3.26	1.48	1.56
5	A	1002	ANP	PG-O3G	-3.26	1.48	1.56
5	A	1003	ANP	PG-O2G	-3.25	1.48	1.56
5	B	1003	ANP	PG-O2G	-3.24	1.48	1.56
5	A	1003	ANP	PG-O3G	-3.23	1.48	1.56
5	B	1003	ANP	PG-O3G	-3.23	1.48	1.56
5	A	1003	ANP	PB-O3A	2.94	1.62	1.59
5	B	1003	ANP	PB-O3A	2.93	1.62	1.59
5	B	1002	ANP	PB-O3A	2.88	1.62	1.59
5	A	1002	ANP	PB-O3A	2.88	1.62	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
5	A	1002	ANP	O2B-PB-O1B	2.48	115.18	109.87
5	B	1002	ANP	O2B-PB-O1B	2.47	115.17	109.87

There are no chirality outliers.

All (14) torsion outliers are listed below:

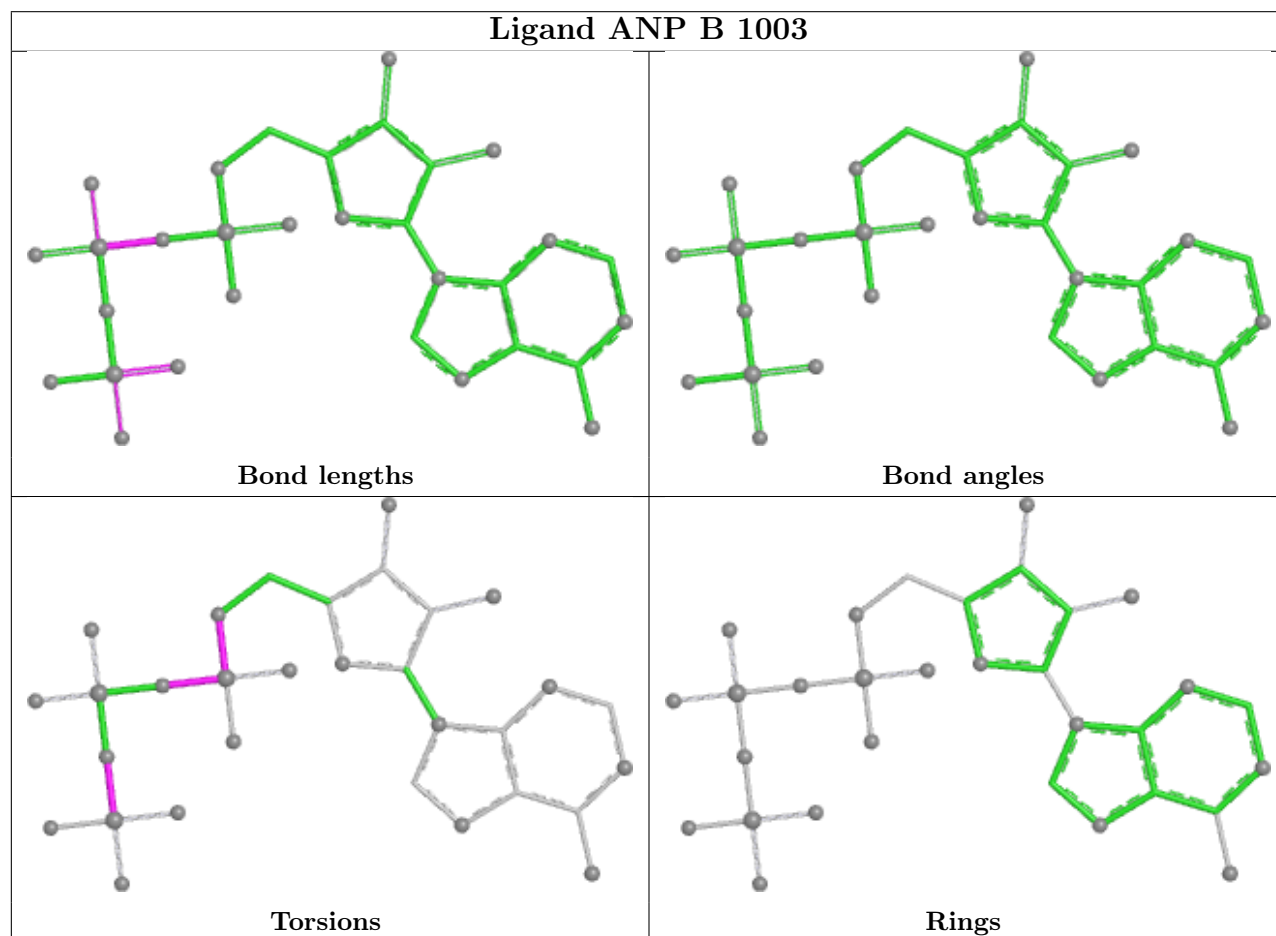
Mol	Chain	Res	Type	Atoms
5	A	1002	ANP	PB-N3B-PG-O1G
5	A	1002	ANP	PA-O3A-PB-O2B
5	A	1003	ANP	PB-N3B-PG-O1G
5	A	1003	ANP	C5'-O5'-PA-O1A
5	B	1002	ANP	PB-N3B-PG-O1G
5	B	1002	ANP	PA-O3A-PB-O2B
5	B	1003	ANP	PB-N3B-PG-O1G
5	B	1003	ANP	C5'-O5'-PA-O1A
5	A	1003	ANP	PB-O3A-PA-O2A
5	B	1003	ANP	PB-O3A-PA-O2A
5	A	1002	ANP	C2'-C1'-N9-C8
5	B	1002	ANP	C2'-C1'-N9-C8
5	A	1002	ANP	PA-O3A-PB-O1B
5	B	1002	ANP	PA-O3A-PB-O1B

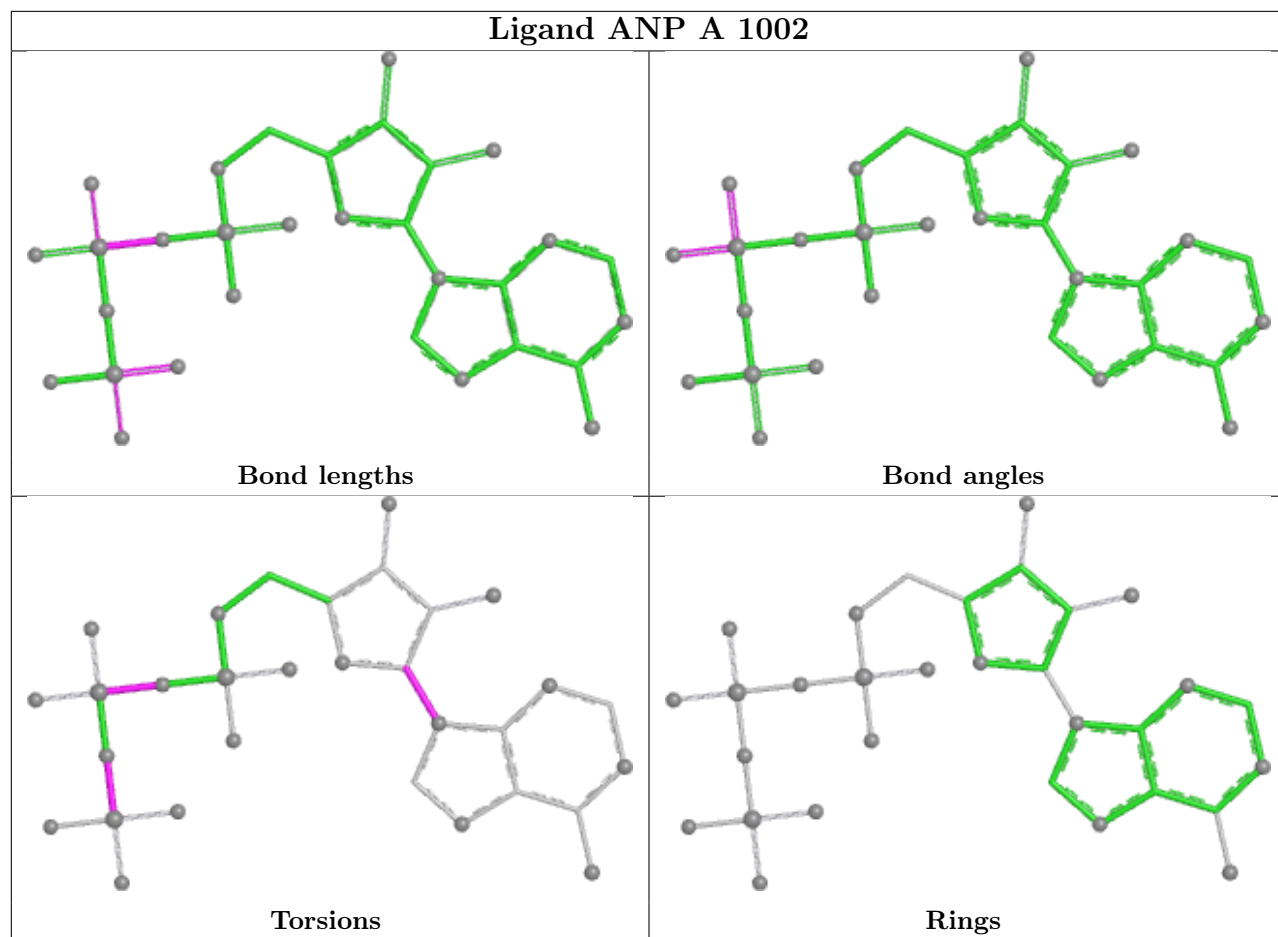
There are no ring outliers.

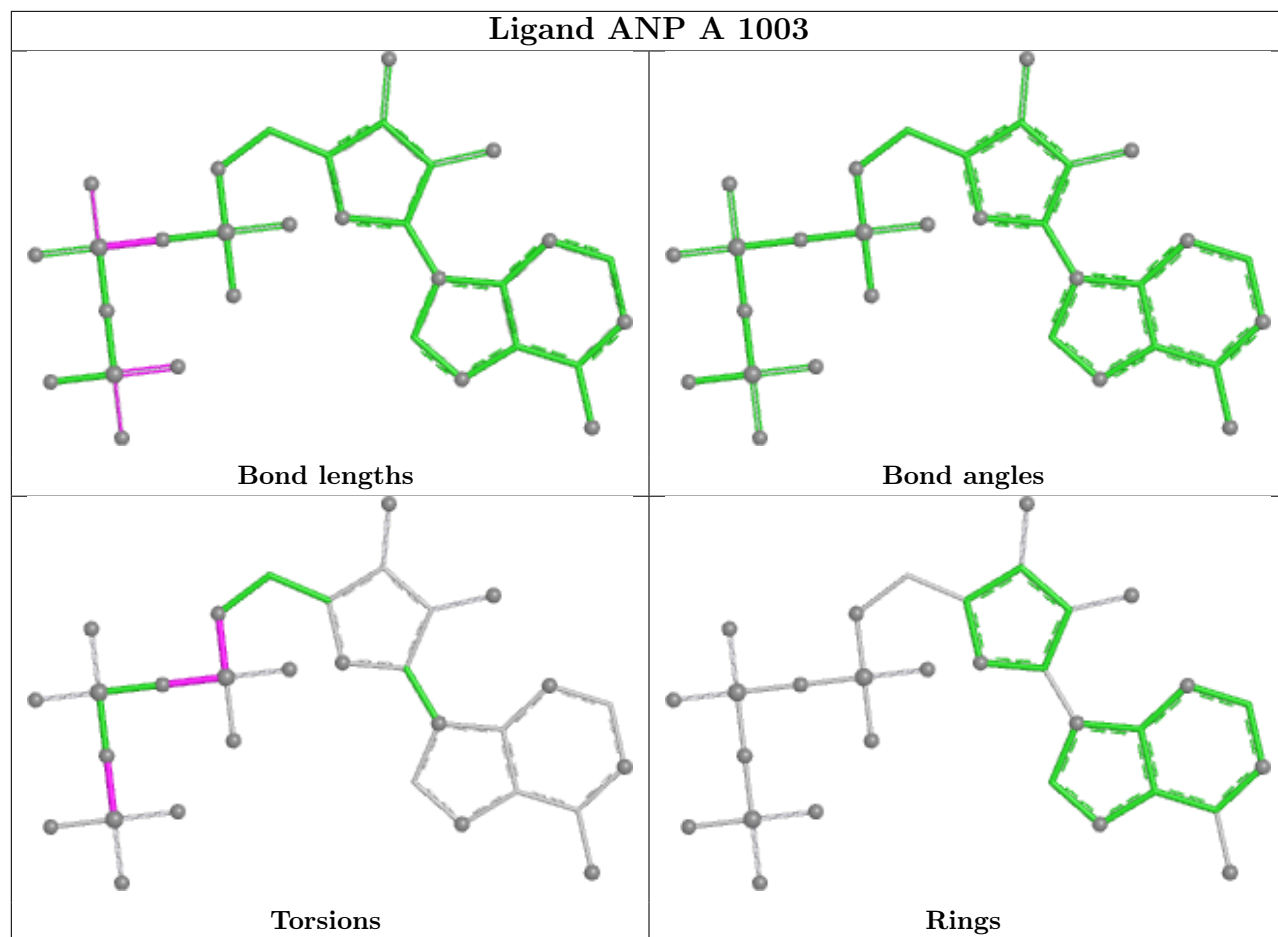
4 monomers are involved in 8 short contacts:

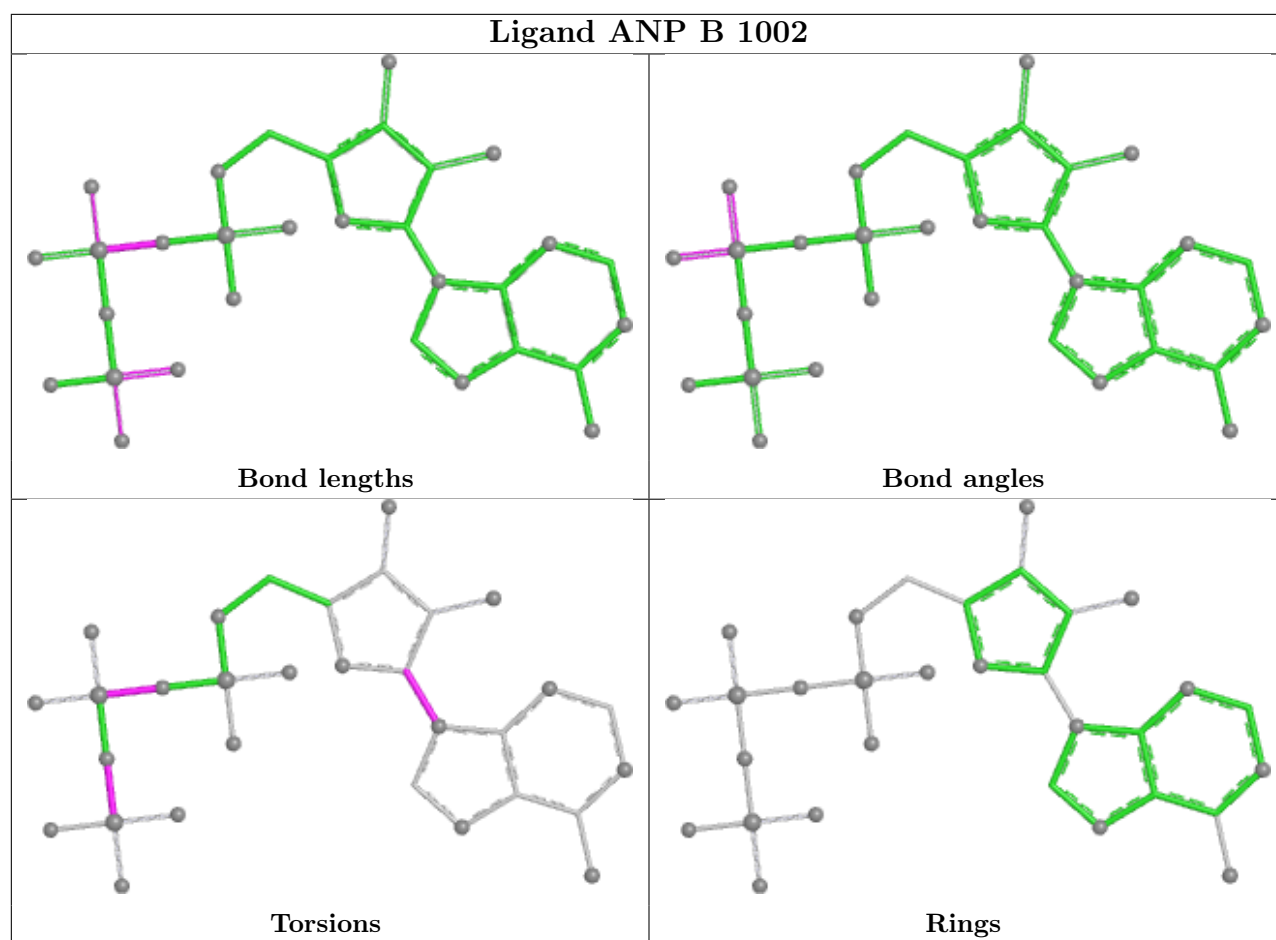
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1003	ANP	2	0
5	A	1002	ANP	2	0
5	A	1003	ANP	2	0
5	B	1002	ANP	2	0

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.

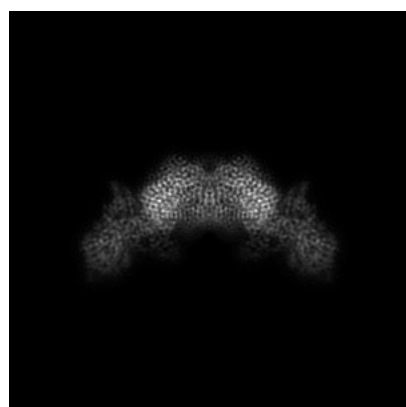
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-51691. These allow visual inspection of the internal detail of the map and identification of artifacts.

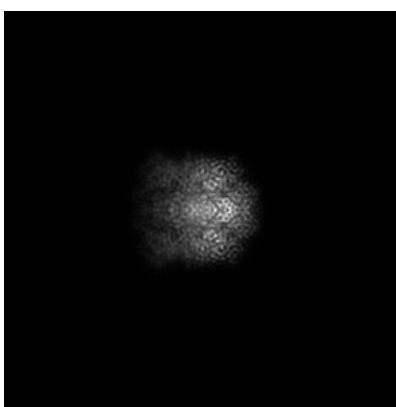
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

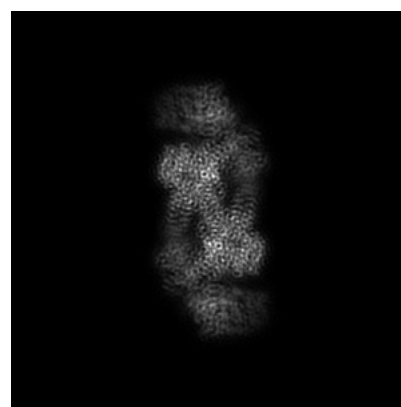
6.1.1 Primary map



X



Y



Z

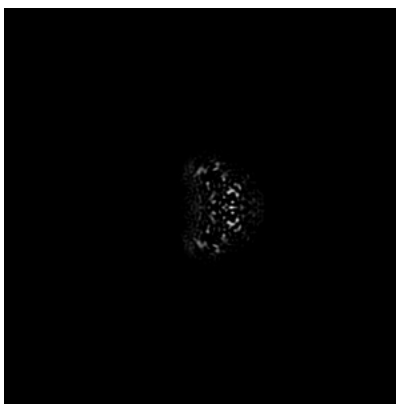
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 192



Y Index: 192



Z Index: 192

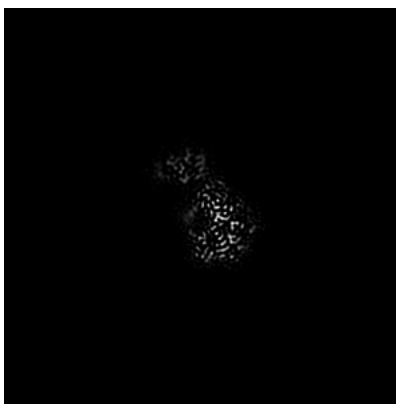
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 194



Y Index: 235

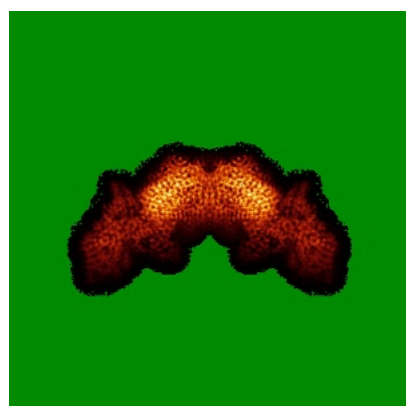


Z Index: 204

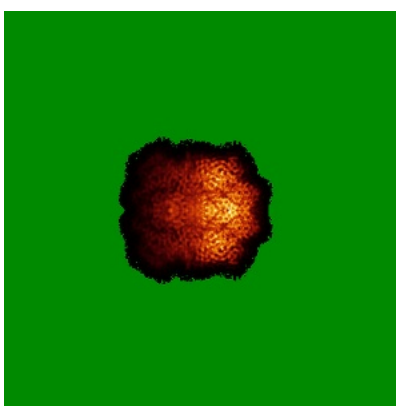
The images above show the largest variance slices of the map in three orthogonal directions.

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

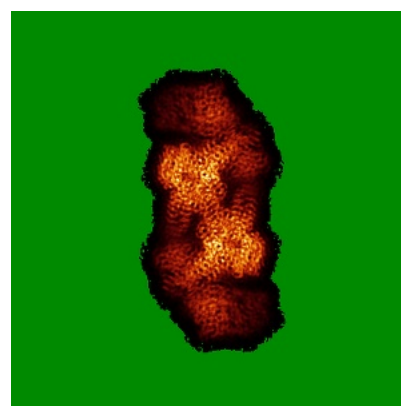
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.35. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

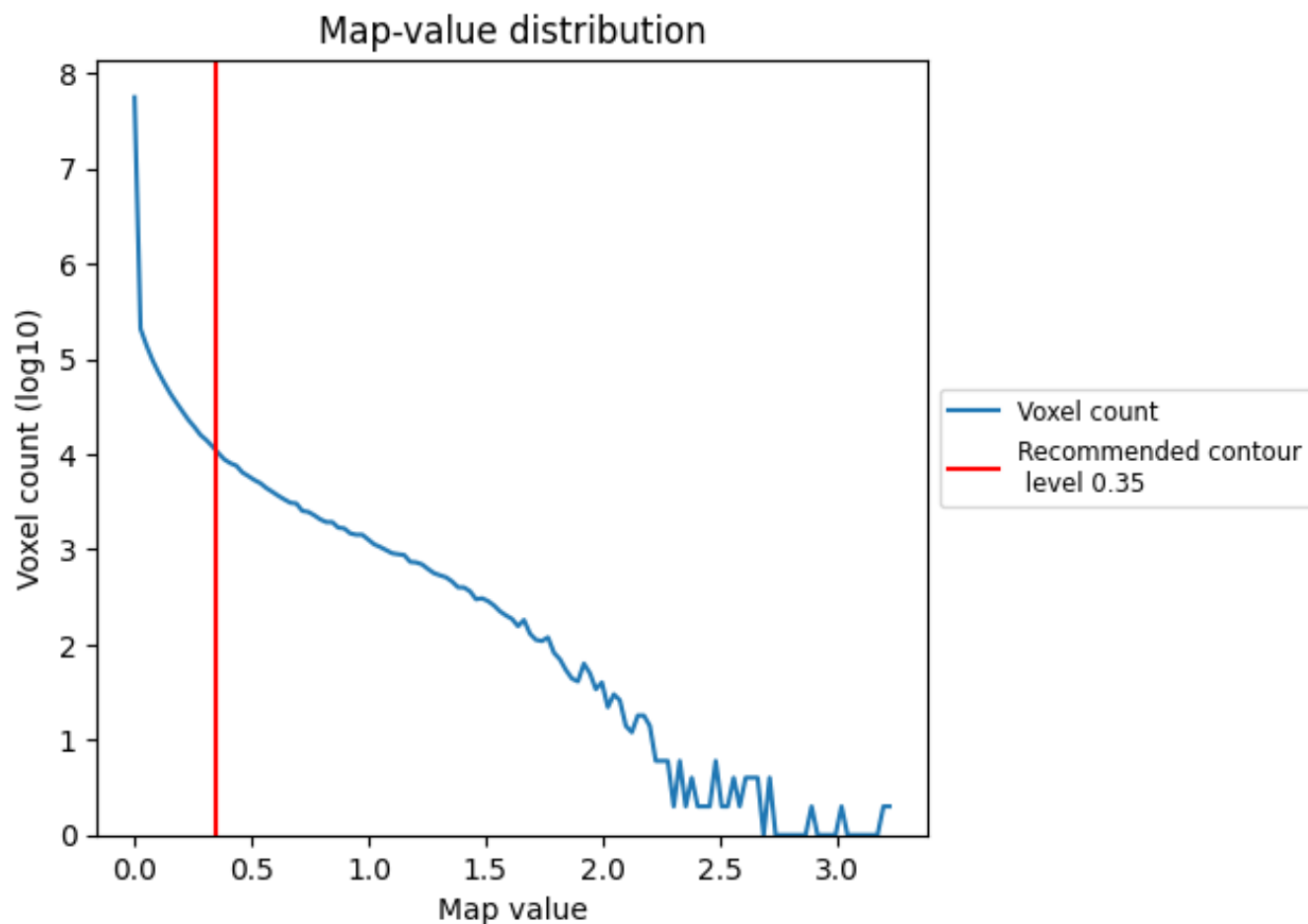
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

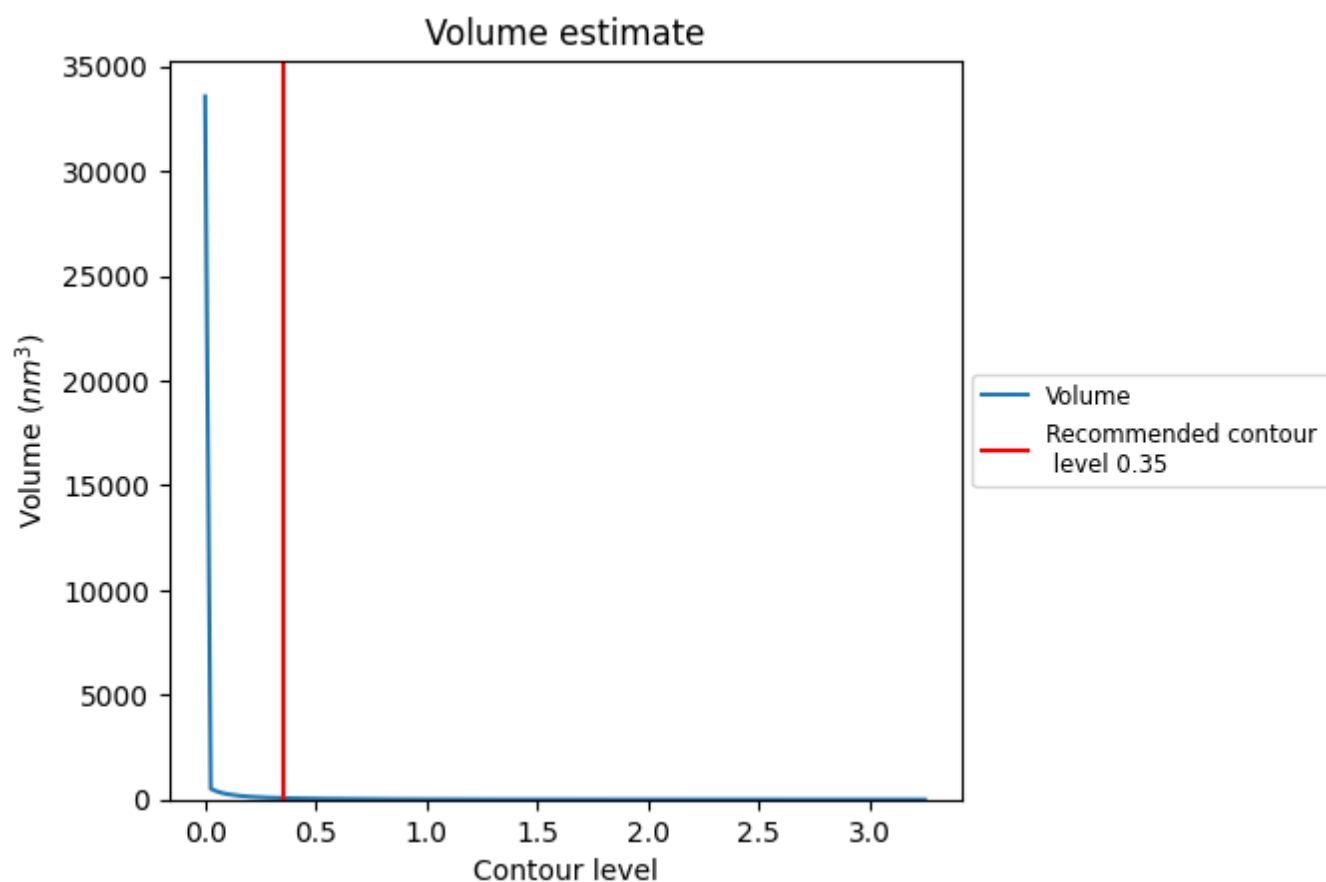
This section contains the results of statistical analysis of the map.

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.

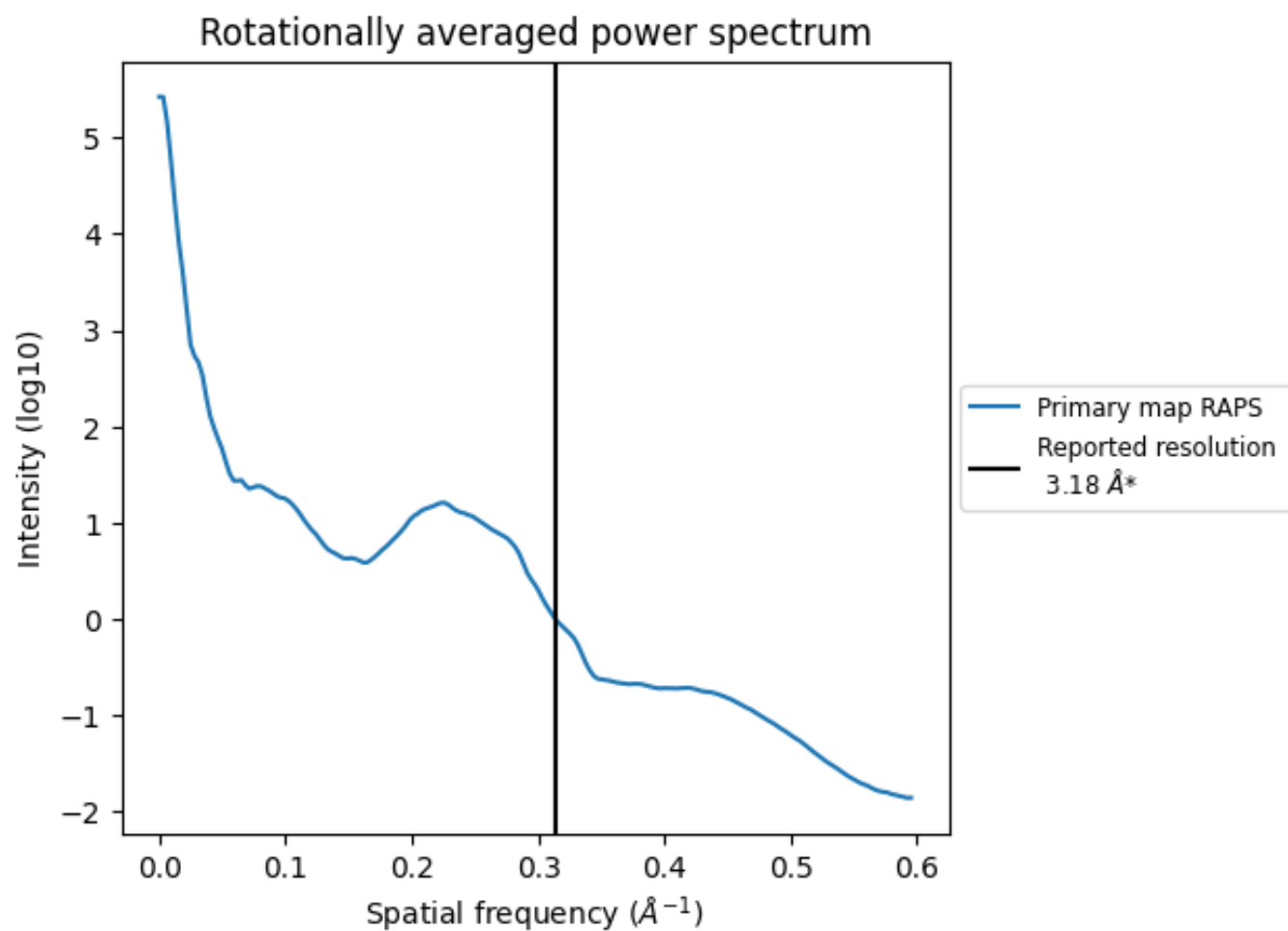
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 71 nm^3 ; this corresponds to an approximate mass of 64 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 0.314 Å⁻¹

8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

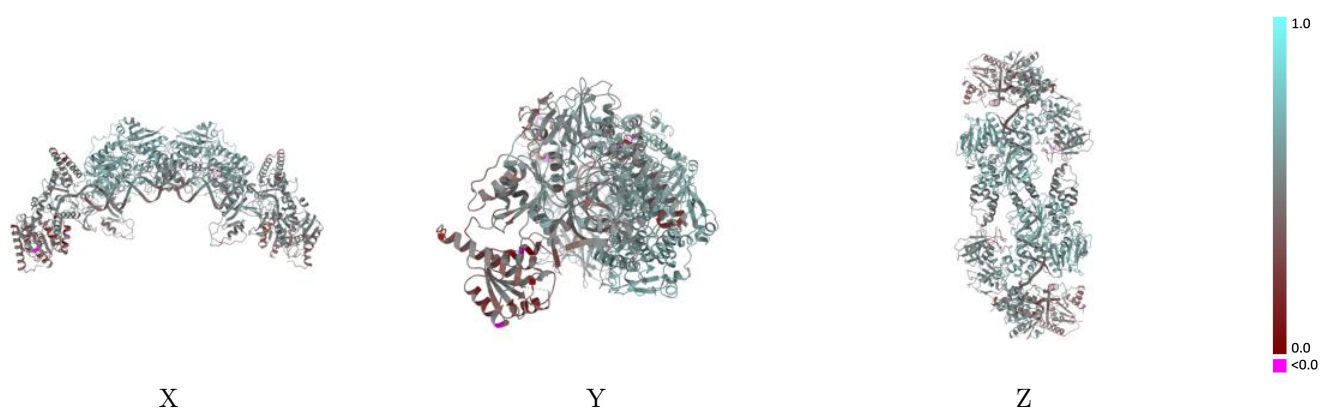
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-51691 and PDB model 9GY7. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)

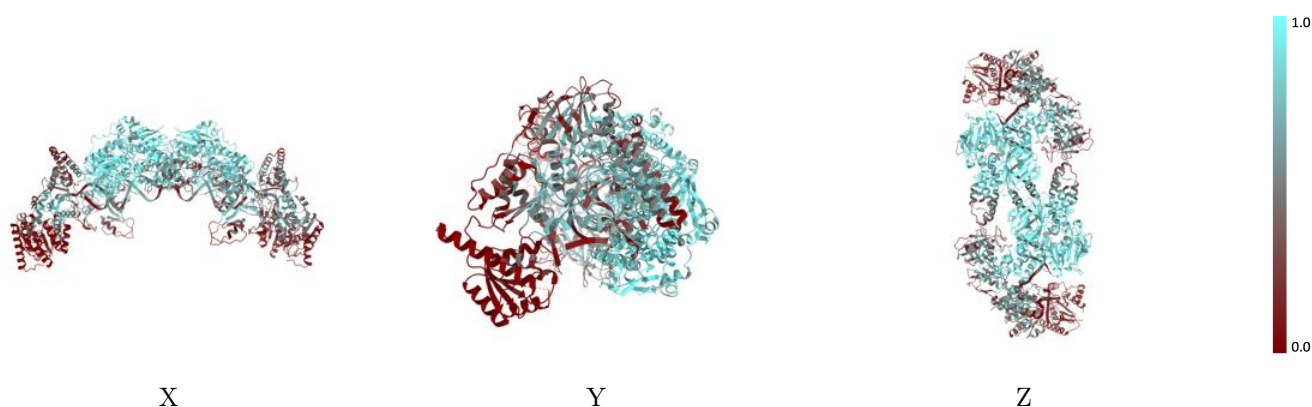
This section was not generated.

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



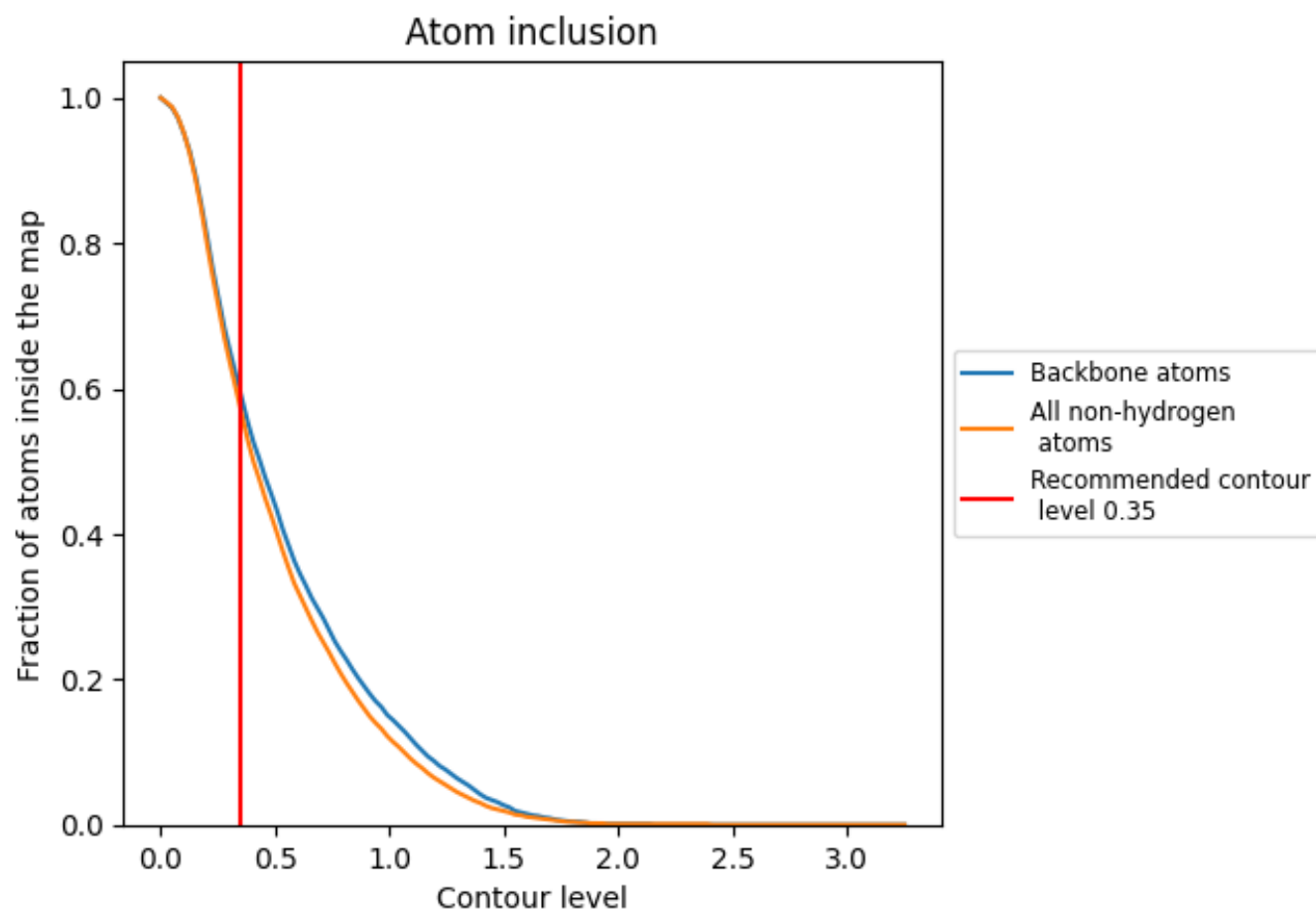
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.35).

9.4 Atom inclusion [i](#)



At the recommended contour level, 59% of all backbone atoms, 57% 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 (0.35) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5720	<div></div> 0.5280
A	<div></div> 0.7440	<div></div> 0.5770
B	<div></div> 0.7430	<div></div> 0.5750
C	<div></div> 0.4810	<div></div> 0.4640
D	<div></div> 0.4810	<div></div> 0.4600
E	<div></div> 0.3560	<div></div> 0.4780
F	<div></div> 0.3600	<div></div> 0.4780

1.0

0.0

<0.0