



Full wwPDB EM Validation Report ⓘ

Mar 23, 2026 – 10:47 PM UTC

PDB ID : 7EJC / pdb_00007ejc
EMDB ID : EMD-31158
Title : human RAD51 presynaptic complex
Authors : Zhao, L.Y.; Xu, J.F.; Wang, H.W.
Deposited on : 2021-04-02
Resolution : 2.97 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

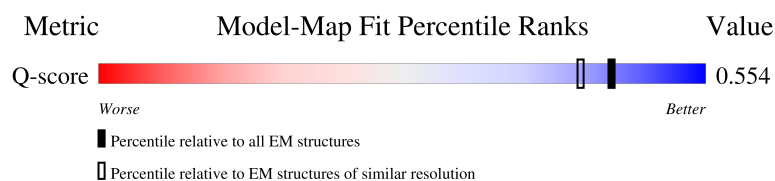
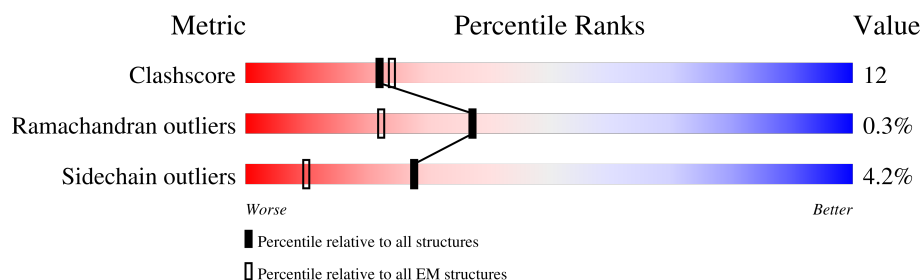
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.97 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	13205 (2.47 - 3.47)

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	G	9	
2	A	339	
2	B	339	
2	C	339	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7485 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 DNA chain called DNA (5'-D(P*TP*TP*TP*TP*TP*TP*TP*T)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
1	G	9	Total	C	N	O	P	0	0
			180	90	18	63	9		

- Molecule 2 is a protein called DNA repair protein RAD51 homolog 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	310	Total	C	N	O	S	0	0
			2375	1488	420	454	13		
2	C	310	Total	C	N	O	S	0	0
			2375	1488	420	454	13		
2	B	310	Total	C	N	O	S	0	0
			2375	1488	420	454	13		

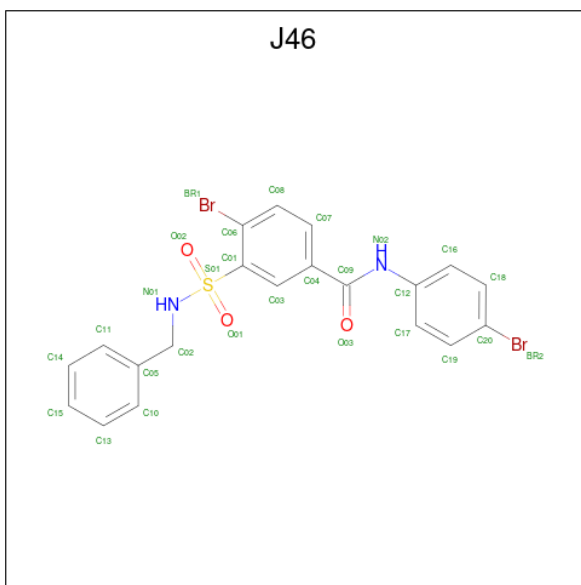
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	313	GLN	LYS	engineered mutation	UNP Q06609
C	313	GLN	LYS	engineered mutation	UNP Q06609
B	313	GLN	LYS	engineered mutation	UNP Q06609

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

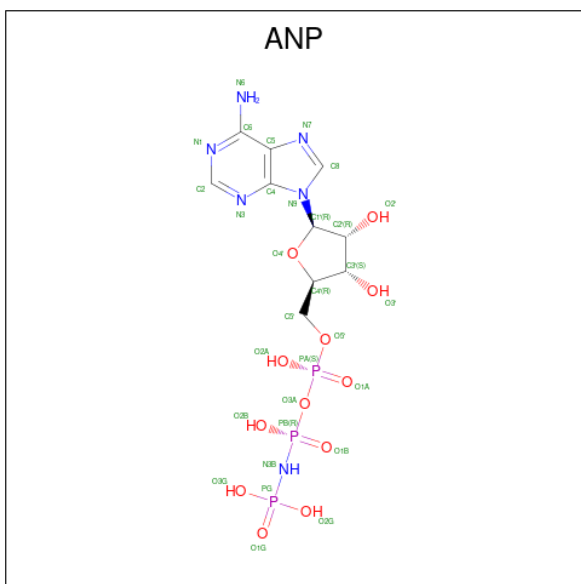
Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	Mg	0
			1	1	
3	C	1	Total	Mg	0
			1	1	
3	B	1	Total	Mg	0
			1	1	

- Molecule 4 is 4-bromanyl-N-(4-bromophenyl)-3-[(phenylmethyl)sulfamoyl]benzamide (CCD ID: J46) (formula: C₂₀H₁₆Br₂N₂O₃S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						AltConf
4	A	1	Total	Br	C	N	O	S	0
			28	2	20	2	3	1	
4	C	1	Total	Br	C	N	O	S	0
			28	2	20	2	3	1	
4	B	1	Total	Br	C	N	O	S	0
			28	2	20	2	3	1	

- Molecule 5 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).

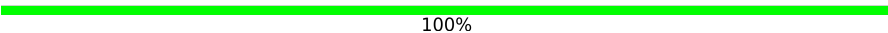


Mol	Chain	Residues	Atoms					AltConf
5	C	1	Total 31	C 10	N 6	O 12	P 3	0
5	C	1	Total 31	C 10	N 6	O 12	P 3	0
5	B	1	Total 31	C 10	N 6	O 12	P 3	0

3 Residue-property plots [i](#)

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

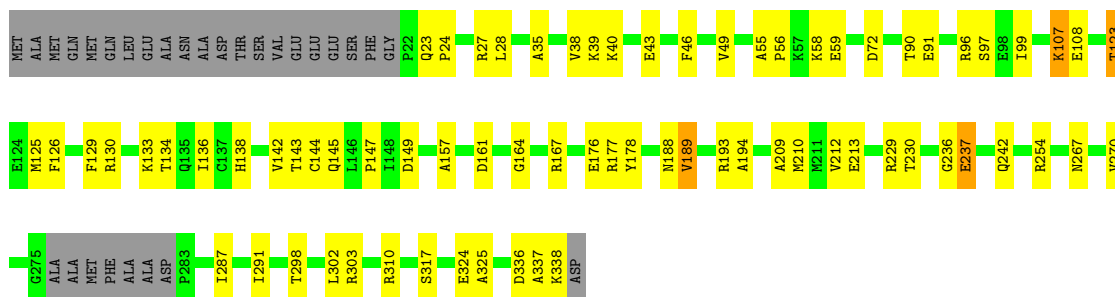
- Molecule 1: DNA (5'-D(P*TP*TP*TP*TP*TP*TP*TP*T)-3')

Chain G:  100%

There are no outlier residues recorded for this chain.

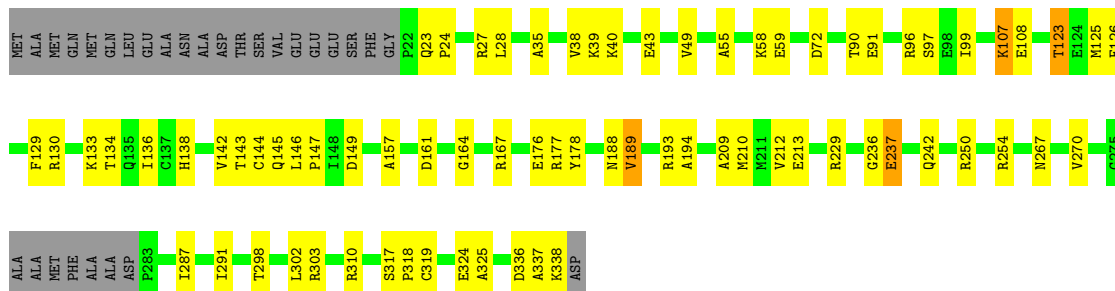
- Molecule 2: DNA repair protein RAD51 homolog 1

Chain A:  70% 20% 9%



- Molecule 2: DNA repair protein RAD51 homolog 1

Chain C:  70% 21% 9%



- Molecule 2: DNA repair protein RAD51 homolog 1

Chain B:  68% 22% 9%



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=56.7°, rise=15.8 Å, axial sym=C1	Depositor
Number of segments used	92403	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	48	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.353	Depositor
Minimum map value	-0.106	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.04	Depositor
Map size (Å)	354.0, 354.0, 354.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.885, 0.885, 0.885	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: J46, ANP, MG

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	G	0.32	0/197	0.54	0/302
2	A	0.28	0/2409	0.64	2/3247 (0.1%)
2	B	0.29	0/2409	0.65	2/3247 (0.1%)
2	C	0.28	0/2409	0.64	2/3247 (0.1%)
All	All	0.29	0/7424	0.64	6/10043 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	236	GLY	N-CA-C	-5.95	107.61	114.69
2	B	236	GLY	N-CA-C	-5.90	107.67	114.69
2	C	237	GLU	N-CA-C	-5.73	106.16	114.12
2	A	236	GLY	N-CA-C	-5.72	107.88	114.69
2	B	237	GLU	N-CA-C	-5.71	106.18	114.12
2	A	237	GLU	N-CA-C	-5.60	106.34	114.12

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	G	180	0	109	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	2375	0	2397	55	0
2	B	2375	0	2398	59	0
2	C	2375	0	2397	54	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	28	0	0	5	0
4	B	28	0	0	7	0
4	C	28	0	0	5	0
5	B	31	0	13	5	0
5	C	62	0	26	7	0
All	All	7485	0	7340	171	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 12.

All (171) 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)
2:C:58:LYS:HE3	4:C:404:J46:BR1	1.94	1.22
2:A:55:ALA:HB1	2:A:59:GLU:OE2	1.50	1.11
2:B:55:ALA:HB1	2:B:59:GLU:OE2	1.60	1.00
2:B:58:LYS:HE3	4:B:403:J46:BR1	2.22	0.94
2:B:134:THR:OG1	2:B:222:ASP:OD2	1.86	0.94
2:A:58:LYS:HE3	4:A:602:J46:BR1	2.26	0.91
2:C:134:THR:OG1	5:C:401:ANP:O2B	1.90	0.88
2:A:55:ALA:CB	2:A:59:GLU:OE2	2.24	0.85
2:A:134:THR:OG1	5:B:401:ANP:O2B	1.95	0.84
2:C:55:ALA:HB1	2:C:59:GLU:OE2	1.78	0.84
2:C:193:ARG:NH1	2:C:194:ALA:O	2.12	0.83
2:B:193:ARG:NH1	2:B:194:ALA:O	2.12	0.82
2:A:193:ARG:NH1	2:A:194:ALA:O	2.12	0.82
2:A:46:PHE:CZ	2:A:59:GLU:OE1	2.32	0.82
2:C:58:LYS:CE	4:C:404:J46:BR1	2.82	0.80
2:B:55:ALA:CB	2:B:59:GLU:OE2	2.32	0.77
2:A:130:ARG:O	2:A:310:ARG:NH1	2.19	0.76
2:B:130:ARG:O	2:B:310:ARG:NH1	2.19	0.76
2:C:130:ARG:O	2:C:310:ARG:NH1	2.19	0.76
2:A:242:GLN:HB3	2:A:291:ILE:HD12	1.72	0.72
2:B:242:GLN:HB3	2:B:291:ILE:HD12	1.72	0.72
2:C:242:GLN:HB3	2:C:291:ILE:HD12	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:134:THR:CB	5:B:401:ANP:O2B	2.41	0.69
2:C:134:THR:CB	5:C:401:ANP:O2B	2.41	0.68
2:B:336:ASP:H	2:B:338:LYS:HD2	1.59	0.67
2:A:336:ASP:H	2:A:338:LYS:HD2	1.59	0.67
2:A:91:GLU:N	2:A:91:GLU:OE1	2.28	0.67
2:C:336:ASP:H	2:C:338:LYS:HD2	1.59	0.66
2:C:91:GLU:N	2:C:91:GLU:OE1	2.28	0.66
2:C:270:VAL:HG13	2:C:287:ILE:HG22	1.78	0.65
2:B:91:GLU:OE1	2:B:91:GLU:N	2.28	0.65
2:A:46:PHE:HZ	2:A:59:GLU:OE1	1.79	0.65
2:A:270:VAL:HG13	2:A:287:ILE:HG22	1.78	0.64
2:C:138:HIS:CE1	2:C:167:ARG:H	2.16	0.64
2:C:254:ARG:HD2	4:C:404:J46:C18	2.28	0.64
2:B:270:VAL:HG13	2:B:287:ILE:HG22	1.78	0.64
2:A:138:HIS:CE1	2:A:167:ARG:H	2.16	0.63
2:B:138:HIS:CE1	2:B:167:ARG:H	2.16	0.63
2:B:298:THR:HG23	2:B:317:SER:HB2	1.82	0.62
2:A:298:THR:HG23	2:A:317:SER:HB2	1.82	0.61
2:C:298:THR:HG23	2:C:317:SER:HB2	1.82	0.61
2:B:303:ARG:NH2	2:B:324:GLU:OE1	2.33	0.61
2:A:303:ARG:NH2	2:A:324:GLU:OE1	2.33	0.61
2:B:223:SER:H	2:B:266:THR:HG1	1.46	0.61
2:C:303:ARG:NH2	2:C:324:GLU:OE1	2.33	0.61
2:C:250:ARG:HH11	2:B:230:THR:HG23	1.65	0.60
2:C:337:ALA:N	2:C:338:LYS:HB2	2.17	0.60
2:B:337:ALA:N	2:B:338:LYS:HB2	2.17	0.60
2:A:134:THR:HB	5:B:401:ANP:O2B	2.03	0.59
2:A:337:ALA:N	2:A:338:LYS:HB2	2.17	0.59
2:A:230:THR:HG23	2:B:250:ARG:HH11	1.67	0.58
2:C:55:ALA:CB	2:C:59:GLU:OE2	2.51	0.58
2:A:123:THR:HG23	2:A:298:THR:HB	1.86	0.58
2:B:123:THR:HG23	2:B:298:THR:HB	1.86	0.58
2:B:254:ARG:HD2	4:B:403:J46:C18	2.35	0.57
2:C:24:PRO:HD2	2:C:27:ARG:HE	1.69	0.57
2:A:24:PRO:HD2	2:A:27:ARG:HE	1.69	0.57
2:C:123:THR:HG23	2:C:298:THR:HB	1.86	0.57
2:B:24:PRO:HD2	2:B:27:ARG:HE	1.69	0.57
2:A:108:GLU:HG3	2:A:325:ALA:HB1	1.87	0.56
2:C:108:GLU:HG3	2:C:325:ALA:HB1	1.86	0.56
2:A:129:PHE:O	2:A:130:ARG:HG2	2.05	0.56
2:C:319:CYS:O	5:C:403:ANP:N6	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:108:GLU:HG3	2:B:325:ALA:HB1	1.86	0.56
2:B:129:PHE:O	2:B:130:ARG:HG2	2.05	0.56
2:A:46:PHE:HZ	2:A:59:GLU:CD	2.13	0.56
2:B:319:CYS:O	5:B:401:ANP:N6	2.39	0.56
2:C:23:GLN:HB3	2:C:27:ARG:HD2	1.89	0.55
2:C:134:THR:HB	5:C:401:ANP:O2B	2.06	0.55
2:C:129:PHE:O	2:C:130:ARG:HG2	2.05	0.55
2:A:23:GLN:HB3	2:A:27:ARG:HD2	1.89	0.55
2:C:229:ARG:HH21	2:C:287:ILE:HG21	1.72	0.55
2:A:229:ARG:HH21	2:A:287:ILE:HG21	1.72	0.55
2:B:23:GLN:HB3	2:B:27:ARG:HD2	1.89	0.54
5:C:403:ANP:O3G	5:C:403:ANP:O2B	2.26	0.54
5:C:401:ANP:O2B	5:C:401:ANP:O3G	2.26	0.54
5:B:401:ANP:O2B	5:B:401:ANP:O3G	2.26	0.54
2:C:35:ALA:O	2:C:39:LYS:HD3	2.09	0.53
2:A:35:ALA:O	2:A:39:LYS:HD3	2.09	0.53
4:A:602:J46:BR1	4:A:602:J46:O01	2.82	0.53
2:B:229:ARG:HH21	2:B:287:ILE:HG21	1.72	0.53
2:A:157:ALA:O	2:A:189:VAL:HA	2.09	0.53
2:B:254:ARG:HD2	4:B:403:J46:C16	2.38	0.53
2:C:336:ASP:N	2:C:338:LYS:HD2	2.24	0.52
2:B:35:ALA:O	2:B:39:LYS:HD3	2.09	0.52
4:B:403:J46:BR1	4:B:403:J46:O01	2.82	0.52
2:C:157:ALA:O	2:C:189:VAL:HA	2.09	0.52
4:C:404:J46:BR1	4:C:404:J46:O01	2.82	0.51
2:B:336:ASP:N	2:B:338:LYS:HD2	2.24	0.51
2:B:157:ALA:O	2:B:189:VAL:HA	2.09	0.51
2:A:336:ASP:N	2:A:338:LYS:HD2	2.24	0.51
2:A:161:ASP:OD1	2:A:164:GLY:N	2.22	0.50
2:A:134:THR:HG22	2:A:134:THR:O	2.11	0.50
2:B:337:ALA:H	2:B:338:LYS:HB2	1.77	0.49
2:A:107:LYS:HA	2:A:107:LYS:HE2	1.95	0.49
2:A:337:ALA:H	2:A:338:LYS:HB2	1.77	0.49
2:B:107:LYS:HA	2:B:107:LYS:HE2	1.95	0.49
2:C:107:LYS:HE2	2:C:107:LYS:HA	1.95	0.49
2:B:134:THR:O	2:B:134:THR:HG23	2.11	0.49
2:C:138:HIS:HE1	2:C:167:ARG:H	1.61	0.48
2:C:147:PRO:HB2	2:C:149:ASP:OD1	2.14	0.48
2:C:337:ALA:H	2:C:338:LYS:HB2	1.77	0.48
2:A:126:PHE:HB3	2:A:267:ASN:HB3	1.96	0.48
2:B:126:PHE:HB3	2:B:267:ASN:HB3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:147:PRO:HB2	2:B:149:ASP:OD1	2.14	0.48
2:C:126:PHE:HB3	2:C:267:ASN:HB3	1.96	0.47
2:B:59:GLU:CD	2:B:254:ARG:HH22	2.21	0.47
2:C:28:LEU:HD22	2:C:49:VAL:HG12	1.96	0.47
2:A:46:PHE:CZ	2:A:59:GLU:CD	2.90	0.47
2:A:147:PRO:HB2	2:A:149:ASP:OD1	2.14	0.47
2:B:28:LEU:HD22	2:B:49:VAL:HG12	1.96	0.47
2:A:59:GLU:HG3	4:A:602:J46:C03	2.43	0.47
2:A:28:LEU:HD22	2:A:49:VAL:HG12	1.96	0.47
2:C:134:THR:O	2:C:134:THR:HG22	2.13	0.47
2:A:138:HIS:HE1	2:A:167:ARG:H	1.61	0.46
2:B:138:HIS:HE1	2:B:167:ARG:H	1.61	0.46
2:B:58:LYS:CE	4:B:403:J46:BR1	3.08	0.45
2:B:59:GLU:HG3	4:B:403:J46:C01	2.47	0.45
2:C:133:LYS:HB3	2:C:302:LEU:HD12	1.99	0.45
2:A:254:ARG:HD2	4:A:602:J46:C18	2.46	0.45
2:A:254:ARG:HD2	4:A:602:J46:C16	2.47	0.45
2:A:40:LYS:HA	2:A:43:GLU:OE1	2.18	0.44
2:C:145:GLN:HE21	2:C:188:ASN:HB3	1.83	0.44
2:C:40:LYS:HA	2:C:43:GLU:OE1	2.18	0.44
2:B:133:LYS:HB3	2:B:302:LEU:HD12	1.99	0.44
2:B:134:THR:OG1	2:B:222:ASP:CG	2.57	0.44
2:C:318:PRO:HA	5:C:403:ANP:H8	1.99	0.44
2:B:40:LYS:HA	2:B:43:GLU:OE1	2.18	0.44
2:B:145:GLN:HE21	2:B:188:ASN:HB3	1.83	0.44
2:A:125:MET:HE1	2:A:136:ILE:HB	2.00	0.44
2:C:125:MET:HE1	2:C:136:ILE:HB	2.00	0.44
2:A:133:LYS:HB3	2:A:302:LEU:HD12	1.99	0.43
2:C:96:ARG:O	2:C:99:ILE:HG12	2.18	0.43
2:B:56:PRO:HG2	2:B:59:GLU:HB2	2.00	0.43
2:B:96:ARG:O	2:B:99:ILE:HG12	2.18	0.43
2:B:125:MET:HE1	2:B:136:ILE:HB	2.00	0.43
2:B:188:ASN:O	2:B:189:VAL:C	2.61	0.43
2:A:96:ARG:O	2:A:99:ILE:HG12	2.18	0.43
2:A:270:VAL:HG13	2:A:287:ILE:CG2	2.46	0.43
2:C:161:ASP:OD1	2:C:164:GLY:N	2.22	0.43
2:C:188:ASN:O	2:C:189:VAL:C	2.61	0.43
2:C:270:VAL:HG13	2:C:287:ILE:CG2	2.47	0.43
2:B:270:VAL:HG13	2:B:287:ILE:CG2	2.46	0.43
2:B:59:GLU:HG3	4:B:403:J46:C03	2.48	0.43
2:A:188:ASN:O	2:A:189:VAL:C	2.61	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:145:GLN:HE21	2:A:188:ASN:HB3	1.83	0.43
2:C:254:ARG:HD2	4:C:404:J46:C16	2.48	0.43
2:B:204:LEU:HD23	2:B:204:LEU:HA	1.85	0.42
2:B:209:ALA:O	2:B:212:VAL:HG12	2.19	0.42
2:A:209:ALA:O	2:A:212:VAL:HG12	2.19	0.42
2:C:209:ALA:O	2:C:212:VAL:HG12	2.19	0.42
2:A:144:CYS:SG	2:A:157:ALA:HB2	2.60	0.42
2:B:177:ARG:HH21	2:B:178:TYR:HE2	1.67	0.42
2:B:210:MET:O	2:B:213:GLU:HG3	2.20	0.42
2:B:146:LEU:HD23	2:B:146:LEU:HA	1.92	0.42
2:C:146:LEU:HD23	2:C:146:LEU:HA	1.92	0.41
2:C:177:ARG:HH21	2:C:178:TYR:HE2	1.67	0.41
2:B:144:CYS:SG	2:B:157:ALA:HB2	2.60	0.41
2:C:250:ARG:NH1	2:B:230:THR:HG23	2.35	0.41
2:A:177:ARG:HH21	2:A:178:TYR:HE2	1.67	0.41
2:C:210:MET:O	2:C:213:GLU:HG3	2.20	0.41
2:B:161:ASP:OD1	2:B:164:GLY:N	2.22	0.41
2:A:210:MET:O	2:A:213:GLU:HG3	2.20	0.41
2:B:28:LEU:HD23	2:B:28:LEU:H	1.85	0.41
2:C:144:CYS:SG	2:C:157:ALA:HB2	2.60	0.41
2:B:109:LEU:HD23	2:B:109:LEU:HA	1.90	0.41
2:A:28:LEU:HD23	2:A:28:LEU:H	1.86	0.41
2:A:56:PRO:HG2	2:A:59:GLU:HB2	2.02	0.41
2:C:28:LEU:HD23	2:C:28:LEU:H	1.86	0.41
2:A:157:ALA:HB3	2:A:189:VAL:HG22	2.03	0.40
2:C:157:ALA:HB3	2:C:189:VAL:HG22	2.03	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	
2	A	306/339 (90%)	289 (94%)	16 (5%)	1 (0%)	36	67
2	B	306/339 (90%)	289 (94%)	16 (5%)	1 (0%)	36	67
2	C	306/339 (90%)	289 (94%)	16 (5%)	1 (0%)	36	67
All	All	918/1017 (90%)	867 (94%)	48 (5%)	3 (0%)	37	67

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	189	VAL
2	C	189	VAL
2	B	189	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	248/269 (92%)	238 (96%)	10 (4%)	28	60
2	B	248/269 (92%)	237 (96%)	11 (4%)	25	58
2	C	248/269 (92%)	238 (96%)	10 (4%)	28	60
All	All	744/807 (92%)	713 (96%)	31 (4%)	28	59

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

Mol	Chain	Res	Type
2	A	38	VAL
2	A	72	ASP
2	A	90	THR
2	A	97	SER
2	A	107	LYS
2	A	123	THR
2	A	142	VAL
2	A	143	THR
2	A	176	GLU
2	A	237	GLU

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Mol	Chain	Res	Type
2	C	38	VAL
2	C	72	ASP
2	C	90	THR
2	C	97	SER
2	C	107	LYS
2	C	123	THR
2	C	142	VAL
2	C	143	THR
2	C	176	GLU
2	C	237	GLU
2	B	38	VAL
2	B	72	ASP
2	B	90	THR
2	B	97	SER
2	B	107	LYS
2	B	123	THR
2	B	134	THR
2	B	142	VAL
2	B	143	THR
2	B	176	GLU
2	B	237	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	34	ASN
2	A	114	GLN
2	A	138	HIS
2	C	34	ASN
2	C	93	HIS
2	C	114	GLN
2	C	138	HIS
2	C	290	ASN
2	B	34	ASN
2	B	93	HIS
2	B	114	GLN
2	B	138	HIS
2	B	290	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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$
4	J46	A	602	-	30,30,30	4.58	21 (70%)	41,42,42	3.52	9 (21%)
4	J46	B	403	-	30,30,30	4.57	21 (70%)	41,42,42	3.52	9 (21%)
5	ANP	C	403	3	33,33,33	0.87	3 (9%)	45,52,52	0.69	1 (2%)
5	ANP	B	401	3	33,33,33	0.88	3 (9%)	45,52,52	0.68	1 (2%)
5	ANP	C	401	3	33,33,33	0.87	3 (9%)	45,52,52	0.68	1 (2%)
4	J46	C	404	-	30,30,30	4.57	21 (70%)	41,42,42	3.53	9 (21%)

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
4	J46	A	602	-	-	1/20/20/20	0/3/3/3
4	J46	B	403	-	-	1/20/20/20	0/3/3/3
5	ANP	C	403	3	-	7/18/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ANP	B	401	3	-	7/18/38/38	0/3/3/3
5	ANP	C	401	3	-	7/18/38/38	0/3/3/3
4	J46	C	404	-	-	1/20/20/20	0/3/3/3

All (72) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	602	J46	C01-C06	-11.17	1.33	1.39
4	B	403	J46	C01-C06	-11.03	1.33	1.39
4	C	404	J46	C01-C06	-11.00	1.33	1.39
4	C	404	J46	C07-C04	9.19	1.53	1.39
4	A	602	J46	C07-C04	9.16	1.53	1.39
4	B	403	J46	C07-C04	9.15	1.53	1.39
4	A	602	J46	C03-C01	7.13	1.50	1.39
4	C	404	J46	C03-C01	7.11	1.50	1.39
4	B	403	J46	C03-C01	7.11	1.50	1.39
4	B	403	J46	C18-C20	6.63	1.51	1.38
4	A	602	J46	C16-C12	6.62	1.50	1.39
4	A	602	J46	C18-C20	6.61	1.51	1.38
4	C	404	J46	C16-C12	6.60	1.50	1.39
4	C	404	J46	C18-C20	6.60	1.51	1.38
4	B	403	J46	C16-C12	6.58	1.50	1.39
4	B	403	J46	C08-C06	6.50	1.51	1.38
4	A	602	J46	C08-C06	6.49	1.51	1.38
4	C	404	J46	C08-C06	6.47	1.51	1.38
4	B	403	J46	C14-C15	6.21	1.51	1.38
4	C	404	J46	C14-C15	6.20	1.51	1.38
4	A	602	J46	C14-C15	6.20	1.51	1.38
4	A	602	J46	C13-C10	5.52	1.48	1.38
4	C	404	J46	C13-C10	5.50	1.48	1.38
4	B	403	J46	C13-C10	5.50	1.48	1.38
4	B	403	J46	C19-C17	5.47	1.47	1.38
4	A	602	J46	C19-C17	5.46	1.47	1.38
4	C	404	J46	C19-C17	5.46	1.47	1.38
4	C	404	J46	C18-C16	-4.86	1.31	1.38
4	A	602	J46	C18-C16	-4.85	1.31	1.38
4	B	403	J46	C18-C16	-4.84	1.31	1.38
4	B	403	J46	C03-C04	-4.65	1.32	1.39
4	C	404	J46	C03-C04	-4.62	1.32	1.39
4	A	602	J46	C03-C04	-4.61	1.32	1.39
4	C	404	J46	C11-C05	4.20	1.47	1.38
4	B	403	J46	C11-C05	4.20	1.47	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	602	J46	C11-C05	4.20	1.47	1.38
4	B	403	J46	S01-N01	4.17	1.67	1.61
4	C	404	J46	S01-N01	4.13	1.67	1.61
4	A	602	J46	S01-N01	4.11	1.67	1.61
4	C	404	J46	C17-C12	-3.34	1.33	1.39
4	B	403	J46	C17-C12	-3.34	1.33	1.39
4	A	602	J46	C17-C12	-3.33	1.33	1.39
4	B	403	J46	C19-C20	-3.28	1.31	1.38
4	A	602	J46	C19-C20	-3.25	1.31	1.38
4	C	404	J46	C19-C20	-3.25	1.31	1.38
4	C	404	J46	C09-N02	3.00	1.44	1.35
4	A	602	J46	C09-N02	2.99	1.44	1.35
4	B	403	J46	C09-N02	2.97	1.44	1.35
4	C	404	J46	C14-C11	-2.73	1.34	1.38
4	A	602	J46	C14-C11	-2.72	1.34	1.38
4	B	403	J46	C14-C11	-2.68	1.34	1.38
4	B	403	J46	C15-C13	-2.68	1.32	1.38
4	A	602	J46	C15-C13	-2.64	1.32	1.38
4	C	404	J46	C15-C13	-2.64	1.32	1.38
5	B	401	ANP	PB-O3A	-2.42	1.56	1.59
5	C	401	ANP	PB-O3A	-2.39	1.56	1.59
4	B	403	J46	C08-C07	-2.35	1.35	1.38
5	C	403	ANP	PB-O3A	-2.34	1.56	1.59
4	C	404	J46	C08-C07	-2.32	1.35	1.38
4	A	602	J46	C08-C07	-2.27	1.35	1.38
5	B	401	ANP	PG-N3B	2.14	1.68	1.63
5	C	401	ANP	PG-N3B	2.12	1.68	1.63
5	C	403	ANP	PG-N3B	2.11	1.68	1.63
4	A	602	J46	BR1-C06	-2.09	1.85	1.89
4	C	404	J46	BR1-C06	-2.07	1.85	1.89
4	B	403	J46	BR1-C06	-2.07	1.85	1.89
4	B	403	J46	C02-C05	2.06	1.56	1.51
5	C	401	ANP	PG-O1G	2.05	1.49	1.46
5	B	401	ANP	PG-O1G	2.04	1.49	1.46
5	C	403	ANP	PG-O1G	2.04	1.49	1.46
4	C	404	J46	C02-C05	2.04	1.56	1.51
4	A	602	J46	C02-C05	2.02	1.55	1.51

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	602	J46	O02-S01-O01	-15.95	100.15	119.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	403	J46	O02-S01-O01	-15.93	100.18	119.52
4	C	404	J46	O02-S01-O01	-15.92	100.18	119.52
4	C	404	J46	BR1-C06-C01	-12.56	115.16	122.71
4	B	403	J46	BR1-C06-C01	-12.51	115.19	122.71
4	A	602	J46	BR1-C06-C01	-12.50	115.20	122.71
4	B	403	J46	O01-S01-C01	4.37	114.90	107.68
4	C	404	J46	O01-S01-C01	4.37	114.89	107.68
4	A	602	J46	O01-S01-C01	4.34	114.84	107.68
4	A	602	J46	C08-C06-C01	3.44	122.04	120.11
4	C	404	J46	C08-C06-C01	3.43	122.03	120.11
4	B	403	J46	C08-C06-C01	3.42	122.02	120.11
4	A	602	J46	O01-S01-N01	3.02	111.73	107.03
4	C	404	J46	O01-S01-N01	3.02	111.73	107.03
4	B	403	J46	O01-S01-N01	3.01	111.72	107.03
4	A	602	J46	O02-S01-N01	2.87	111.50	107.03
4	C	404	J46	O02-S01-N01	2.85	111.46	107.03
4	B	403	J46	O02-S01-N01	2.84	111.45	107.03
4	C	404	J46	BR1-C06-C08	2.61	122.80	117.82
4	B	403	J46	BR1-C06-C08	2.60	122.78	117.82
4	A	602	J46	BR1-C06-C08	2.59	122.76	117.82
5	C	403	ANP	O1B-PB-N3B	-2.54	108.03	111.77
4	C	404	J46	C16-C12-C17	2.53	122.40	119.04
5	B	401	ANP	O1B-PB-N3B	-2.53	108.05	111.77
4	B	403	J46	C16-C12-C17	2.52	122.39	119.04
5	C	401	ANP	O1B-PB-N3B	-2.52	108.06	111.77
4	A	602	J46	C16-C12-C17	2.50	122.37	119.04
4	A	602	J46	C08-C07-C04	-2.15	118.50	120.80
4	C	404	J46	C08-C07-C04	-2.14	118.51	120.80
4	B	403	J46	C08-C07-C04	-2.12	118.53	120.80

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	401	ANP	PB-N3B-PG-O1G
5	C	401	ANP	PG-N3B-PB-O3A
5	C	401	ANP	C5'-O5'-PA-O1A
5	C	401	ANP	C5'-O5'-PA-O3A
5	C	403	ANP	PB-N3B-PG-O1G
5	C	403	ANP	PG-N3B-PB-O3A
5	C	403	ANP	C5'-O5'-PA-O1A
5	C	403	ANP	C5'-O5'-PA-O3A

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Mol	Chain	Res	Type	Atoms
5	B	401	ANP	PB-N3B-PG-O1G
5	B	401	ANP	PG-N3B-PB-O3A
5	B	401	ANP	C5'-O5'-PA-O1A
5	B	401	ANP	C5'-O5'-PA-O3A
5	C	401	ANP	C3'-C4'-C5'-O5'
5	C	403	ANP	C3'-C4'-C5'-O5'
5	B	401	ANP	C3'-C4'-C5'-O5'
5	C	401	ANP	C5'-O5'-PA-O2A
5	C	403	ANP	C5'-O5'-PA-O2A
5	B	401	ANP	C5'-O5'-PA-O2A
5	C	403	ANP	O4'-C4'-C5'-O5'
5	B	401	ANP	O4'-C4'-C5'-O5'
5	C	401	ANP	O4'-C4'-C5'-O5'
4	A	602	J46	C02-N01-S01-O01
4	C	404	J46	C02-N01-S01-O01
4	B	403	J46	C02-N01-S01-O01

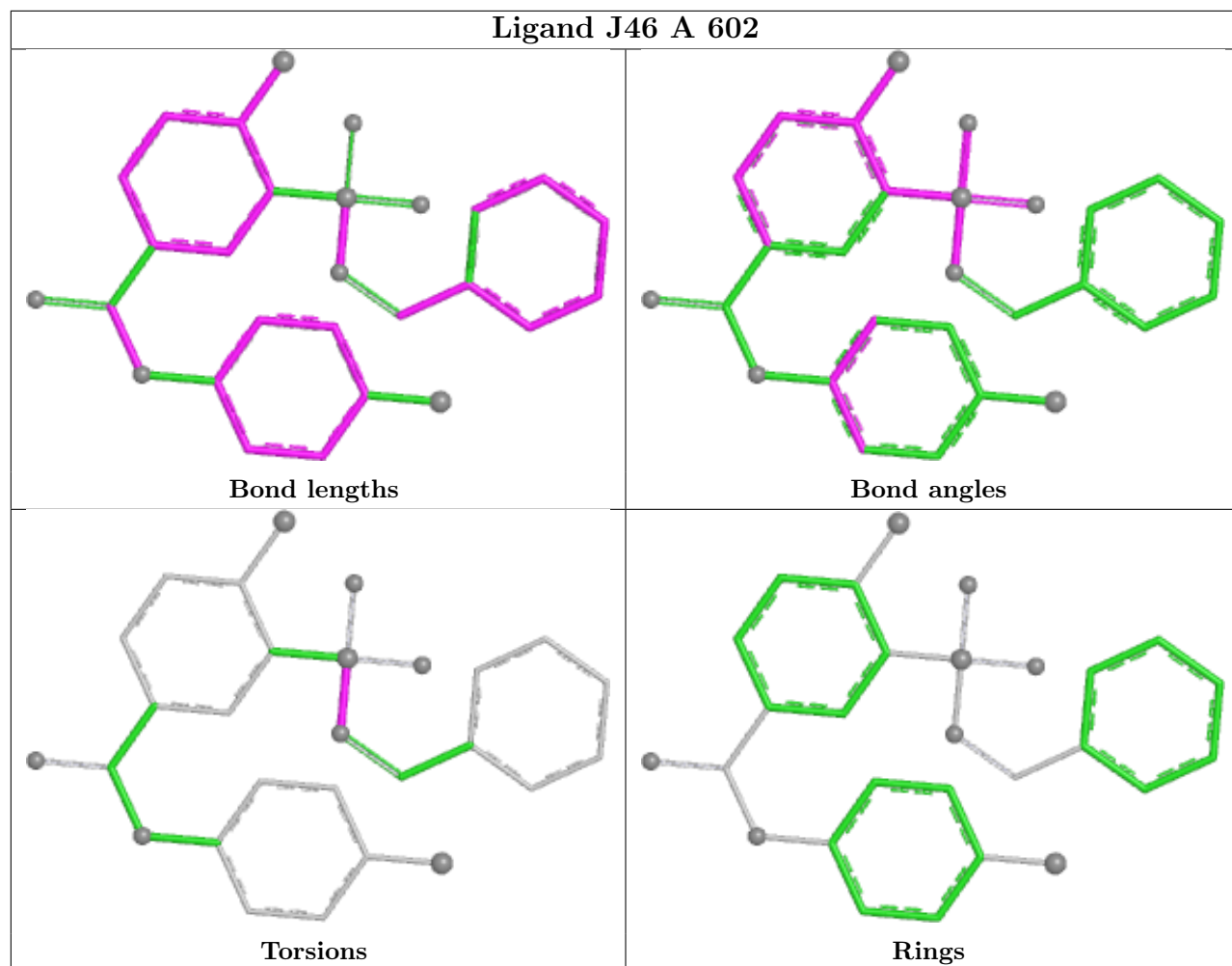
There are no ring outliers.

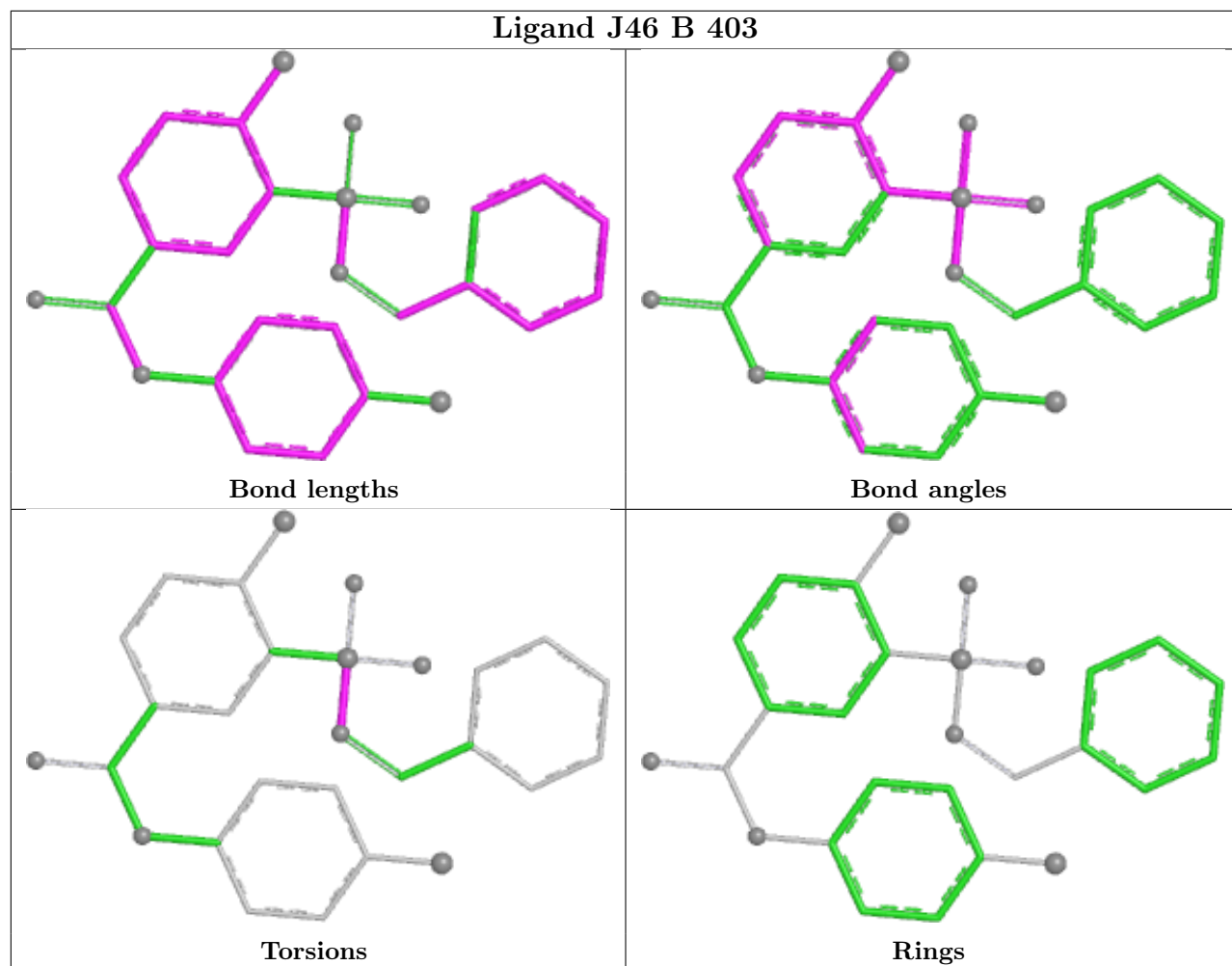
6 monomers are involved in 29 short contacts:

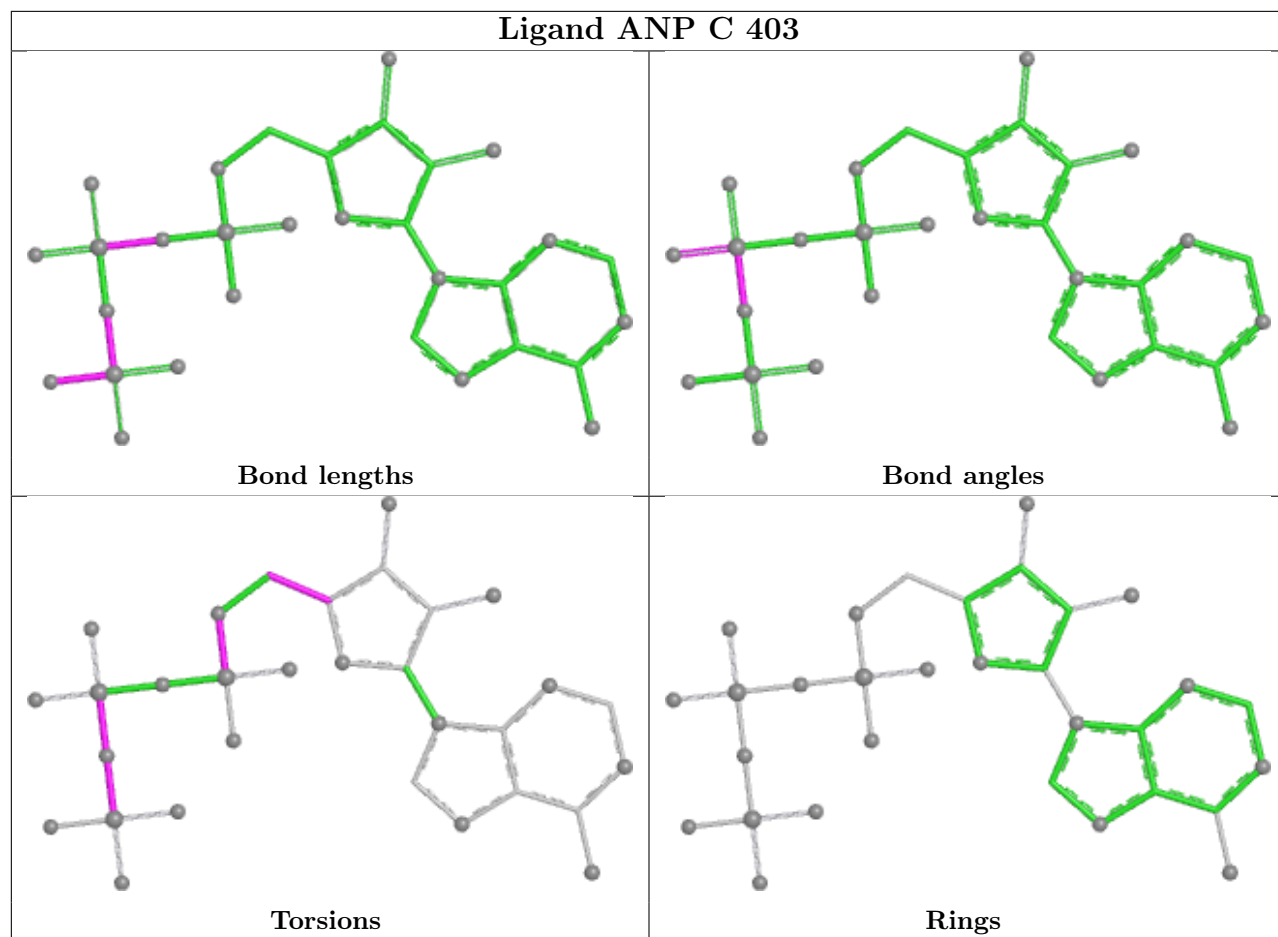
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	602	J46	5	0
4	B	403	J46	7	0
5	C	403	ANP	3	0
5	B	401	ANP	5	0
5	C	401	ANP	4	0
4	C	404	J46	5	0

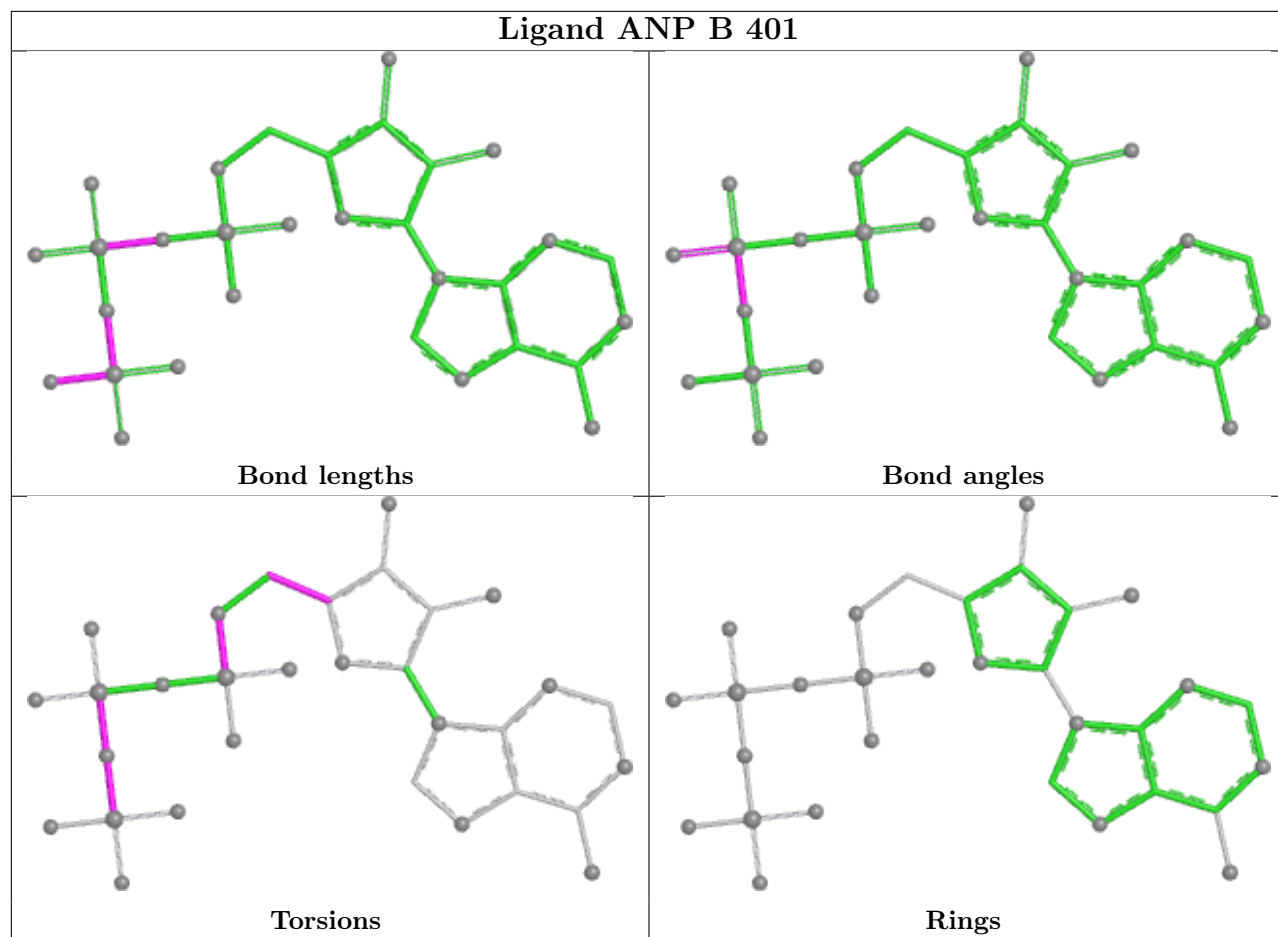
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

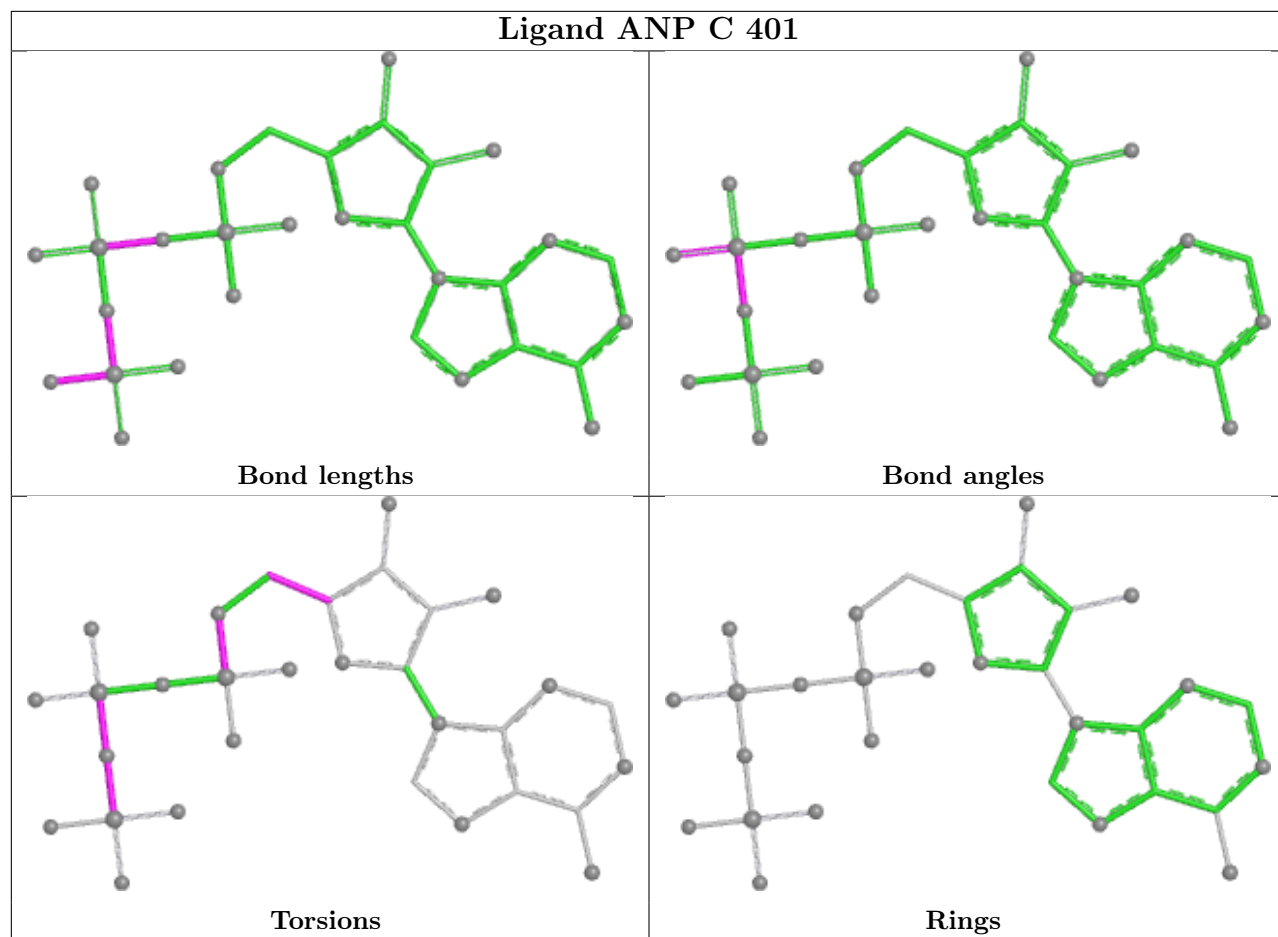
Ligand J46 A 602

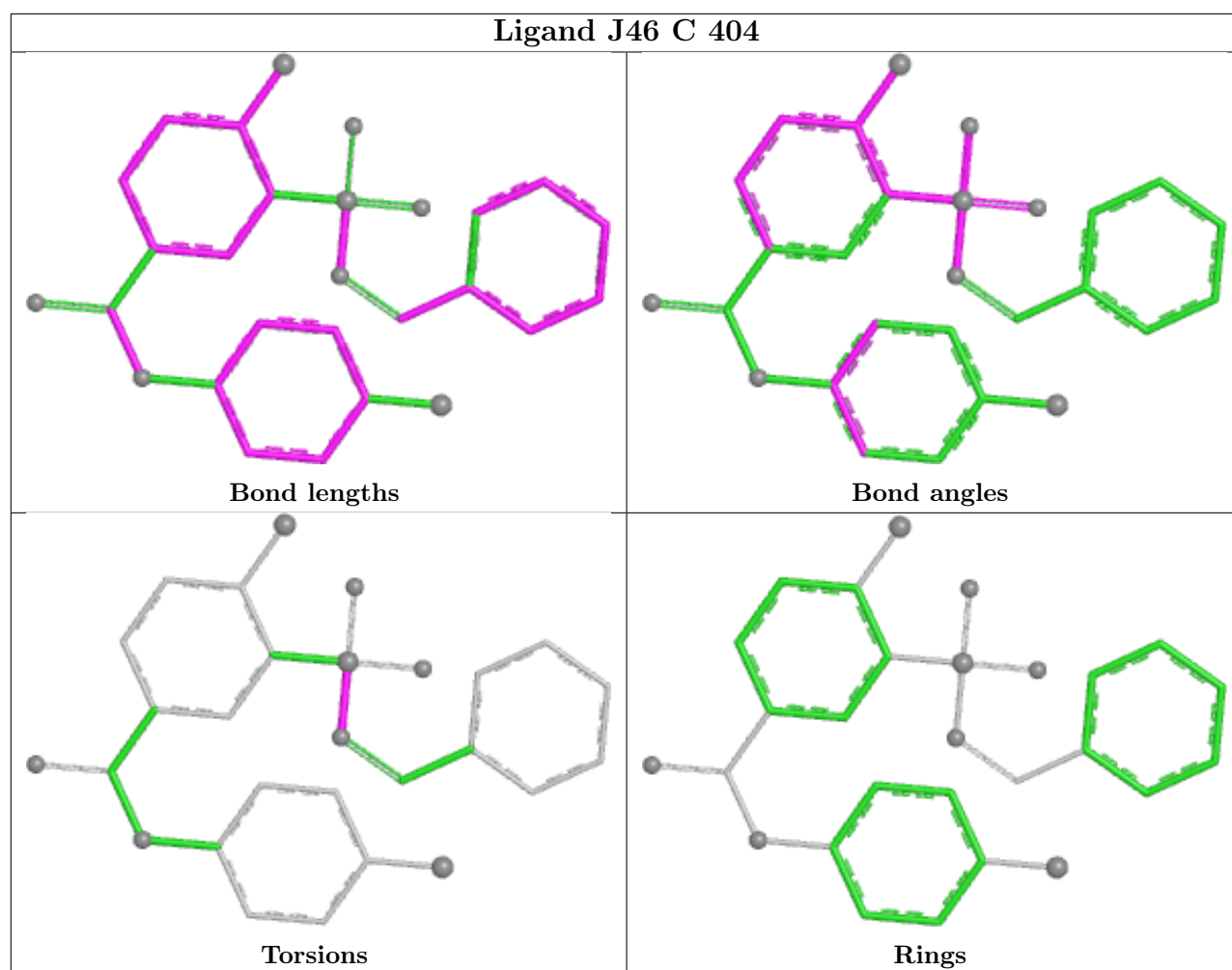












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

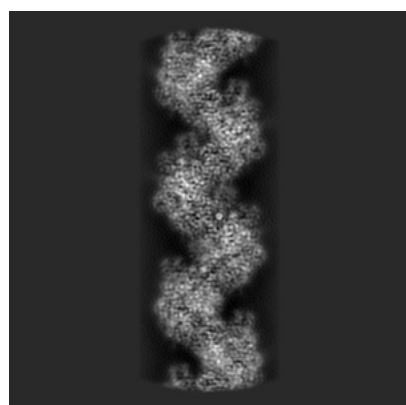
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31158. These allow visual inspection of the internal detail of the map and identification of artifacts.

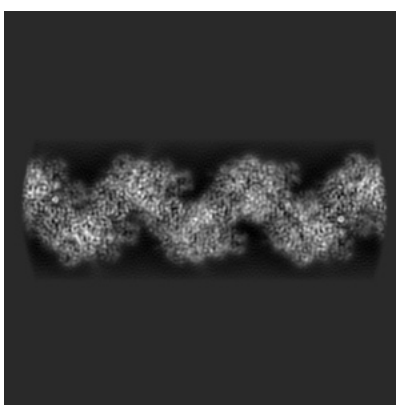
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

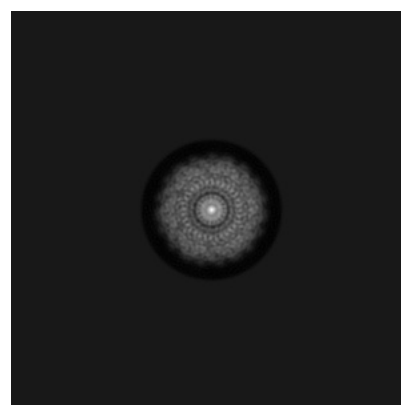
6.1.1 Primary map



X



Y

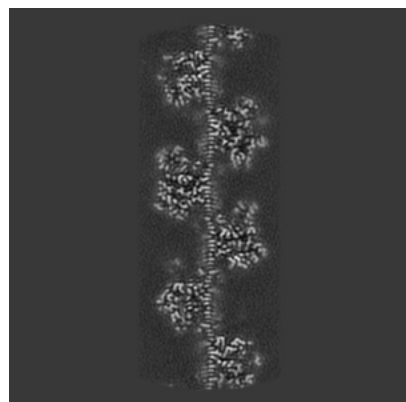


Z

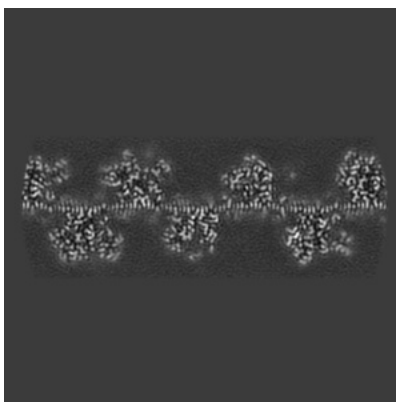
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

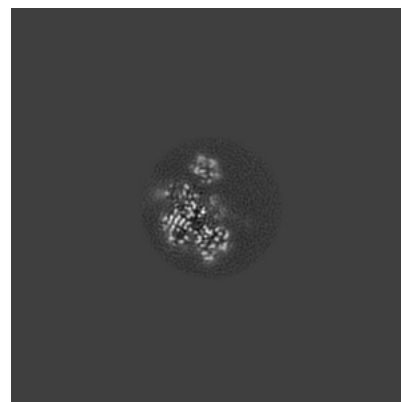
6.2.1 Primary map



X Index: 200



Y Index: 200

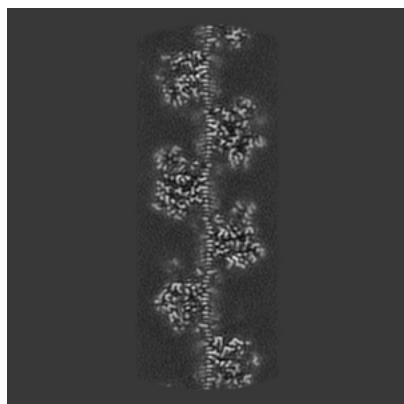


Z Index: 200

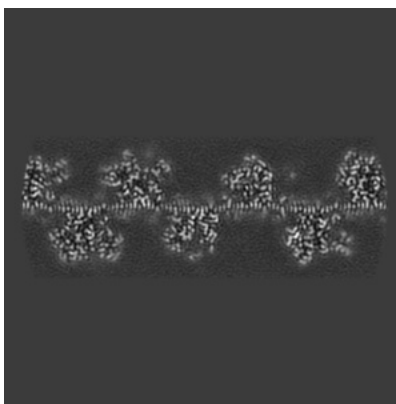
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

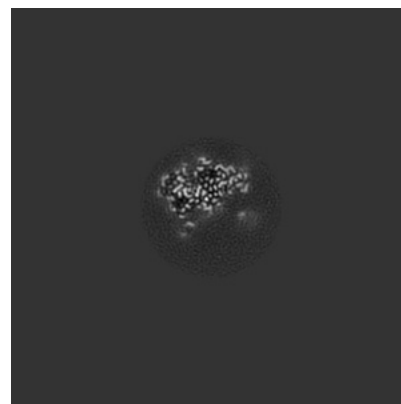
6.3.1 Primary map



X Index: 200



Y Index: 200

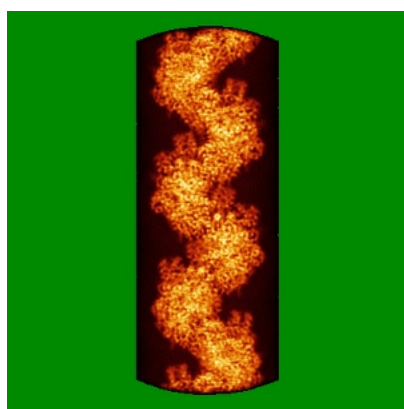


Z Index: 171

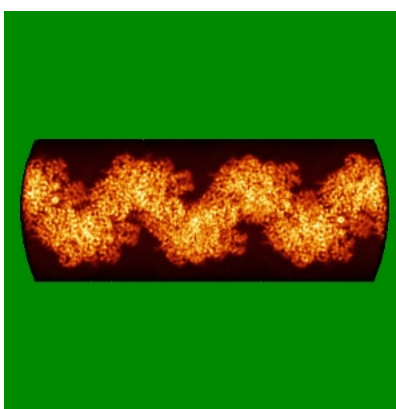
The images above show the largest variance slices of the map in three orthogonal directions.

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

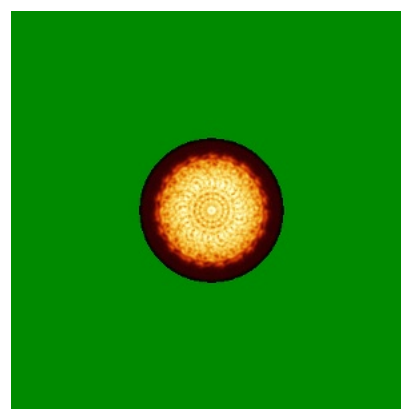
6.4.1 Primary map



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](#)

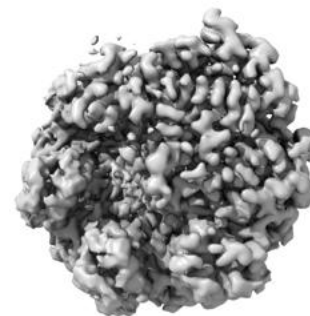
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.04. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

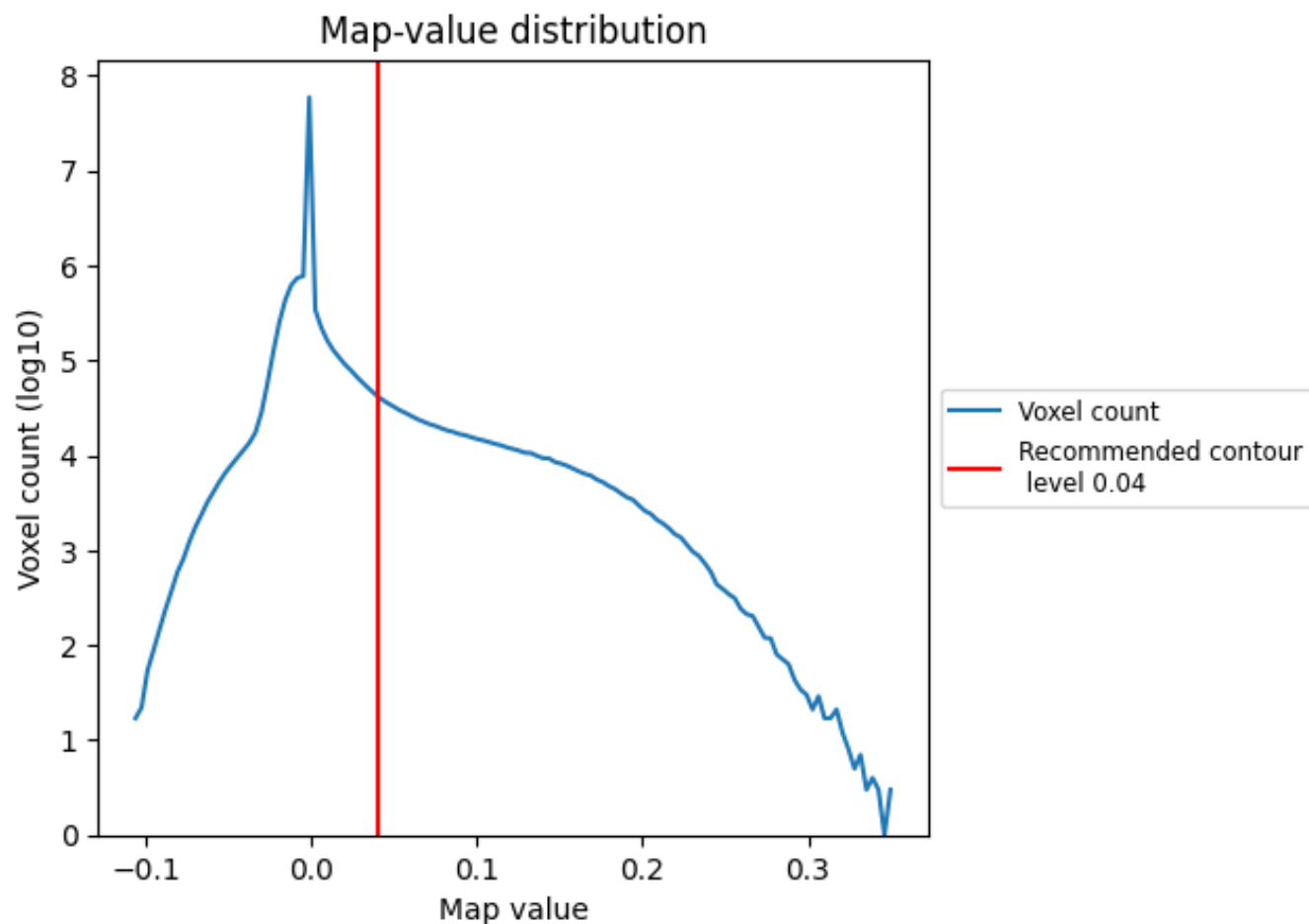
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

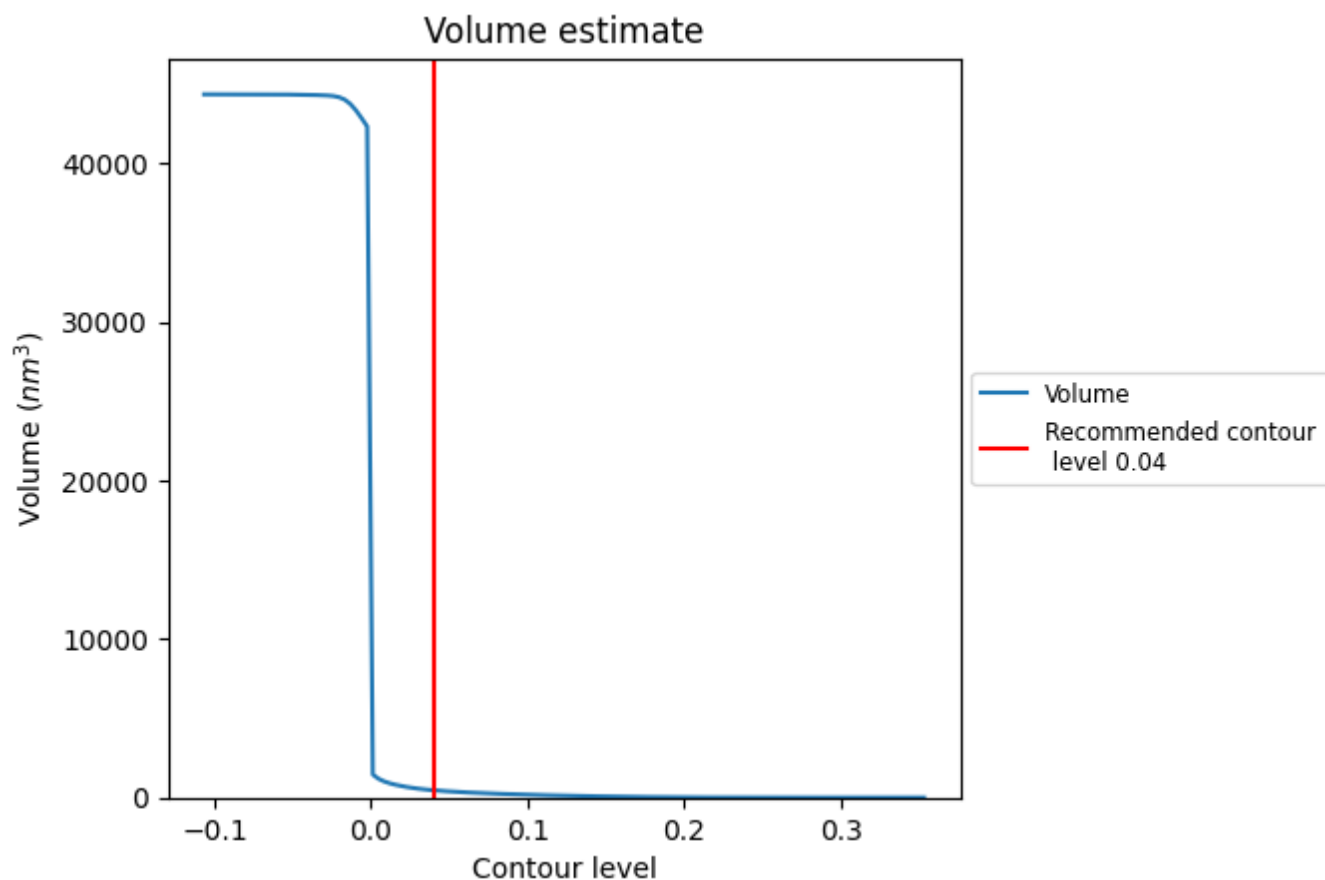
This section contains the results of statistical analysis of the map.

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.

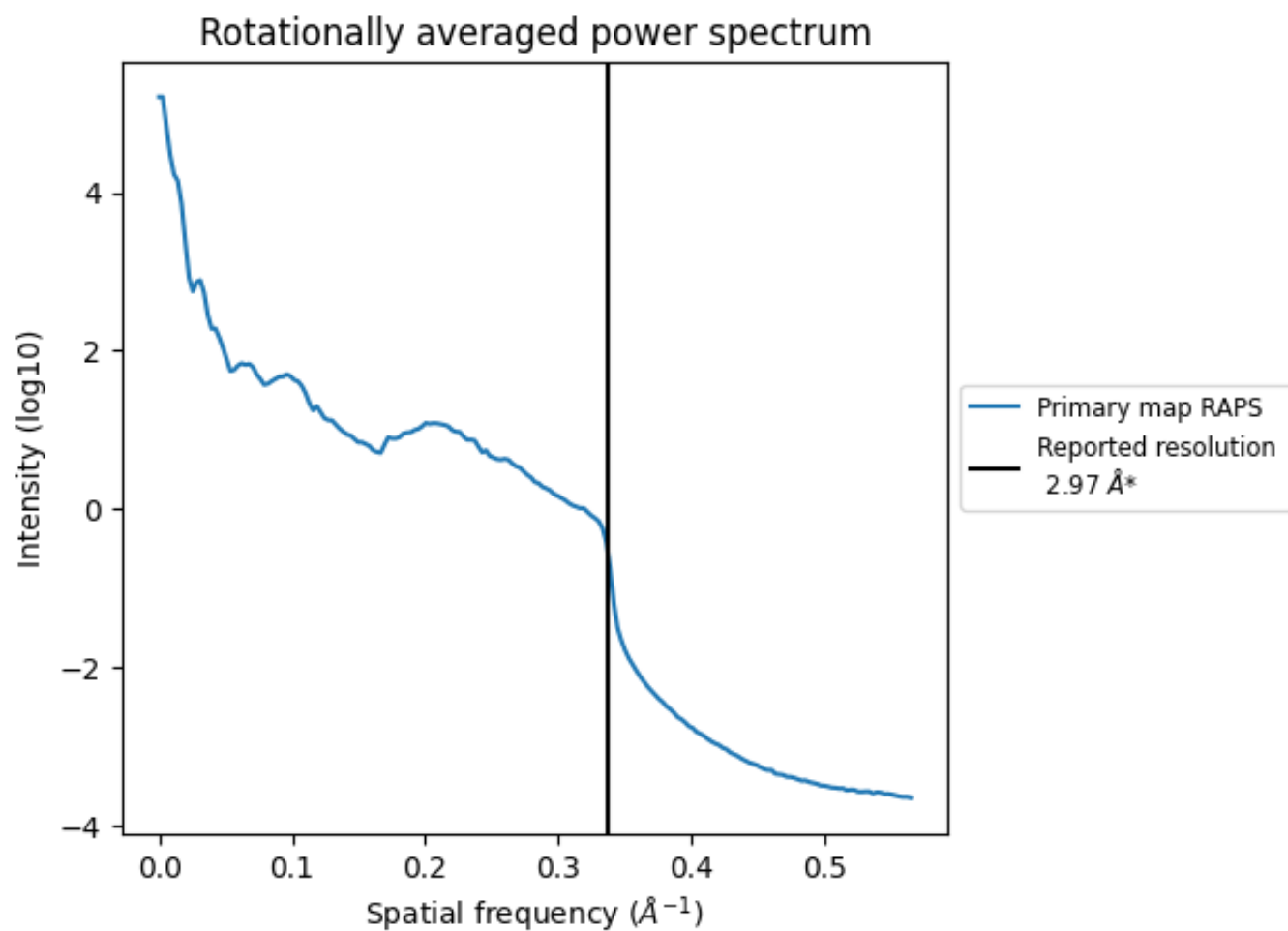
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 462 nm^3 ; this corresponds to an approximate mass of 417 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 0.337 Å⁻¹

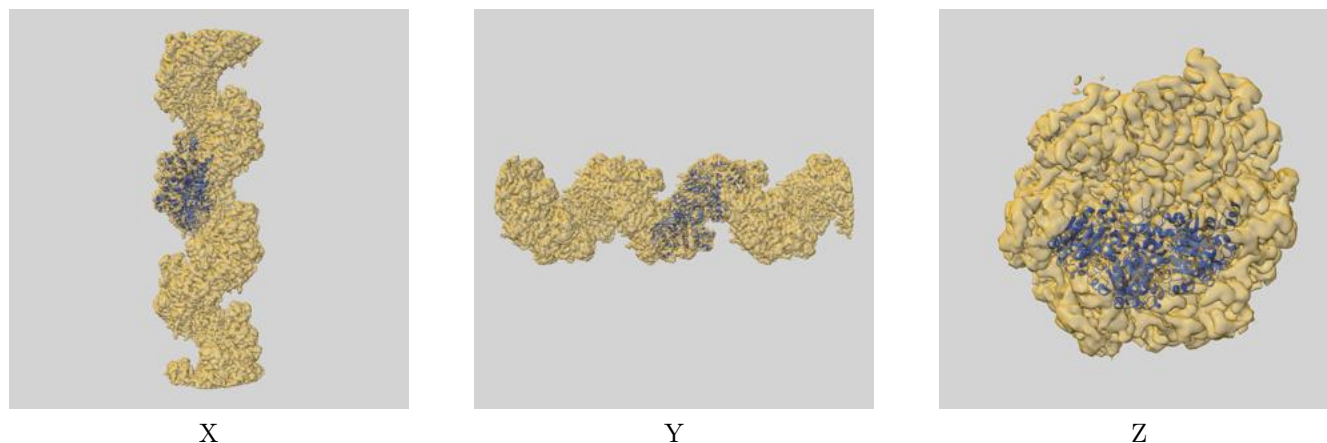
8 Fourier-Shell correlation ⓘ

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-31158 and PDB model 7EJC. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



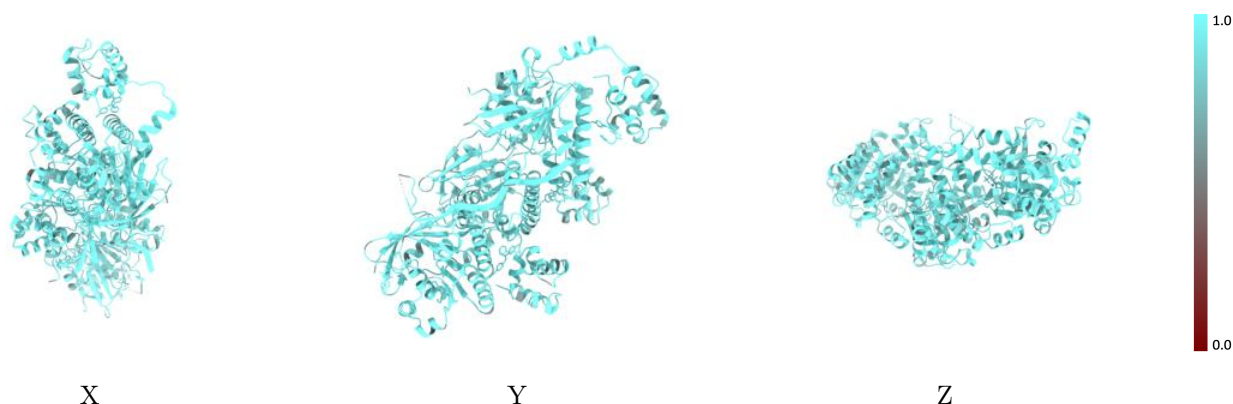
The images above show the 3D surface view of the map at the recommended contour level 0.04 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



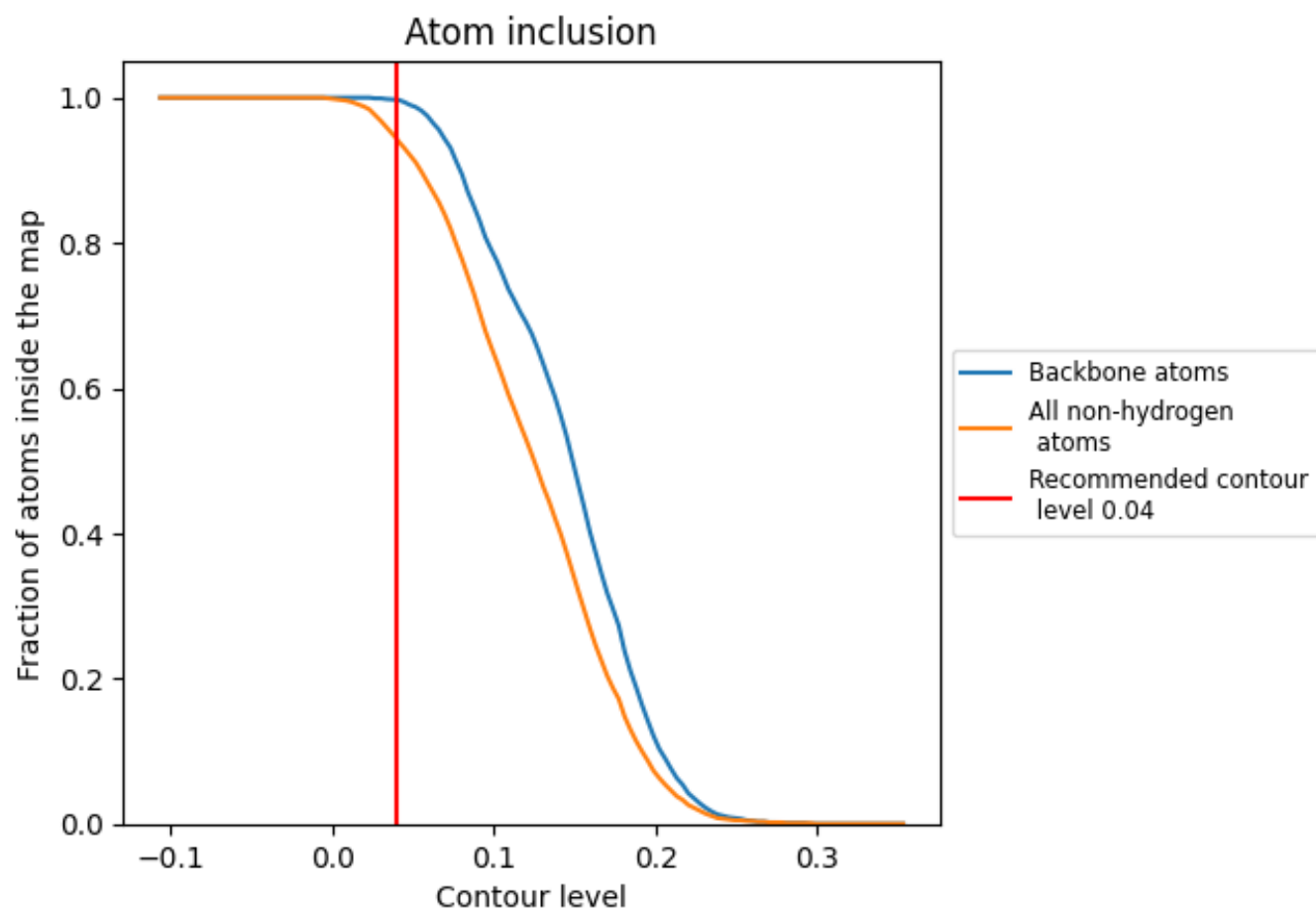
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.04).

9.4 Atom inclusion ⓘ



At the recommended contour level, 100% of all backbone atoms, 94% 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 (0.04) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9430	0.5540
A	0.9410	0.5500
B	0.9410	0.5530
C	0.9430	0.5550
G	1.0000	0.6130

