



Full wwPDB EM Validation Report ⓘ

Mar 14, 2026 – 04:25 AM UTC

PDB ID : 8G6E / pdb_00008g6e
EMDB ID : EMD-29764
Title : Structure of the Plasmodium falciparum 20S proteasome complexed with inhibitor TDI-8304
Authors : Hsu, H.-C.; Li, H.
Deposited on : 2023-02-15
Resolution : 2.18 Å(reported)
Based on initial model : 6MUW

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:

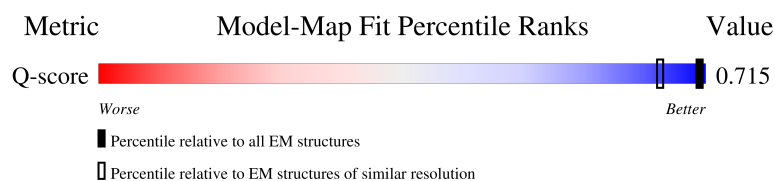
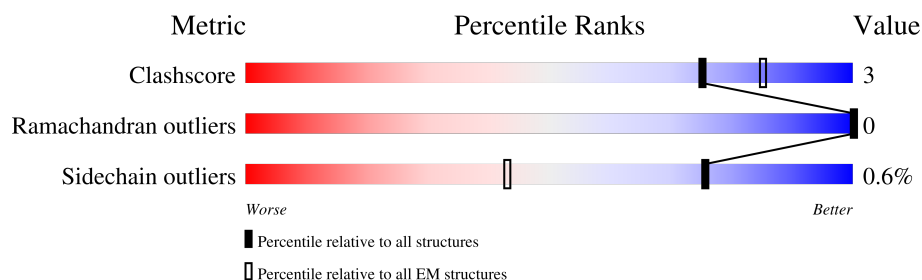
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 2.18 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.



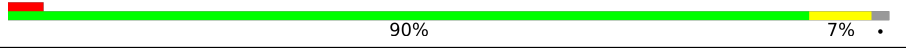
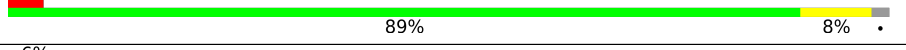
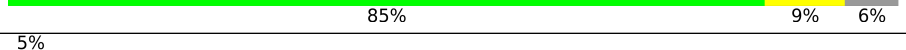
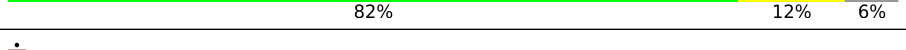
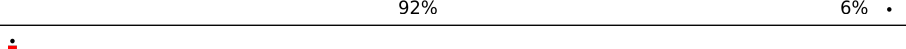
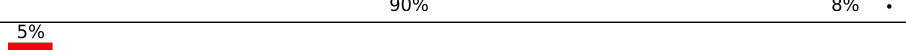
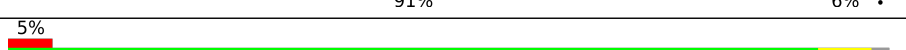
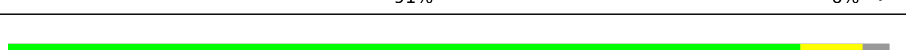
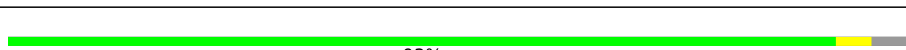
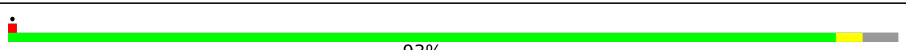
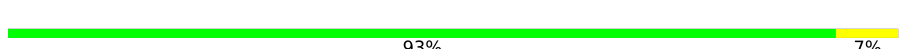
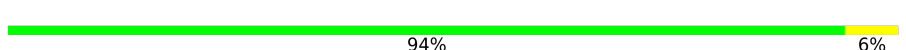
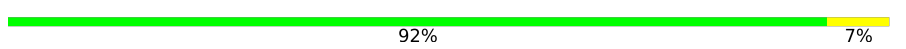
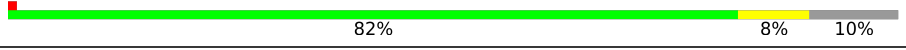
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2701 (1.70 - 2.68)

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	260	
1	O	260	
2	B	235	
2	P	235	

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Mol	Chain	Length	Quality of chain
3	C	246	 91% 8%
3	Q	246	 89% 9%
4	D	241	 90% 7%
4	R	241	 89% 8%
5	E	256	 85% 9% 6%
5	S	256	 85% 9% 6%
6	F	253	 85% 9% 6%
6	T	253	 82% 12% 6%
7	G	252	 92% 6%
7	U	252	 90% 8%
8	H	252	 91% 6%
8	V	252	 91% 6%
9	I	229	 89% 7%
9	W	229	 88% 9%
10	J	218	 93%
10	X	218	 93%
11	K	195	 93% 7%
11	Y	195	 94% 6%
12	L	211	 92% 7%
12	Z	211	 92% 7%
13	M	240	 81% 9% 10%
13	a	240	 82% 8% 10%
14	N	265	 83% 14%
14	b	265	 83% 13%

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 52266 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 Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	253	Total	C	N	O	S	0	0
			2006	1259	337	395	15		
1	O	253	Total	C	N	O	S	0	0
			2006	1259	337	395	15		

- Molecule 2 is a protein called Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	233	Total	C	N	O	S	0	0
			1851	1189	302	353	7		
2	P	233	Total	C	N	O	S	0	0
			1851	1189	302	353	7		

- Molecule 3 is a protein called Proteasome subunit alpha type.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	242	Total	C	N	O	S	0	0
			1938	1241	315	379	3		
3	Q	242	Total	C	N	O	S	0	0
			1938	1241	315	379	3		

- Molecule 4 is a protein called Proteasome subunit alpha type.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	236	Total	C	N	O	S	0	0
			1870	1192	318	352	8		
4	R	236	Total	C	N	O	S	0	0
			1870	1192	318	352	8		

- Molecule 5 is a protein called Proteasome subunit alpha type.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	241	Total	C	N	O	S	0	0
			1867	1175	312	369	11		
5	S	241	Total	C	N	O	S	0	0
			1867	1175	312	369	11		

- Molecule 6 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	239	Total	C	N	O	S	0	0
			1897	1204	313	369	11		
6	T	239	Total	C	N	O	S	0	0
			1897	1204	313	369	11		

- Molecule 7 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	247	Total	C	N	O	S	0	0
			2025	1289	338	385	13		
7	U	247	Total	C	N	O	S	0	0
			2025	1289	338	385	13		

- Molecule 8 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	246	Total	C	N	O	S	0	0
			1991	1257	339	383	12		
8	V	246	Total	C	N	O	S	0	0
			1991	1257	339	383	12		

- Molecule 9 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	221	Total	C	N	O	S	1	0
			1703	1076	297	316	14		
9	W	221	Total	C	N	O	S	1	0
			1703	1076	297	316	14		

- Molecule 10 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	210	Total	C	N	O	S	0	0
			1653	1055	269	315	14		

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Mol	Chain	Residues	Atoms					AltConf	Trace
10	X	210	Total	C	N	O	S	0	0
			1653	1055	269	315	14		

- Molecule 11 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	195	Total	C	N	O	S	0	0
			1614	1042	266	298	8		
11	Y	195	Total	C	N	O	S	0	0
			1614	1042	266	298	8		

- Molecule 12 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	211	Total	C	N	O	S	0	0
			1663	1060	275	320	8		
12	Z	211	Total	C	N	O	S	0	0
			1663	1060	275	320	8		

- Molecule 13 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	215	Total	C	N	O	S	0	0
			1711	1096	285	323	7		
13	a	215	Total	C	N	O	S	0	0
			1711	1096	285	323	7		

- Molecule 14 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	256	Total	C	N	O	S	1	0
			2106	1342	357	397	10		
14	b	256	Total	C	N	O	S	1	0
			2106	1342	357	397	10		

- Molecule 15 is (7S,10S,13S)-N-cyclopentyl-10-[2-(morpholin-4-yl)ethyl]-9,12-dioxo-13-(2-oxo pyrrolidin-1-yl)-2-oxa-8,11-diazabicyclo[13.3.1]nonadeca-1(19),15,17-triene-7-carboxamide (CCD ID: YRE) (formula: C₃₂H₄₇N₅O₆) (labeled as "Ligand of Interest" by depositor).



- Molecule 16 is water.

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Mol	Chain	Residues	Atoms		AltConf
16	G	7	Total 7	O 7	0
16	H	14	Total 14	O 14	0
16	I	7	Total 7	O 7	0
16	J	9	Total 9	O 9	0
16	K	8	Total 8	O 8	0
16	L	5	Total 5	O 5	0
16	M	6	Total 6	O 6	0
16	N	13	Total 13	O 13	0
16	O	5	Total 5	O 5	0
16	P	3	Total 3	O 3	0
16	Q	11	Total 11	O 11	0
16	R	9	Total 9	O 9	0
16	S	7	Total 7	O 7	0
16	T	11	Total 11	O 11	0
16	U	9	Total 9	O 9	0
16	V	9	Total 9	O 9	0
16	W	7	Total 7	O 7	0
16	X	10	Total 10	O 10	0
16	Y	7	Total 7	O 7	0
16	Z	3	Total 3	O 3	0
16	a	7	Total 7	O 7	0

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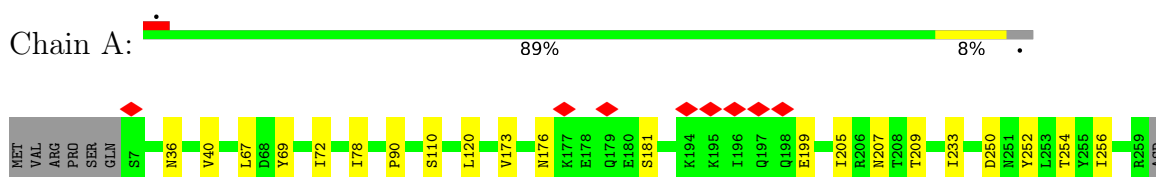
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Mol	Chain	Residues	Atoms		AltConf
16	b	11	Total	O	0
			11	11	

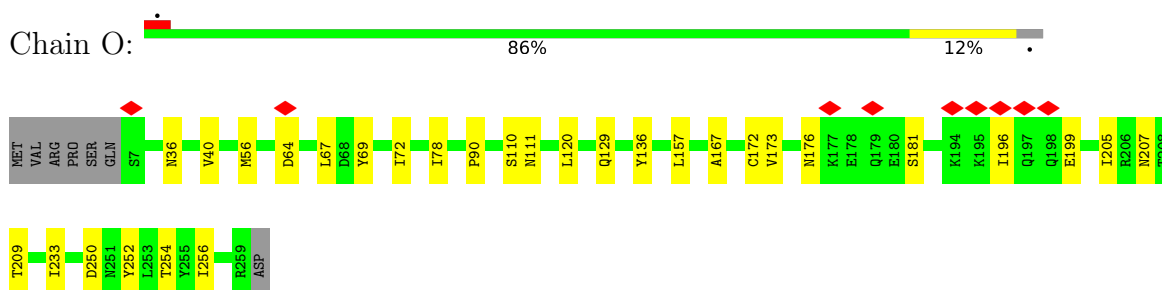
3 Residue-property plots [i](#)

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.

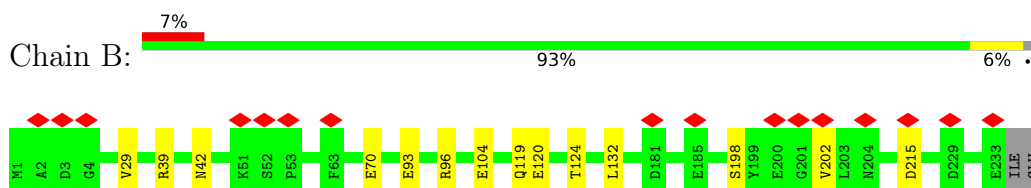
- Molecule 1: Proteasome subunit alpha type-6



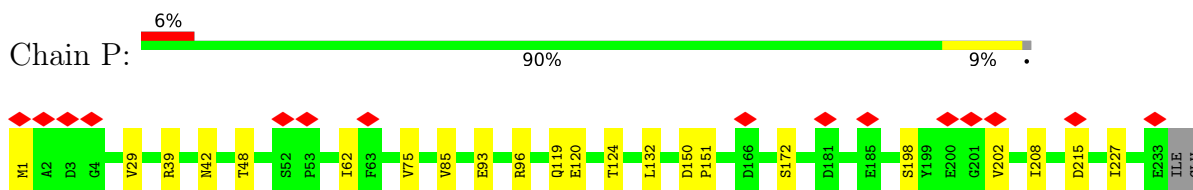
- Molecule 1: Proteasome subunit alpha type-6



- Molecule 2: Proteasome subunit alpha type-2



- Molecule 2: Proteasome subunit alpha type-2

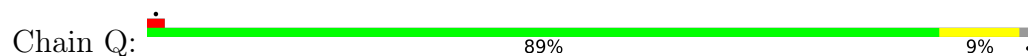


- Molecule 3: Proteasome subunit alpha type





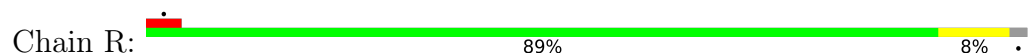
- Molecule 3: Proteasome subunit alpha type



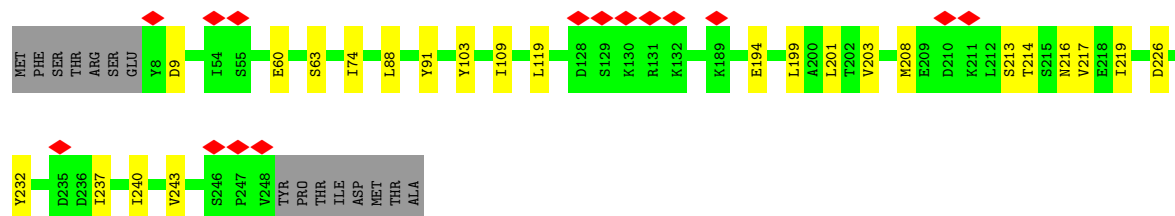
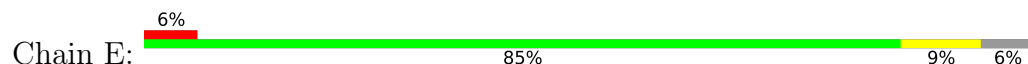
- Molecule 4: Proteasome subunit alpha type



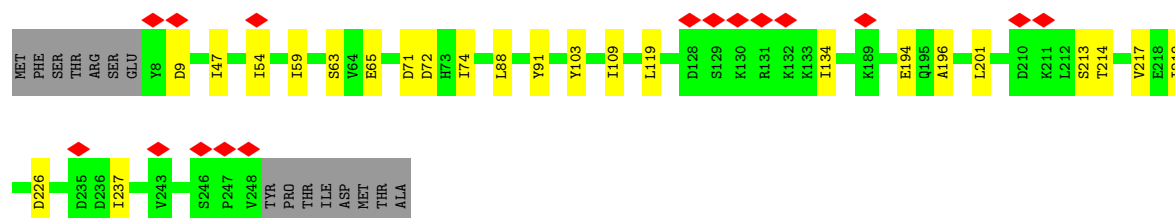
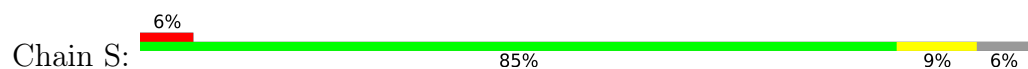
- Molecule 4: Proteasome subunit alpha type



- Molecule 5: Proteasome subunit alpha type

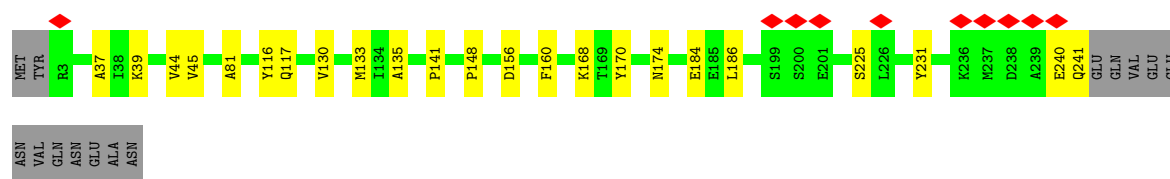


- Molecule 5: Proteasome subunit alpha type



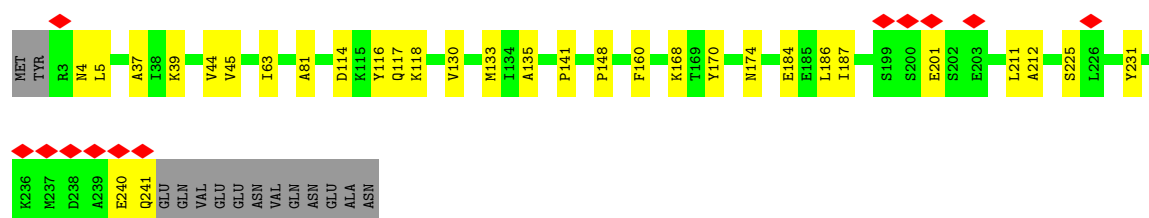
- Molecule 6: Proteasome subunit alpha type-1

Chain F: 

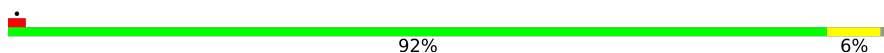


- Molecule 6: Proteasome subunit alpha type-1

Chain T: 

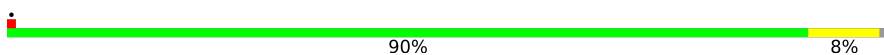


- Molecule 7: Proteasome subunit alpha type-3

Chain G: 

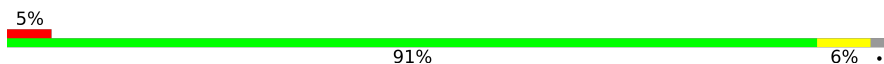


- Molecule 7: Proteasome subunit alpha type-3

Chain U: 

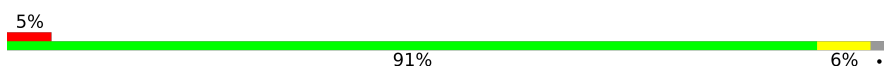


- Molecule 8: Proteasome subunit beta type-6

Chain H: 



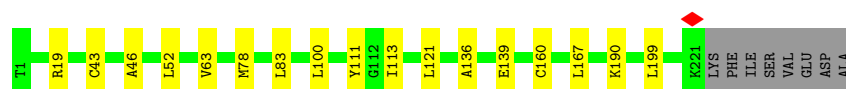
- Molecule 8: Proteasome subunit beta type-6

Chain V: 



- Molecule 9: Proteasome subunit beta

Chain I:  89% 7%



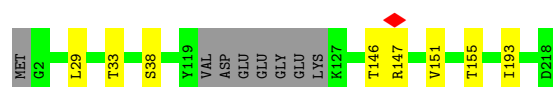
- Molecule 9: Proteasome subunit beta

Chain W:  88% 9%

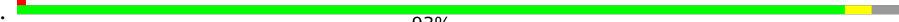


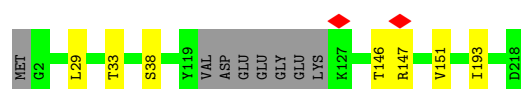
- Molecule 10: Proteasome subunit beta

Chain J:  93%

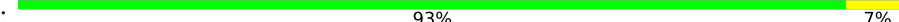


- Molecule 10: Proteasome subunit beta

Chain X:  93%

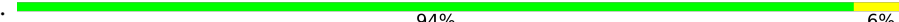


- Molecule 11: Proteasome subunit beta

Chain K:  93% 7%



- Molecule 11: Proteasome subunit beta

Chain Y:  94% 6%

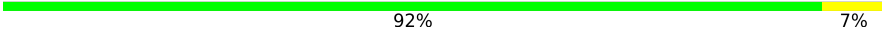


- Molecule 12: Proteasome subunit beta

Chain L:  92% 7%



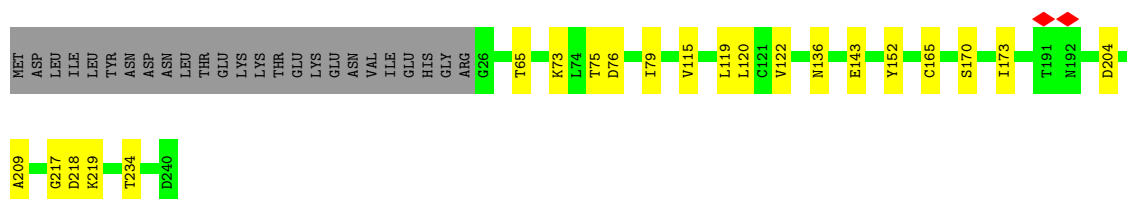
- Molecule 12: Proteasome subunit beta

Chain Z:  92% 7%



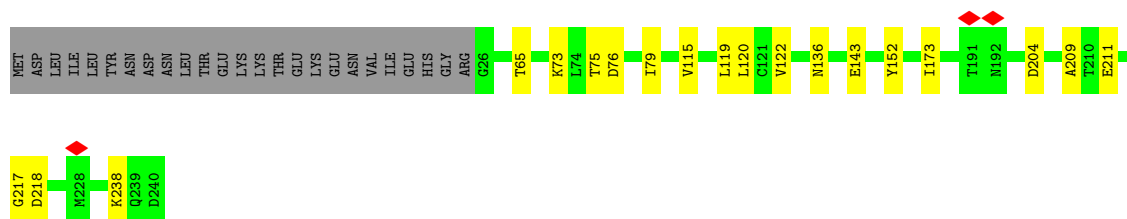
• Molecule 13: Proteasome subunit beta

Chain M:  81% 9% 10%



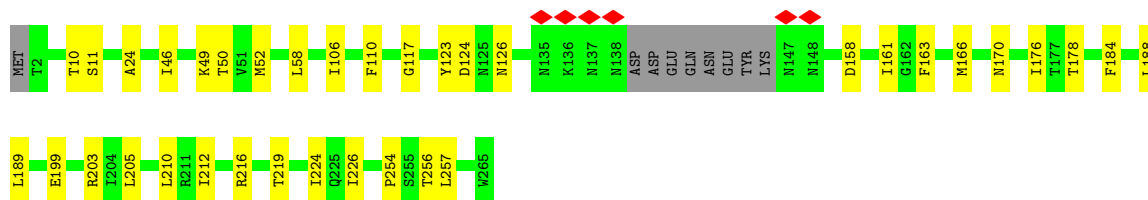
• Molecule 13: Proteasome subunit beta

Chain a:  82% 8% 10%



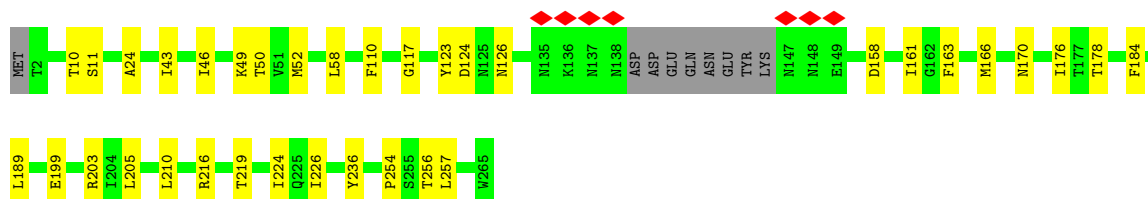
• Molecule 14: Proteasome subunit beta

Chain N:  83% 14% .



• Molecule 14: Proteasome subunit beta

Chain b:  83% 13% .



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	305581	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 ($e^-/\text{\AA}^2$)	66	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.200	Depositor
Minimum map value	-0.094	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.025	Depositor
Map size (Å)	317.952, 317.952, 317.952	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.828, 0.828, 0.828	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: YRE

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	0.18	0/2033	0.31	0/2741
1	O	0.18	0/2033	0.31	0/2741
2	B	0.19	0/1885	0.32	0/2545
2	P	0.19	0/1885	0.32	0/2545
3	C	0.19	0/1974	0.33	0/2673
3	Q	0.19	0/1974	0.33	0/2673
4	D	0.18	0/1900	0.31	0/2563
4	R	0.18	0/1900	0.31	0/2563
5	E	0.18	0/1893	0.32	0/2556
5	S	0.18	0/1893	0.32	0/2556
6	F	0.18	0/1931	0.31	0/2600
6	T	0.18	0/1931	0.31	0/2600
7	G	0.19	0/2069	0.31	0/2796
7	U	0.19	0/2069	0.31	0/2796
8	H	0.22	0/2024	0.33	0/2720
8	V	0.22	0/2024	0.33	0/2720
9	I	0.21	0/1740	0.35	0/2365
9	W	0.21	0/1740	0.35	0/2365
10	J	0.20	0/1681	0.35	0/2268
10	X	0.20	0/1681	0.35	0/2268
11	K	0.20	0/1649	0.34	0/2223
11	Y	0.23	0/1649	0.38	0/2223
12	L	0.21	0/1697	0.35	0/2286
12	Z	0.22	0/1697	0.35	0/2286
13	M	0.20	0/1744	0.36	0/2360
13	a	0.20	0/1744	0.37	0/2360
14	N	0.20	0/2152	0.34	0/2910
14	b	0.20	0/2152	0.34	0/2910
All	All	0.20	0/52744	0.33	0/71212

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	2006	0	2003	12	0
1	O	2006	0	2003	20	0
2	B	1851	0	1868	10	0
2	P	1851	0	1868	16	0
3	C	1938	0	1932	12	0
3	Q	1938	0	1932	13	0
4	D	1870	0	1903	13	0
4	R	1870	0	1903	16	0
5	E	1867	0	1876	19	0
5	S	1867	0	1876	17	0
6	F	1897	0	1892	13	0
6	T	1897	0	1892	17	0
7	G	2025	0	1973	10	0
7	U	2025	0	1973	13	0
8	H	1991	0	1964	13	0
8	V	1991	0	1964	14	0
9	I	1703	0	1704	11	0
9	W	1703	0	1704	14	0
10	J	1653	0	1645	4	0
10	X	1653	0	1645	3	0
11	K	1614	0	1584	8	0
11	Y	1614	0	1584	7	0
12	L	1663	0	1618	9	0
12	Z	1663	0	1618	9	0
13	M	1711	0	1719	12	0
13	a	1711	0	1719	12	0
14	N	2106	0	2062	24	0
14	b	2106	0	2062	22	0
15	C	43	0	0	0	0
15	I	43	0	0	0	0
15	L	43	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	Q	43	0	0	0	0
15	W	43	0	0	0	0
15	Z	43	0	0	0	0
16	A	4	0	0	0	0
16	B	4	0	0	0	0
16	C	8	0	0	0	0
16	D	6	0	0	0	0
16	E	7	0	0	0	0
16	F	11	0	0	0	0
16	G	7	0	0	0	0
16	H	14	0	0	0	0
16	I	7	0	0	0	0
16	J	9	0	0	0	0
16	K	8	0	0	0	0
16	L	5	0	0	0	0
16	M	6	0	0	0	0
16	N	13	0	0	0	0
16	O	5	0	0	0	0
16	P	3	0	0	0	0
16	Q	11	0	0	0	0
16	R	9	0	0	0	0
16	S	7	0	0	0	0
16	T	11	0	0	0	0
16	U	9	0	0	0	0
16	V	9	0	0	0	0
16	W	7	0	0	0	0
16	X	10	0	0	0	0
16	Y	7	0	0	0	0
16	Z	3	0	0	0	0
16	a	7	0	0	0	0
16	b	11	0	0	0	0
All	All	52266	0	51486	319	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 (319) 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)
8:V:104:ASN:OD1	8:V:104:ASN:O	1.94	0.84
8:V:174:ILE:HD11	14:b:58:LEU:HD21	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:174:ILE:HD11	14:N:58:LEU:HD21	1.73	0.69
5:S:88:LEU:HD23	5:S:119:LEU:HD23	1.72	0.69
3:C:35:ILE:HD13	3:C:195:THR:CG2	2.23	0.68
3:Q:35:ILE:HD13	3:Q:195:THR:CG2	2.22	0.68
6:T:44:VAL:HG23	6:T:141:PRO:HB3	1.76	0.67
14:N:124:ASP:OD2	14:N:126:ASN:ND2	2.28	0.67
4:D:83:LEU:CD2	4:D:129:ILE:HD11	2.25	0.67
5:S:91:TYR:CG	5:S:119:LEU:HD22	2.31	0.66
14:N:46:ILE:HD12	14:N:50:THR:HG22	1.78	0.66
9:I:43:CYS:HB3	9:I:100:LEU:HD12	1.78	0.66
4:R:83:LEU:CD2	4:R:129:ILE:HD11	2.26	0.66
6:T:117:GLN:NE2	7:U:86:ASP:OD1	2.28	0.66
5:E:91:TYR:CG	5:E:119:LEU:HD22	2.31	0.65
7:U:207:MET:HE1	7:U:215:LEU:HB2	1.77	0.65
4:D:83:LEU:HD22	4:D:129:ILE:HD11	1.79	0.65
4:R:83:LEU:HD22	4:R:129:ILE:HD11	1.78	0.64
2:P:39:ARG:NH1	3:Q:57:ASP:OD2	2.30	0.64
5:E:88:LEU:HD23	5:E:119:LEU:HD23	1.78	0.64
7:G:207:MET:HE1	7:G:215:LEU:HB2	1.79	0.64
6:F:156:ASP:OD1	7:G:62:LYS:NZ	2.31	0.64
6:F:117:GLN:NE2	7:G:86:ASP:OD1	2.31	0.64
14:b:199:GLU:OE2	14:b:203:ARG:NH2	2.31	0.63
14:b:124:ASP:OD2	14:b:126:ASN:ND2	2.32	0.62
9:W:43:CYS:HB3	9:W:100:LEU:HD12	1.81	0.62
14:N:199:GLU:OE2	14:N:203:ARG:NH2	2.32	0.62
14:b:254:PRO:HD2	14:b:257:LEU:HD12	1.80	0.61
6:F:44:VAL:HG23	6:F:141:PRO:HB3	1.81	0.61
2:P:208:ILE:O	2:P:208:ILE:HD12	2.02	0.60
5:E:194:GLU:N	5:E:194:GLU:OE1	2.35	0.59
14:b:46:ILE:HD12	14:b:50:THR:HG22	1.84	0.59
14:N:254:PRO:HD2	14:N:257:LEU:HD12	1.84	0.59
1:A:67:LEU:CD2	1:A:72:ILE:HD11	2.33	0.59
14:b:46:ILE:HD11	14:b:52:MET:HB2	1.84	0.59
8:V:89:ASP:OD1	8:V:90:GLU:N	2.36	0.59
1:O:67:LEU:CD2	1:O:72:ILE:HD11	2.33	0.59
5:S:194:GLU:N	5:S:194:GLU:OE1	2.35	0.59
3:C:35:ILE:HD13	3:C:195:THR:HG23	1.85	0.58
5:S:9:ASP:OD1	5:S:9:ASP:N	2.35	0.58
12:L:58:TYR:CE1	12:L:62:ILE:HD11	2.37	0.58
14:N:46:ILE:HD11	14:N:52:MET:HB2	1.86	0.58
8:H:89:ASP:OD1	8:H:90:GLU:N	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:35:ILE:HD13	3:Q:195:THR:HG23	1.85	0.57
2:B:39:ARG:NH1	3:C:57:ASP:OD2	2.38	0.57
5:E:9:ASP:OD1	5:E:9:ASP:N	2.36	0.57
6:F:45:VAL:HG23	6:F:186:LEU:HD23	1.86	0.57
12:Z:58:TYR:CE1	12:Z:62:ILE:HD11	2.40	0.56
1:A:173:VAL:HG21	1:A:181:SER:HB3	1.88	0.56
8:H:24:THR:HG23	14:b:219:THR:HG23	1.88	0.56
2:B:42:ASN:ND2	2:B:215:ASP:OD1	2.35	0.55
6:T:168:LYS:HG2	7:U:60:MET:HE1	1.88	0.55
13:a:75:THR:HG22	13:a:76:ASP:N	2.22	0.55
14:N:10:THR:HG23	14:N:11:SER:N	2.22	0.55
2:B:93:GLU:OE2	2:B:96:ARG:NH2	2.37	0.55
11:Y:1:MET:HG2	11:Y:134:TYR:H	1.72	0.54
13:M:75:THR:HG22	13:M:76:ASP:N	2.22	0.54
13:a:143:GLU:OE1	13:a:143:GLU:N	2.38	0.54
2:B:120:GLU:O	2:B:124:THR:HG23	2.08	0.54
5:S:54:ILE:HG23	5:S:59:ILE:CG2	2.36	0.54
14:b:10:THR:HG23	14:b:11:SER:N	2.22	0.54
7:G:94:ALA:HA	7:G:118:ILE:HD11	1.90	0.54
7:U:94:ALA:HA	7:U:118:ILE:HD11	1.90	0.54
14:N:219:THR:HG23	8:V:24:THR:HG23	1.90	0.54
1:A:78:ILE:HG21	1:A:120:LEU:HD21	1.90	0.53
14:N:123:TYR:OH	14:N:158:ASP:OD2	2.25	0.53
8:V:118:TYR:CE1	8:V:238:VAL:HG23	2.44	0.53
2:P:120:GLU:O	2:P:124:THR:HG23	2.09	0.53
5:E:232:TYR:HE2	5:E:240:ILE:HD11	1.73	0.53
3:C:176:LYS:NZ	4:D:51:PRO:O	2.42	0.53
1:O:90:PRO:HG2	7:U:159:ALA:HB2	1.91	0.53
14:N:163:PHE:HE1	14:N:178:THR:HG22	1.73	0.53
8:H:59:ILE:HD11	8:H:139:ILE:HD12	1.91	0.53
2:P:208:ILE:HD11	2:P:227:ILE:HG21	1.91	0.53
13:M:143:GLU:OE1	13:M:143:GLU:N	2.41	0.52
2:P:208:ILE:CD1	2:P:227:ILE:HG21	2.39	0.52
14:b:163:PHE:HE1	14:b:178:THR:HG22	1.74	0.52
6:F:81:ALA:HB2	6:F:130:VAL:HG21	1.91	0.52
6:T:45:VAL:HG23	6:T:186:LEU:HD23	1.90	0.52
11:K:1:MET:HG2	11:K:134:TYR:H	1.74	0.52
8:V:59:ILE:HD11	8:V:139:ILE:HD12	1.91	0.52
1:A:90:PRO:HG2	7:G:159:ALA:HB2	1.92	0.52
5:E:201:LEU:HD21	5:E:219:ILE:HD11	1.92	0.52
8:H:118:TYR:CE1	8:H:238:VAL:HG23	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:250:ASP:O	1:O:254:THR:HG23	2.10	0.52
6:T:81:ALA:HB2	6:T:130:VAL:HG21	1.91	0.52
1:A:69:TYR:HA	1:A:72:ILE:HD12	1.93	0.51
1:O:167:ALA:HB2	2:P:62:ILE:HD11	1.90	0.51
9:I:46:ALA:C	9:I:52:LEU:HD11	2.36	0.51
9:I:83:LEU:HD23	9:I:113:ILE:HD12	1.92	0.51
14:N:216:ARG:NE	9:W:139:GLU:OE2	2.42	0.51
4:R:93:ARG:NH1	4:R:97:ASN:OD1	2.43	0.51
6:T:37:ALA:HB2	6:T:133:MET:SD	2.51	0.51
8:V:24:THR:HG22	8:V:24:THR:O	2.11	0.51
13:M:219:LYS:HD3	13:M:234:THR:HG21	1.92	0.51
1:A:36:ASN:OD1	1:A:176:ASN:ND2	2.40	0.50
1:O:78:ILE:HG21	1:O:120:LEU:HD21	1.93	0.50
14:b:161:ILE:HG22	14:b:176:ILE:HG13	1.93	0.50
10:J:33:THR:O	10:J:193:ILE:HG22	2.10	0.50
12:L:18:SER:O	12:L:31:VAL:HG22	2.11	0.50
1:O:205:ILE:HG23	1:O:233:ILE:CD1	2.41	0.50
2:P:93:GLU:OE2	2:P:96:ARG:NH2	2.38	0.50
11:K:151:LEU:HD21	11:K:159:ILE:HD12	1.93	0.50
14:N:189:LEU:HD23	14:N:205:LEU:HD21	1.93	0.50
1:A:250:ASP:O	1:A:254:THR:HG23	2.11	0.50
8:H:24:THR:HG22	8:H:24:THR:O	2.11	0.50
8:V:118:TYR:HE1	8:V:238:VAL:HG23	1.76	0.50
6:F:44:VAL:HG11	6:F:135:ALA:HB1	1.94	0.50
1:O:173:VAL:HG21	1:O:181:SER:HB3	1.93	0.50
10:X:33:THR:O	10:X:193:ILE:HG22	2.11	0.50
1:A:205:ILE:HG23	1:A:233:ILE:CD1	2.41	0.50
4:D:193:ILE:HG21	4:D:204:VAL:HG21	1.94	0.50
5:E:213:SER:O	5:E:217:VAL:HG12	2.12	0.50
8:V:62:TYR:CE2	8:V:139:ILE:HD11	2.46	0.50
12:Z:201:THR:HG23	12:Z:209:TYR:CE1	2.47	0.50
2:P:29:VAL:HG22	2:P:132:LEU:HA	1.94	0.49
8:H:62:TYR:CE2	8:H:139:ILE:HD11	2.47	0.49
1:O:199:GLU:OE2	1:O:207:ASN:ND2	2.42	0.49
10:X:146:THR:HG22	10:X:147:ARG:N	2.27	0.49
14:N:161:ILE:HG22	14:N:176:ILE:HG13	1.94	0.49
1:A:67:LEU:HD21	1:A:72:ILE:HD11	1.94	0.49
8:H:9:ASP:OD1	8:H:10:ASN:N	2.45	0.49
1:O:69:TYR:HA	1:O:72:ILE:HD12	1.94	0.49
6:F:37:ALA:HB2	6:F:133:MET:SD	2.53	0.49
8:V:9:ASP:OD1	8:V:10:ASN:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:W:46:ALA:C	9:W:52:LEU:HD11	2.38	0.49
10:J:146:THR:HG22	10:J:147:ARG:N	2.27	0.49
5:S:91:TYR:CD2	5:S:119:LEU:HD22	2.47	0.49
2:B:29:VAL:HG22	2:B:132:LEU:HA	1.93	0.49
3:C:149:HIS:O	3:C:156:TYR:HA	2.13	0.48
9:I:139:GLU:OE2	14:b:216:ARG:NE	2.45	0.48
12:L:201:THR:HG23	12:L:209:TYR:CE1	2.48	0.48
3:Q:206:LYS:NZ	3:Q:208:GLU:OE2	2.35	0.48
6:T:44:VAL:HG11	6:T:135:ALA:HB1	1.95	0.48
14:b:210:LEU:HD13	14:b:224:ILE:HG21	1.95	0.48
5:E:232:TYR:CE2	5:E:240:ILE:HD11	2.49	0.48
8:H:118:TYR:HE1	8:H:238:VAL:HG23	1.77	0.48
6:F:168:LYS:HG2	7:G:60:MET:HE1	1.95	0.48
1:A:209:THR:HG21	1:A:252:TYR:CE2	2.49	0.48
10:X:29:LEU:HB3	10:X:38:SER:HB3	1.96	0.48
3:Q:149:HIS:O	3:Q:156:TYR:HA	2.14	0.48
4:R:199:LEU:HD23	4:R:230:ASP:OD1	2.14	0.48
4:D:83:LEU:HD23	4:D:129:ILE:HD11	1.96	0.48
6:F:39:LYS:HB3	6:F:44:VAL:HG22	1.95	0.48
9:I:160:CYS:SG	9:I:190:LYS:NZ	2.85	0.48
5:S:201:LEU:HD21	5:S:219:ILE:HD11	1.95	0.48
1:O:209:THR:HG21	1:O:252:TYR:CE2	2.49	0.48
7:U:201:LEU:HD21	7:U:212:LEU:HD11	1.95	0.48
12:Z:160:ILE:HG21	12:Z:174:VAL:HG23	1.95	0.48
14:b:49:LYS:O	14:b:117:GLY:HA3	2.14	0.48
12:L:40:ASN:OD1	12:L:73:LYS:NZ	2.45	0.47
14:N:210:LEU:HD13	14:N:224:ILE:HG21	1.95	0.47
10:J:155:THR:O	10:J:155:THR:HG22	2.14	0.47
12:Z:18:SER:O	12:Z:31:VAL:HG22	2.14	0.47
13:a:75:THR:HG22	13:a:76:ASP:H	1.77	0.47
4:D:148:GLU:OE2	4:D:152:ILE:HD11	2.14	0.47
1:O:67:LEU:HD21	1:O:72:ILE:HD11	1.95	0.47
5:S:213:SER:O	5:S:217:VAL:HG12	2.14	0.47
8:H:104:ASN:OD1	8:H:104:ASN:O	2.33	0.47
8:H:169:LEU:O	8:H:170:ASN:HB2	2.15	0.47
9:I:63:VAL:HG22	9:I:78:MET:HE3	1.96	0.47
10:J:29:LEU:HB3	10:J:38:SER:HB3	1.97	0.47
11:K:173:LEU:HD21	11:Y:172:LEU:HD11	1.97	0.47
14:N:24:ALA:HB2	14:N:226:ILE:HD13	1.97	0.47
1:O:157:LEU:HD21	1:O:172:CYS:SG	2.55	0.47
5:E:60:GLU:O	5:E:63:SER:OG	2.32	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:111:LEU:HD11	7:U:140:ILE:HG22	1.96	0.47
5:E:91:TYR:CD2	5:E:119:LEU:HD22	2.49	0.47
4:D:199:LEU:HD23	4:D:230:ASP:OD1	2.15	0.46
12:L:58:TYR:HE1	12:L:62:ILE:HD11	1.80	0.46
8:V:169:LEU:O	8:V:170:ASN:HB2	2.15	0.46
14:b:24:ALA:HB2	14:b:226:ILE:HD13	1.97	0.46
5:E:226:ASP:OD1	5:E:226:ASP:N	2.49	0.46
12:L:160:ILE:HG21	12:L:174:VAL:HG23	1.95	0.46
3:Q:6:ASP:OD2	4:R:3:TYR:OH	2.29	0.46
11:Y:151:LEU:HD21	11:Y:159:ILE:HD12	1.98	0.46
5:E:214:THR:HA	5:E:237:ILE:HG21	1.97	0.46
14:N:49:LYS:O	14:N:117:GLY:HA3	2.14	0.46
9:W:111:TYR:CZ	9:W:121:LEU:HD13	2.51	0.46
2:P:42:ASN:ND2	2:P:215:ASP:OD1	2.39	0.46
7:U:94:ALA:CA	7:U:118:ILE:HD11	2.46	0.46
7:G:111:LEU:HD11	7:G:140:ILE:HG22	1.98	0.46
9:W:83:LEU:HD23	9:W:113:ILE:HD12	1.98	0.46
11:Y:21:SER:OG	11:Y:32:ASP:OD2	2.34	0.46
6:T:184:GLU:OE2	6:T:231:TYR:OH	2.33	0.46
1:O:36:ASN:OD1	1:O:176:ASN:ND2	2.46	0.46
9:W:63:VAL:HG22	9:W:78:MET:HE3	1.98	0.46
6:T:39:LYS:HB3	6:T:44:VAL:HG22	1.98	0.45
6:T:114:ASP:O	6:T:118:LYS:HG2	2.15	0.45
1:A:40:VAL:HG22	1:A:173:VAL:HG22	1.98	0.45
5:S:214:THR:HA	5:S:237:ILE:HG21	1.99	0.45
14:b:123:TYR:OH	14:b:158:ASP:OD2	2.26	0.45
6:F:170:TYR:O	6:F:174:ASN:ND2	2.49	0.45
9:I:136:ALA:HB2	14:b:184:PHE:HA	1.99	0.45
11:K:21:SER:OG	11:K:32:ASP:OD1	2.34	0.45
11:K:172:LEU:HD11	11:Y:173:LEU:HD21	1.99	0.45
3:Q:179:TRP:O	4:R:52:LYS:NZ	2.42	0.45
1:A:199:GLU:OE2	1:A:207:ASN:ND2	2.43	0.45
5:S:226:ASP:OD1	5:S:226:ASP:N	2.49	0.45
2:B:119:GLN:NE2	3:C:82:ASP:OD1	2.48	0.44
13:M:75:THR:HG22	13:M:76:ASP:H	1.82	0.44
14:N:256:THR:OG1	8:V:114:ASN:ND2	2.39	0.44
9:I:19:ARG:NH1	9:I:167:LEU:O	2.50	0.44
1:O:129:GLN:OE1	2:P:85:VAL:HG21	2.17	0.44
7:U:80:TYR:OH	7:U:84:ASP:OD1	2.29	0.44
5:E:208:MET:HE1	5:E:216:ASN:ND2	2.33	0.44
1:O:64:ASP:OD1	1:O:64:ASP:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:193:ILE:HG21	4:R:204:VAL:HG21	1.99	0.44
6:T:4:ASN:O	6:T:5:LEU:HB2	2.17	0.44
4:D:155:ALA:HB3	5:E:63:SER:HB3	1.99	0.44
5:E:74:ILE:HG12	5:E:109:ILE:HD12	2.00	0.44
1:O:111:ASN:OD1	9:W:82:ARG:NH2	2.51	0.44
13:a:119:LEU:O	13:a:122:VAL:HG22	2.18	0.44
6:F:184:GLU:OE2	6:F:231:TYR:OH	2.35	0.44
14:b:189:LEU:HD23	14:b:205:LEU:HD21	2.00	0.44
11:K:8:ARG:HG2	11:K:128:ASN:HA	2.00	0.43
13:M:119:LEU:O	13:M:122:VAL:HG22	2.18	0.43
14:N:184:PHE:HA	9:W:136:ALA:HB2	1.98	0.43
3:Q:31:ALA:O	3:Q:166:THR:N	2.44	0.43
13:a:211:GLU:OE1	13:a:238:LYS:NZ	2.48	0.43
5:E:103:TYR:HB3	13:M:115:VAL:HG13	2.00	0.43
5:E:243:VAL:O	5:E:243:VAL:HG22	2.18	0.43
3:Q:182:ASP:OD1	3:Q:182:ASP:N	2.52	0.43
4:R:33:VAL:HG23	4:R:193:ILE:HD12	1.99	0.43
7:G:94:ALA:CA	7:G:118:ILE:HD11	2.46	0.43
2:P:1:MET:SD	2:P:1:MET:N	2.87	0.43
3:C:182:ASP:N	3:C:182:ASP:OD1	2.51	0.43
4:R:36:LYS:HB2	4:R:144:ILE:HD12	1.99	0.43
4:R:83:LEU:HD23	4:R:129:ILE:HD11	1.97	0.43
4:R:121:ARG:NE	5:S:134:ILE:HG22	2.33	0.43
7:U:201:LEU:CD2	7:U:212:LEU:HD11	2.49	0.43
12:L:201:THR:HG23	12:L:209:TYR:CD1	2.53	0.43
14:N:216:ARG:NH2	9:W:139:GLU:OE2	2.50	0.43
5:S:88:LEU:HD23	5:S:119:LEU:CD2	2.45	0.43
9:W:19:ARG:NH1	9:W:167:LEU:O	2.51	0.43
2:P:48:THR:HB	2:P:75:VAL:HG21	2.01	0.43
3:Q:202:THR:OG1	3:Q:203:GLU:N	2.51	0.43
3:C:202:THR:OG1	3:C:203:GLU:N	2.52	0.43
8:H:114:ASN:ND2	14:b:256:THR:OG1	2.39	0.43
14:N:188:LEU:HD11	14:N:212:ILE:CD1	2.48	0.43
13:a:173:ILE:HD11	13:a:209:ALA:CB	2.49	0.43
9:W:129:SER:OG	9:W:166:ASP:OD2	2.34	0.42
13:M:217:GLY:O	13:M:218:ASP:HB2	2.19	0.42
13:a:217:GLY:O	13:a:218:ASP:HB2	2.19	0.42
2:B:119:GLN:HG3	3:C:81:ALA:HB1	2.01	0.42
3:C:31:ALA:O	3:C:166:THR:N	2.45	0.42
3:Q:118:LYS:O	3:Q:122:THR:HG23	2.19	0.42
6:T:201:GLU:OE1	6:T:201:GLU:N	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:5:SER:OG	7:U:6:ALA:N	2.52	0.42
14:b:110:PHE:CE2	14:b:166:MET:HE1	2.54	0.42
14:b:163:PHE:O	14:b:170:ASN:HA	2.19	0.42
2:B:202:VAL:O	2:B:202:VAL:HG23	2.18	0.42
14:N:163:PHE:O	14:N:170:ASN:HA	2.19	0.42
4:R:208:LEU:HD23	4:R:210:THR:HG23	2.00	0.42
2:P:202:VAL:O	2:P:202:VAL:HG23	2.20	0.42
12:Z:15:ALA:HA	12:Z:175:ARG:O	2.20	0.42
6:T:116:TYR:CD1	6:T:148:PRO:HA	2.54	0.42
13:a:173:ILE:HD11	13:a:209:ALA:HB2	2.02	0.42
1:O:196:ILE:HG22	1:O:196:ILE:O	2.19	0.42
9:W:43:CYS:CB	9:W:100:LEU:HD12	2.47	0.42
7:G:28:TYR:OH	7:G:155:GLU:HB3	2.20	0.41
9:I:111:TYR:CZ	9:I:121:LEU:HD13	2.54	0.41
5:S:74:ILE:HG12	5:S:109:ILE:CD1	2.49	0.41
5:S:103:TYR:HB3	13:a:115:VAL:HG13	2.01	0.41
7:U:28:TYR:OH	7:U:155:GLU:HB3	2.20	0.41
4:D:208:LEU:HD23	4:D:210:THR:HG23	2.02	0.41
6:F:116:TYR:CD1	6:F:148:PRO:HA	2.55	0.41
1:O:136:TYR:CD2	2:P:1:MET:HA	2.55	0.41
8:V:59:ILE:HD11	8:V:136:THR:HA	2.03	0.41
12:Z:201:THR:HG23	12:Z:209:TYR:CD1	2.55	0.41
3:C:118:LYS:O	3:C:122:THR:HG23	2.19	0.41
9:I:199:LEU:HG	13:a:204:ASP:HB2	2.01	0.41
12:L:15:ALA:HA	12:L:175:ARG:O	2.21	0.41
4:D:5:ARG:O	4:D:123:GLY:N	2.47	0.41
7:G:71:HIS:HA	7:G:76:ILE:O	2.20	0.41
14:N:110:PHE:CE2	14:N:166:MET:HE1	2.55	0.41
1:O:40:VAL:HG22	1:O:173:VAL:HG22	2.02	0.41
1:O:56:MET:HE3	1:O:67:LEU:HD21	2.03	0.41
6:T:187:ILE:HG23	6:T:211:LEU:HD21	2.03	0.41
9:I:139:GLU:OE2	14:b:216:ARG:NH2	2.48	0.41
13:M:173:ILE:HD11	13:M:209:ALA:CB	2.50	0.41
6:T:170:TYR:O	6:T:174:ASN:ND2	2.53	0.41
8:H:59:ILE:HD11	8:H:136:THR:HA	2.03	0.41
2:P:119:GLN:HG3	3:Q:81:ALA:HB1	2.03	0.41
4:R:155:ALA:HB3	5:S:63:SER:HB3	2.02	0.41
7:U:66:TYR:OH	7:U:84:ASP:OD1	2.27	0.41
2:P:150:ASP:CG	2:P:151:PRO:HD2	2.46	0.41
5:S:71:ASP:OD1	5:S:72:ASP:N	2.48	0.41
12:Z:20:ALA:HB2	12:Z:31:VAL:HG11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:a:120:LEU:HD23	13:a:152:TYR:CZ	2.56	0.41
13:M:165:CYS:SG	13:M:170:SER:HA	2.61	0.41
13:M:204:ASP:HB2	9:W:199:LEU:HG	2.01	0.41
14:N:10:THR:HG23	14:N:11:SER:H	1.86	0.41
14:N:219:THR:HG23	8:V:24:THR:CG2	2.51	0.41
6:T:63:ILE:HD12	6:T:212:ALA:HB2	2.03	0.41
9:W:220:ILE:HG22	9:W:221:LYS:HG3	2.03	0.41
12:Z:40:ASN:OD1	12:Z:73:LYS:NZ	2.46	0.41
2:B:70:GLU:O	2:B:104:GLU:CD	2.64	0.41
12:L:149:LEU:O	12:L:153:VAL:HG23	2.21	0.41
3:Q:124:TYR:OH	4:R:118:PHE:HZ	2.04	0.41
4:D:92:GLN:HG3	11:K:67:LEU:HB2	2.03	0.40
5:E:199:LEU:O	5:E:203:VAL:HG23	2.21	0.40
13:M:120:LEU:HD23	13:M:152:TYR:CZ	2.55	0.40
12:Z:149:LEU:O	12:Z:153:VAL:HG23	2.21	0.40
2:B:42:ASN:OD1	2:B:42:ASN:N	2.53	0.40
11:K:3:THR:O	11:K:3:THR:HG23	2.21	0.40
4:R:91:CYS:SG	4:R:107:ILE:HD11	2.62	0.40
5:S:47:ILE:HD13	5:S:196:ALA:HB1	2.02	0.40
14:b:43:ILE:O	14:b:236:TYR:OH	2.36	0.40
5:E:88:LEU:HD23	5:E:119:LEU:CD2	2.48	0.40
6:F:240:GLU:O	6:F:241:GLN:C	2.65	0.40
3:C:71:HIS:HA	3:C:222:VAL:HG11	2.04	0.40
13:M:73:LYS:HA	13:M:79:ILE:HG22	2.03	0.40
4:R:92:GLN:HG3	11:Y:67:LEU:HB2	2.02	0.40
4:D:33:VAL:HG23	4:D:193:ILE:HD12	2.03	0.40
4:D:167:ILE:O	4:D:170:GLU:HG3	2.22	0.40
6:T:240:GLU:O	6:T:241:GLN:C	2.65	0.40
11:Y:166:GLU:HA	11:Y:166:GLU:OE1	2.21	0.40
13:a:73:LYS:HA	13:a:79:ILE:HG22	2.03	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	251/260 (96%)	249 (99%)	2 (1%)	0	100	100
1	O	251/260 (96%)	248 (99%)	3 (1%)	0	100	100
2	B	231/235 (98%)	226 (98%)	5 (2%)	0	100	100
2	P	231/235 (98%)	225 (97%)	6 (3%)	0	100	100
3	C	240/246 (98%)	237 (99%)	3 (1%)	0	100	100
3	Q	240/246 (98%)	236 (98%)	4 (2%)	0	100	100
4	D	234/241 (97%)	227 (97%)	7 (3%)	0	100	100
4	R	234/241 (97%)	228 (97%)	6 (3%)	0	100	100
5	E	239/256 (93%)	227 (95%)	12 (5%)	0	100	100
5	S	239/256 (93%)	229 (96%)	10 (4%)	0	100	100
6	F	237/253 (94%)	234 (99%)	3 (1%)	0	100	100
6	T	237/253 (94%)	233 (98%)	4 (2%)	0	100	100
7	G	245/252 (97%)	243 (99%)	2 (1%)	0	100	100
7	U	245/252 (97%)	243 (99%)	2 (1%)	0	100	100
8	H	242/252 (96%)	236 (98%)	6 (2%)	0	100	100
8	V	242/252 (96%)	237 (98%)	5 (2%)	0	100	100
9	I	220/229 (96%)	215 (98%)	5 (2%)	0	100	100
9	W	220/229 (96%)	216 (98%)	4 (2%)	0	100	100
10	J	206/218 (94%)	204 (99%)	2 (1%)	0	100	100
10	X	206/218 (94%)	204 (99%)	2 (1%)	0	100	100
11	K	193/195 (99%)	189 (98%)	4 (2%)	0	100	100
11	Y	193/195 (99%)	188 (97%)	5 (3%)	0	100	100
12	L	209/211 (99%)	206 (99%)	3 (1%)	0	100	100
12	Z	209/211 (99%)	206 (99%)	3 (1%)	0	100	100
13	M	213/240 (89%)	206 (97%)	7 (3%)	0	100	100
13	a	213/240 (89%)	206 (97%)	7 (3%)	0	100	100
14	N	253/265 (96%)	249 (98%)	4 (2%)	0	100	100
14	b	253/265 (96%)	249 (98%)	4 (2%)	0	100	100
All	All	6426/6706 (96%)	6296 (98%)	130 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	224/231 (97%)	222 (99%)	2 (1%)	70	81
1	O	224/231 (97%)	222 (99%)	2 (1%)	70	81
2	B	203/205 (99%)	202 (100%)	1 (0%)	81	89
2	P	203/205 (99%)	201 (99%)	2 (1%)	68	79
3	C	209/213 (98%)	207 (99%)	2 (1%)	68	79
3	Q	209/213 (98%)	204 (98%)	5 (2%)	43	55
4	D	202/207 (98%)	201 (100%)	1 (0%)	81	89
4	R	202/207 (98%)	201 (100%)	1 (0%)	81	89
5	E	209/223 (94%)	209 (100%)	0	100	100
5	S	209/223 (94%)	208 (100%)	1 (0%)	81	89
6	F	213/226 (94%)	211 (99%)	2 (1%)	70	81
6	T	213/226 (94%)	211 (99%)	2 (1%)	70	81
7	G	226/229 (99%)	226 (100%)	0	100	100
7	U	226/229 (99%)	225 (100%)	1 (0%)	84	91
8	H	225/231 (97%)	225 (100%)	0	100	100
8	V	225/231 (97%)	225 (100%)	0	100	100
9	I	188/194 (97%)	188 (100%)	0	100	100
9	W	188/194 (97%)	188 (100%)	0	100	100
10	J	184/191 (96%)	183 (100%)	1 (0%)	81	89
10	X	184/191 (96%)	183 (100%)	1 (0%)	81	89
11	K	174/174 (100%)	173 (99%)	1 (1%)	78	87
11	Y	174/174 (100%)	173 (99%)	1 (1%)	78	87
12	L	176/176 (100%)	173 (98%)	3 (2%)	53	66
12	Z	176/176 (100%)	174 (99%)	2 (1%)	65	77
13	M	192/216 (89%)	190 (99%)	2 (1%)	68	79
13	a	192/216 (89%)	190 (99%)	2 (1%)	68	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	231/239 (97%)	230 (100%)	1 (0%)	84	91
14	b	231/239 (97%)	231 (100%)	0	100	100
All	All	5712/5910 (97%)	5676 (99%)	36 (1%)	76	87

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	SER
1	A	256	ILE
2	B	198	SER
3	C	73	PHE
3	C	223	TYR
4	D	230	ASP
6	F	160	PHE
6	F	225	SER
10	J	151	VAL
11	K	7	LEU
12	L	97	CYS
12	L	182	ASN
12	L	201	THR
13	M	65	THR
13	M	136	ASN
14	N	106	ILE
1	O	110	SER
1	O	256	ILE
2	P	172	SER
2	P	198	SER
3	Q	11	THR
3	Q	73	PHE
3	Q	149	HIS
3	Q	223	TYR
3	Q	238	LEU
4	R	230	ASP
5	S	65	GLU
6	T	160	PHE
6	T	225	SER
7	U	179	ILE
10	X	151	VAL
11	Y	25	ILE
12	Z	97	CYS
12	Z	201	THR

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Mol	Chain	Res	Type
13	a	65	THR
13	a	136	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	59	GLN
2	B	224	GLN
3	C	149	HIS
4	D	77	ASN
4	D	92	GLN
4	D	97	ASN
5	E	43	ASN
6	F	60	GLN
6	F	100	ASN
7	G	188	ASN
8	H	38	ASN
8	H	163	GLN
8	H	168	ASN
9	I	71	ASN
9	I	106	ASN
9	I	185	HIS
10	J	171	GLN
11	K	40	HIS
11	K	72	ASN
12	L	70	ASN
12	L	71	ASN
13	M	231	ASN
14	N	112	ASN
1	O	46	ASN
1	O	77	ASN
2	P	148	GLN
2	P	207	ASN
2	P	224	GLN
4	R	92	GLN
4	R	179	ASN
6	T	60	GLN
6	T	100	ASN
6	T	138	HIS
6	T	218	HIS
7	U	188	ASN

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Mol	Chain	Res	Type
8	V	38	ASN
8	V	104	ASN
8	V	163	GLN
8	V	168	ASN
9	W	35	HIS
9	W	185	HIS
11	Y	72	ASN
12	Z	70	ASN
12	Z	71	ASN
13	a	151	ASN
13	a	186	GLN
14	b	112	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](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
15	YRE	W	301	-	47,47,47	0.94	1 (2%)	62,63,63	1.33	5 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
15	YRE	Z	301	-	47,47,47	0.87	2 (4%)	62,63,63	1.14	4 (6%)
15	YRE	L	301	-	47,47,47	0.88	2 (4%)	62,63,63	1.14	4 (6%)
15	YRE	I	301	-	47,47,47	0.90	1 (2%)	62,63,63	1.30	6 (9%)
15	YRE	C	301	-	47,47,47	0.91	2 (4%)	62,63,63	1.50	7 (11%)
15	YRE	Q	301	-	47,47,47	0.92	2 (4%)	62,63,63	1.38	6 (9%)

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
15	YRE	W	301	-	-	15/45/70/70	0/4/5/5
15	YRE	Z	301	-	-	4/45/70/70	0/4/5/5
15	YRE	L	301	-	-	4/45/70/70	0/4/5/5
15	YRE	I	301	-	-	11/45/70/70	0/4/5/5
15	YRE	C	301	-	-	12/45/70/70	0/4/5/5
15	YRE	Q	301	-	-	12/45/70/70	0/4/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	Q	301	YRE	C30-N29	3.46	1.39	1.35
15	W	301	YRE	C17-C16	3.26	1.60	1.54
15	C	301	YRE	C30-N29	2.99	1.39	1.35
15	I	301	YRE	C17-C16	2.80	1.59	1.54
15	Z	301	YRE	C30-N29	2.57	1.38	1.35
15	L	301	YRE	C30-N29	2.48	1.38	1.35
15	Q	301	YRE	C05-N06	2.30	1.39	1.34
15	C	301	YRE	C05-N06	2.26	1.38	1.34
15	L	301	YRE	C17-C18	-2.04	1.46	1.51
15	Z	301	YRE	C17-C18	-2.00	1.46	1.51

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	Q	301	YRE	C16-N29-C30	5.90	129.93	123.08
15	C	301	YRE	C16-N29-C30	5.63	129.61	123.08
15	Z	301	YRE	C16-N29-C30	5.62	129.60	123.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	L	301	YRE	C16-N29-C30	5.57	129.54	123.08
15	W	301	YRE	C33-N29-C30	-5.04	110.28	113.47
15	C	301	YRE	C33-N29-C30	-4.97	110.32	113.47
15	I	301	YRE	C33-N29-C30	-4.97	110.32	113.47
15	Q	301	YRE	C33-N29-C30	-4.95	110.34	113.47
15	I	301	YRE	C17-C16-N29	4.37	117.16	112.65
15	C	301	YRE	C42-C43-N38	3.96	116.14	110.12
15	W	301	YRE	C17-C16-N29	3.74	116.51	112.65
15	C	301	YRE	C40-C39-N38	3.68	115.72	110.12
15	W	301	YRE	C18-C17-C16	3.53	121.50	113.31
15	I	301	YRE	C18-C17-C16	3.35	121.09	113.31
15	W	301	YRE	C16-N29-C30	3.29	126.90	123.08
15	Z	301	YRE	C33-N29-C30	-3.07	111.53	113.47
15	L	301	YRE	C33-N29-C30	-3.03	111.55	113.47
15	I	301	YRE	C16-N29-C30	2.99	126.55	123.08
15	Q	301	YRE	O35-C15-N14	-2.83	117.89	122.96
15	C	301	YRE	O35-C15-N14	-2.81	117.92	122.96
15	Q	301	YRE	C42-C43-N38	2.65	114.15	110.12
15	Z	301	YRE	C33-N29-C16	-2.53	118.87	123.09
15	L	301	YRE	C33-N29-C16	-2.51	118.90	123.09
15	W	301	YRE	C40-C39-N38	2.43	113.81	110.12
15	Q	301	YRE	C18-C17-C16	2.23	118.49	113.31
15	C	301	YRE	C18-C17-C16	2.20	118.41	113.31
15	I	301	YRE	C36-C13-N14	-2.13	106.69	110.91
15	Z	301	YRE	O35-C15-N14	-2.13	119.15	122.96
15	C	301	YRE	C13-N14-C15	-2.11	117.12	121.65
15	I	301	YRE	O35-C15-N14	-2.10	119.19	122.96
15	L	301	YRE	O35-C15-N14	-2.10	119.20	122.96
15	Q	301	YRE	C33-N29-C16	-2.02	119.73	123.09

There are no chirality outliers.

All (58) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
15	C	301	YRE	C02-C13-C36-C37
15	L	301	YRE	C02-C13-C36-C37
15	L	301	YRE	N14-C13-C36-C37
15	Q	301	YRE	C02-C13-C36-C37
15	W	301	YRE	C23-C24-C25-C04
15	W	301	YRE	C13-C36-C37-N38
15	Z	301	YRE	C02-C13-C36-C37
15	Z	301	YRE	N14-C13-C36-C37

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Mol	Chain	Res	Type	Atoms
15	C	301	YRE	N03-C04-C25-C24
15	I	301	YRE	C36-C37-N38-C43
15	C	301	YRE	C05-C04-C25-C24
15	Q	301	YRE	C05-C04-C25-C24
15	Q	301	YRE	N03-C04-C25-C24
15	Q	301	YRE	O21-C22-C23-C24
15	C	301	YRE	O21-C22-C23-C24
15	I	301	YRE	C23-C24-C25-C04
15	L	301	YRE	C23-C24-C25-C04
15	Z	301	YRE	C23-C24-C25-C04
15	C	301	YRE	N14-C13-C36-C37
15	Q	301	YRE	N14-C13-C36-C37
15	W	301	YRE	N14-C13-C36-C37
15	Q	301	YRE	C36-C37-N38-C39
15	W	301	YRE	C22-C23-C24-C25
15	I	301	YRE	C36-C37-N38-C39
15	Q	301	YRE	C19-C20-O21-C22
15	C	301	YRE	C19-C20-O21-C22
15	W	301	YRE	O21-C22-C23-C24
15	W	301	YRE	C36-C37-N38-C43
15	I	301	YRE	C23-C22-O21-C20
15	C	301	YRE	C36-C37-N38-C39
15	Q	301	YRE	C26-C20-O21-C22
15	W	301	YRE	C23-C22-O21-C20
15	C	301	YRE	C26-C20-O21-C22
15	L	301	YRE	C36-C37-N38-C43
15	Z	301	YRE	C36-C37-N38-C43
15	I	301	YRE	C22-C23-C24-C25
15	W	301	YRE	N03-C04-C25-C24
15	I	301	YRE	O35-C15-C16-C17
15	C	301	YRE	C22-C23-C24-C25
15	Q	301	YRE	C23-C22-O21-C20
15	C	301	YRE	C23-C22-O21-C20
15	Q	301	YRE	C22-C23-C24-C25
15	I	301	YRE	O21-C22-C23-C24
15	W	301	YRE	C05-C04-C25-C24
15	C	301	YRE	N03-C04-C05-N06
15	I	301	YRE	C11-C07-N06-C05
15	C	301	YRE	N03-C04-C05-O12
15	W	301	YRE	O35-C15-C16-C17
15	Q	301	YRE	N03-C04-C05-N06
15	Q	301	YRE	N03-C04-C05-O12

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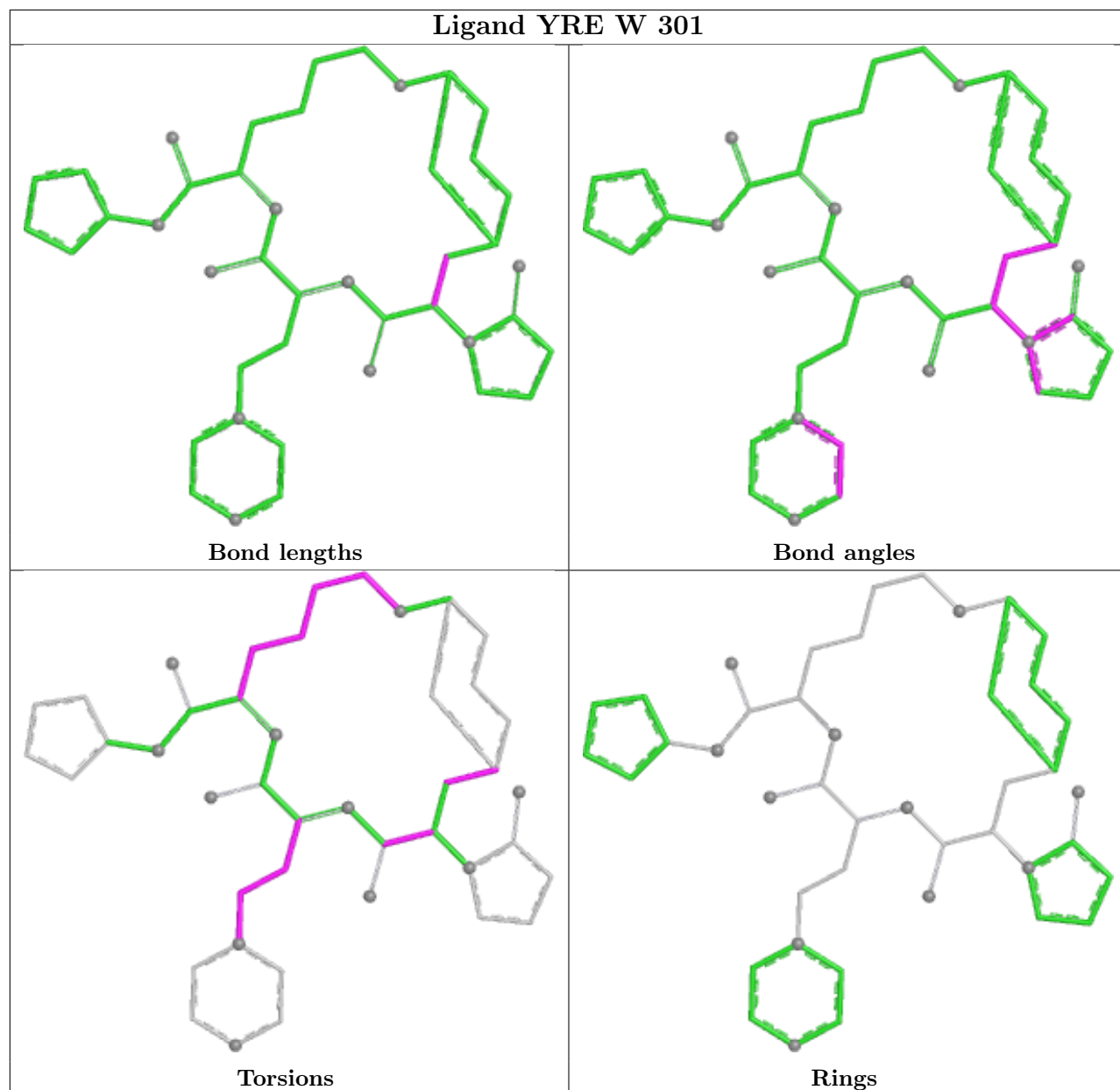
Continued from previous page...

Mol	Chain	Res	Type	Atoms
15	W	301	YRE	C16-C17-C18-C19
15	I	301	YRE	N14-C15-C16-C17
15	I	301	YRE	N14-C15-C16-N29
15	I	301	YRE	O35-C15-C16-N29
15	W	301	YRE	O35-C15-C16-N29
15	W	301	YRE	C16-C17-C18-C28
15	W	301	YRE	N14-C15-C16-C17
15	W	301	YRE	N14-C15-C16-N29

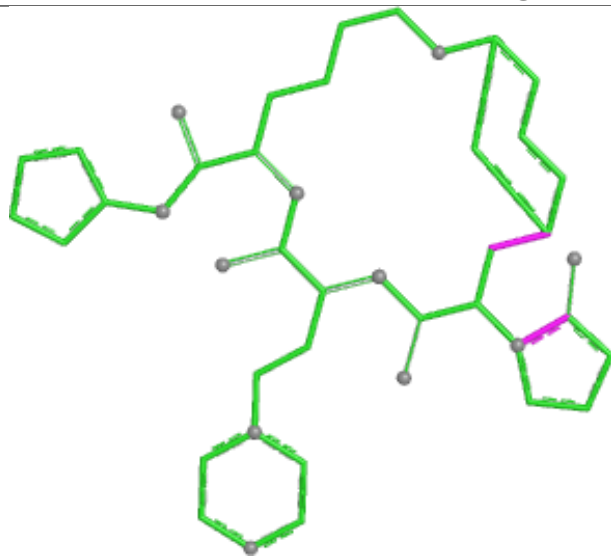
There are no ring outliers.

No monomer is involved in short contacts.

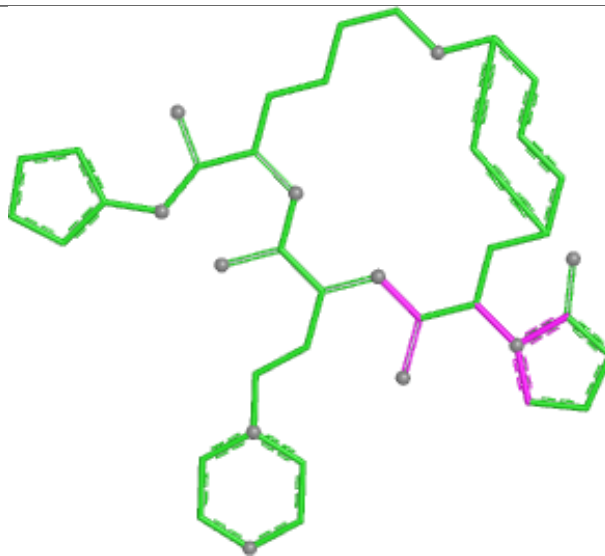
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



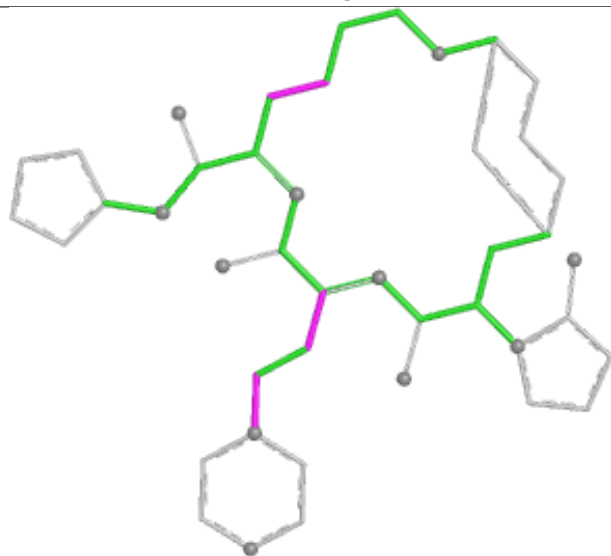
Ligand YRE Z 301



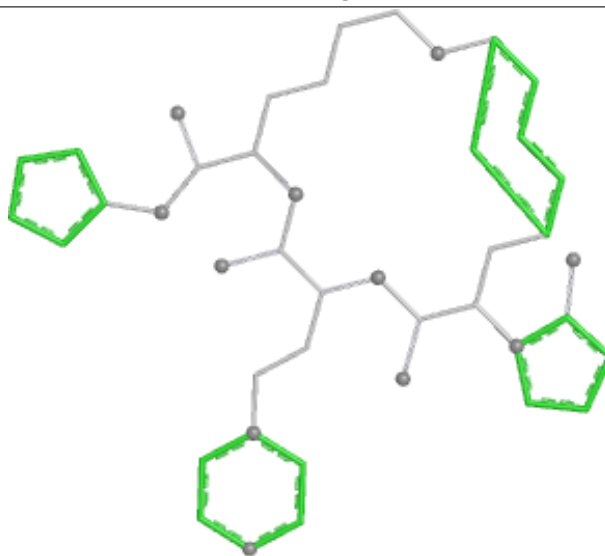
Bond lengths



Bond angles

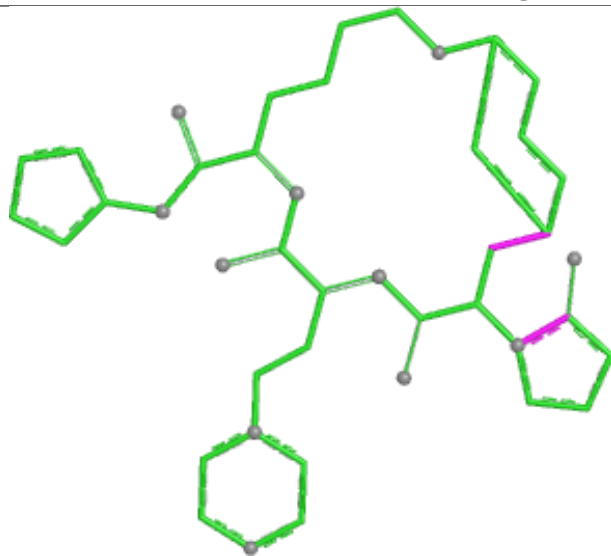


Torsions

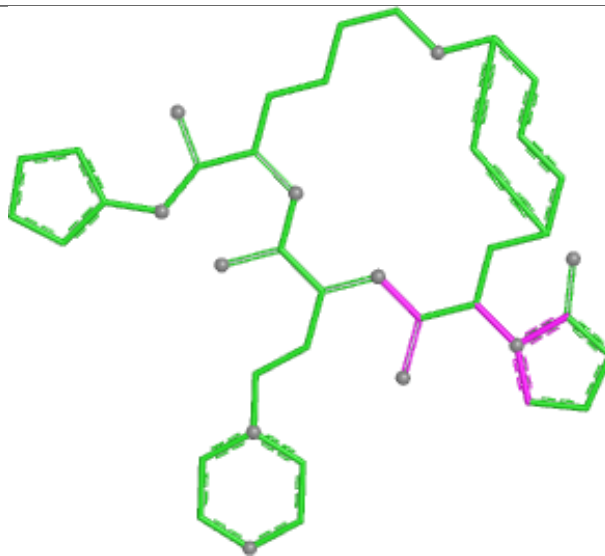


Rings

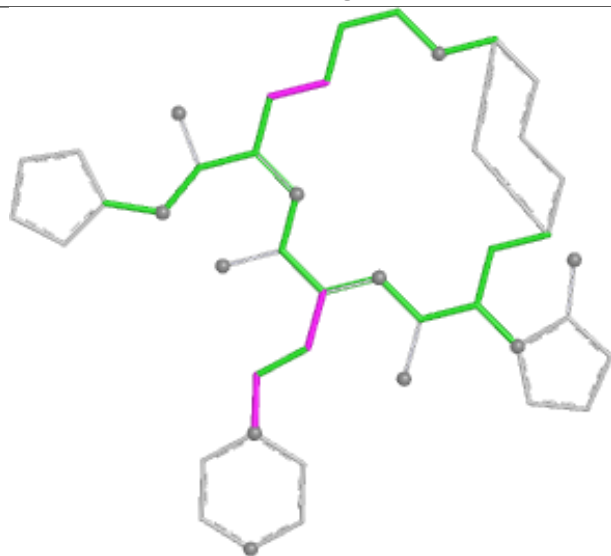
Ligand YRE L 301



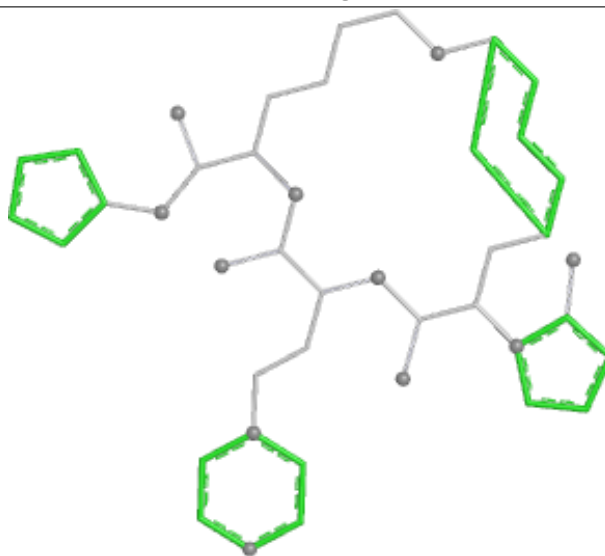
Bond lengths



Bond angles

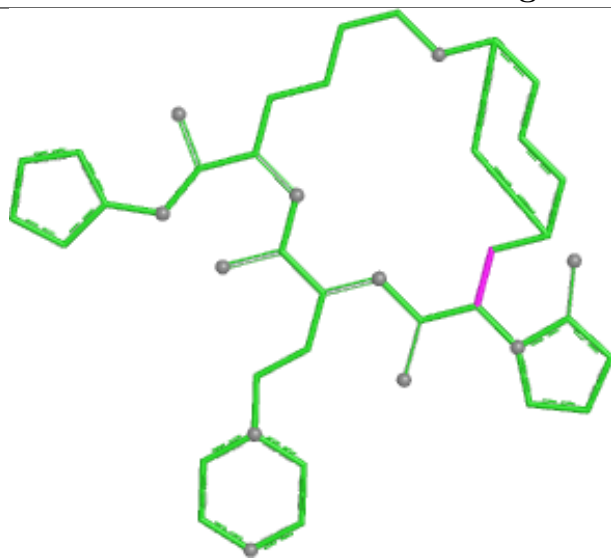


Torsions

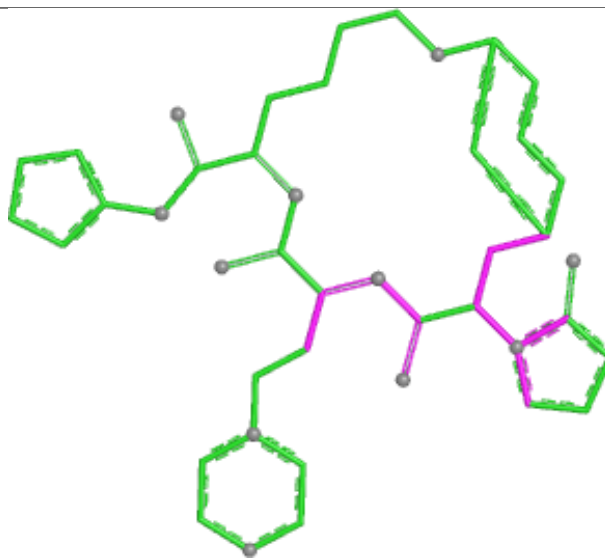


Rings

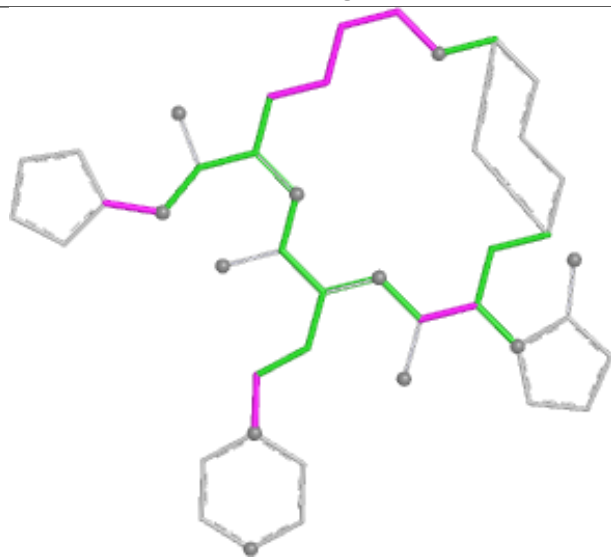
Ligand YRE I 301



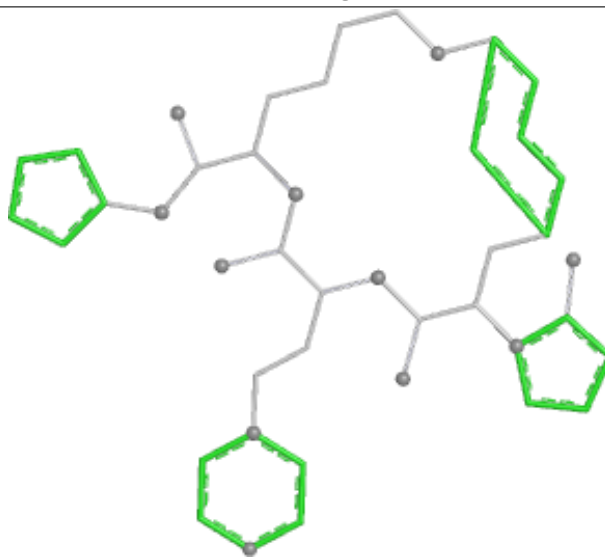
Bond lengths



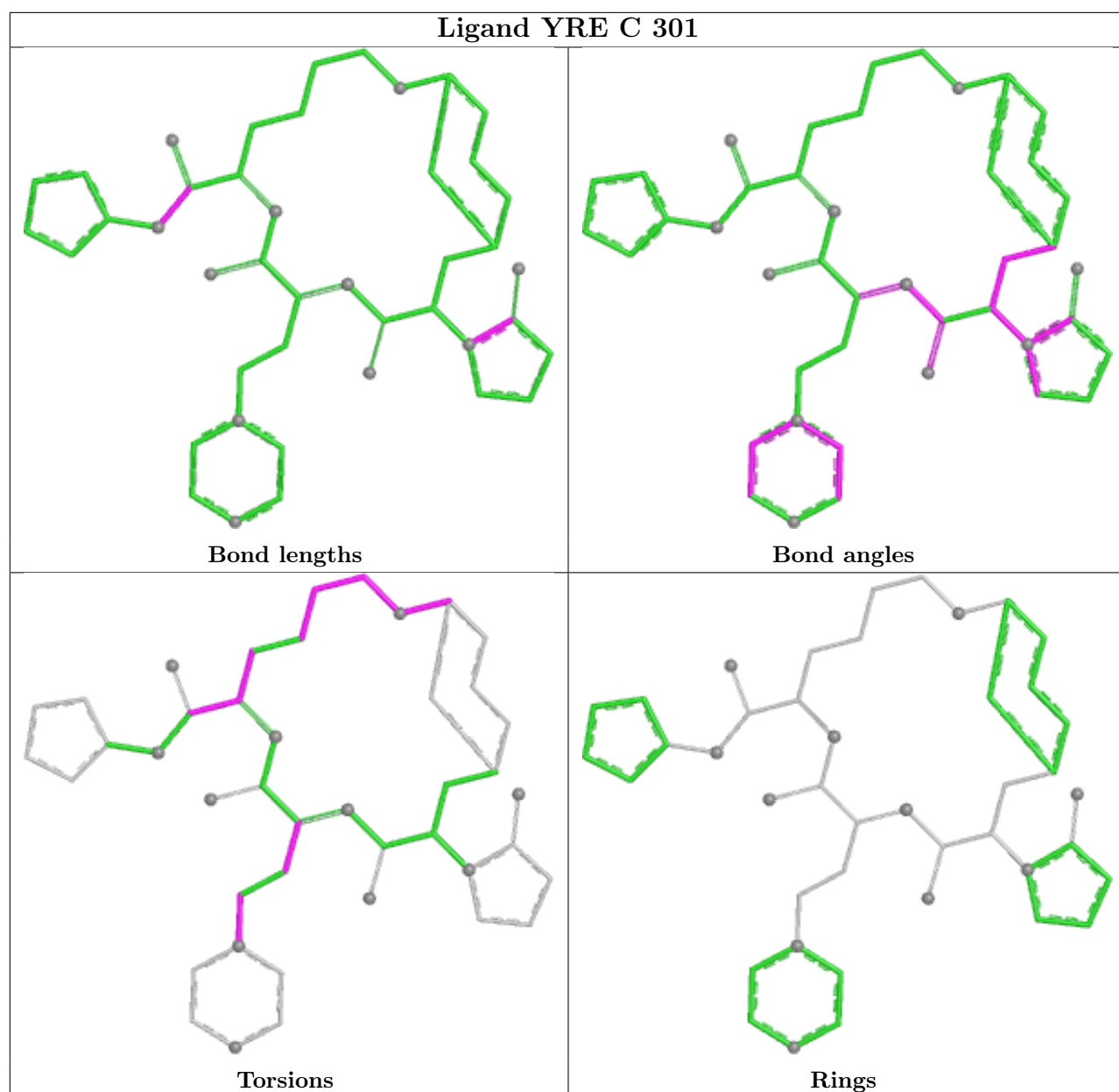
Bond angles

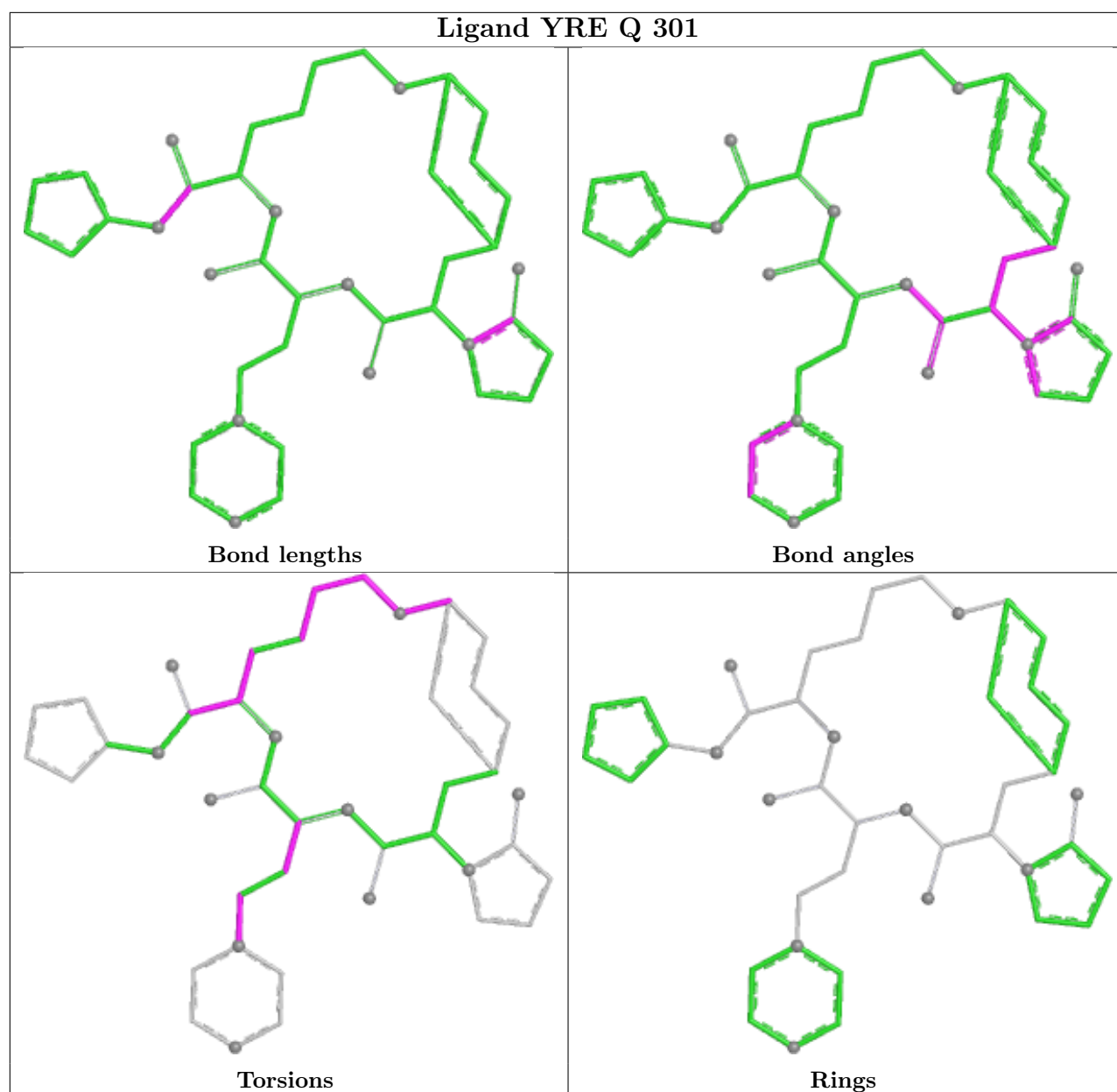


Torsions



Rings





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.

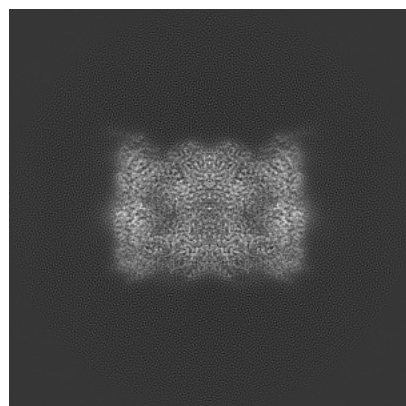
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29764. These allow visual inspection of the internal detail of the map and identification of artifacts.

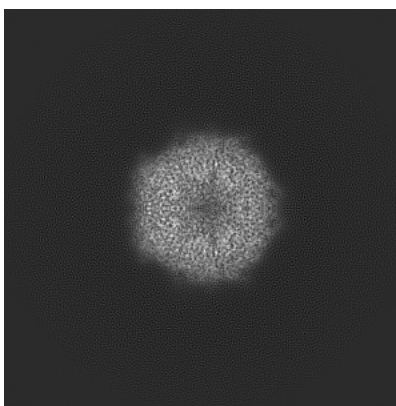
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

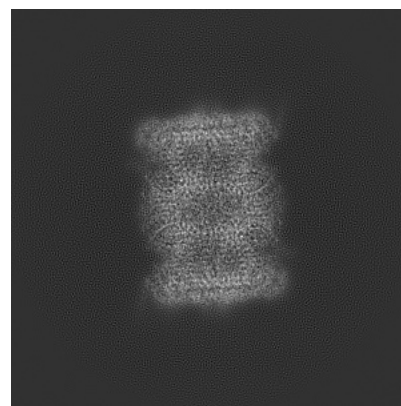
6.1.1 Primary map



X

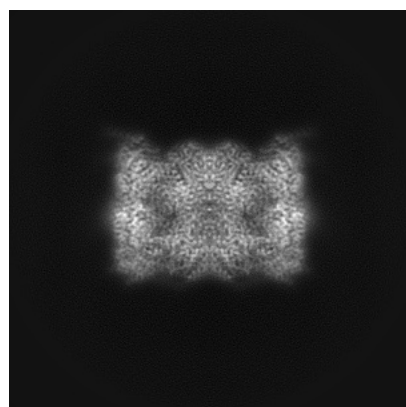


Y

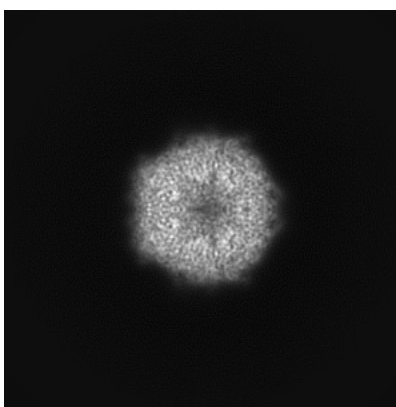


Z

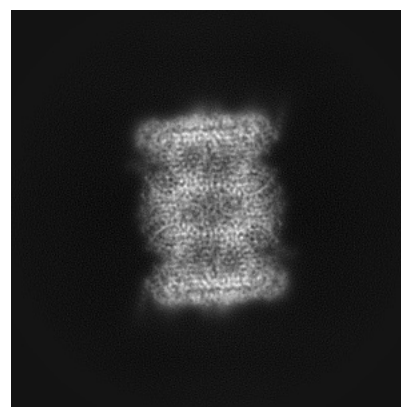
6.1.2 Raw map



X



Y

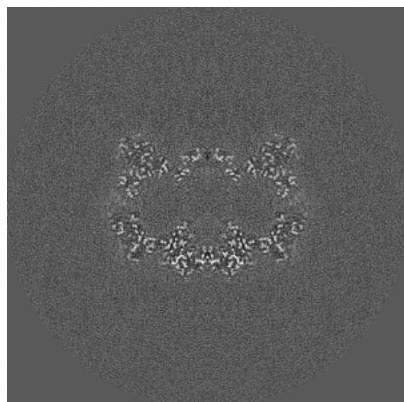


Z

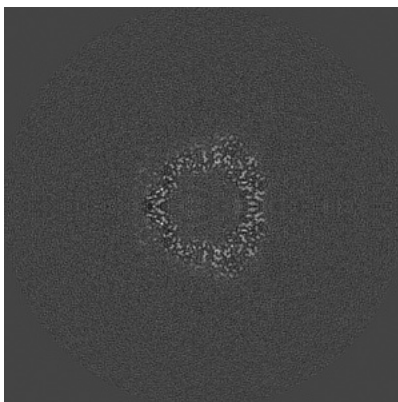
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

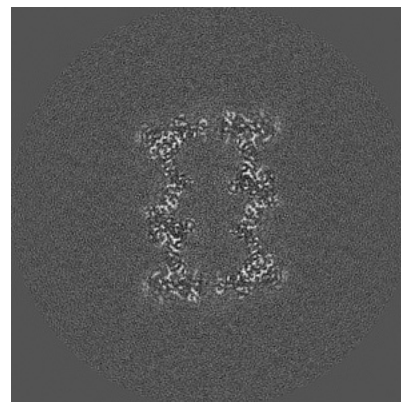
6.2.1 Primary map



X Index: 192

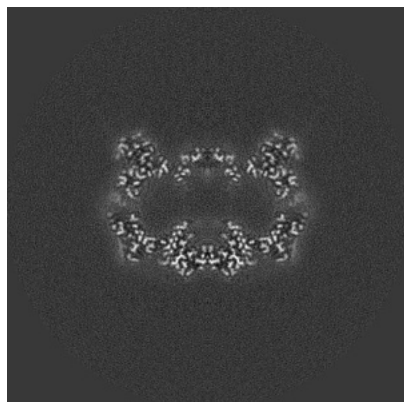


Y Index: 192

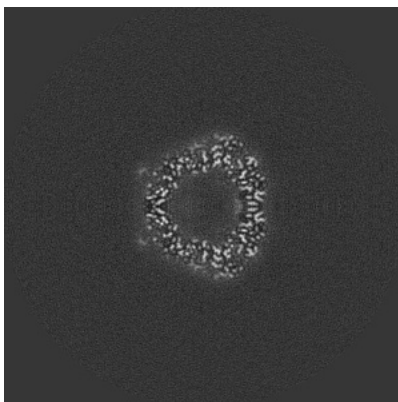


Z Index: 192

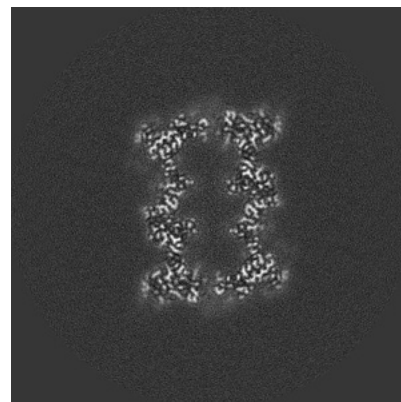
6.2.2 Raw map



X Index: 192



Y Index: 192

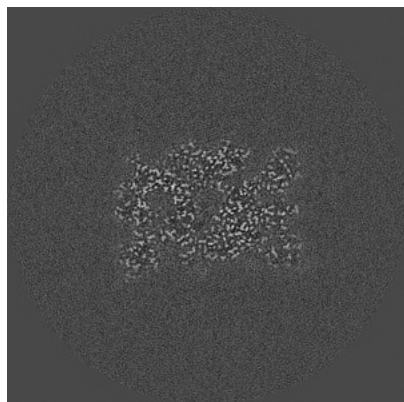


Z Index: 192

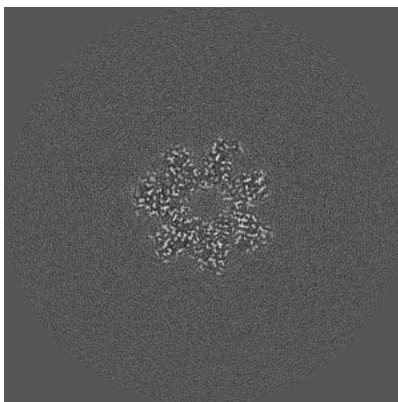
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

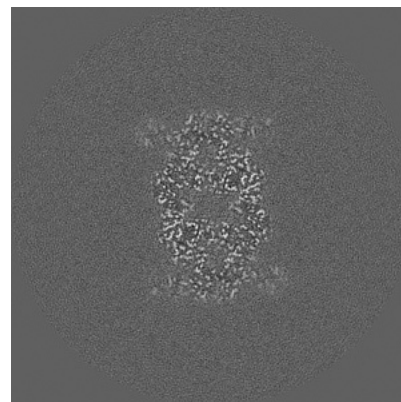
6.3.1 Primary map



X Index: 223

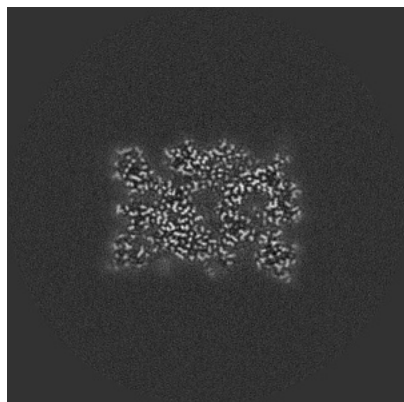


Y Index: 169

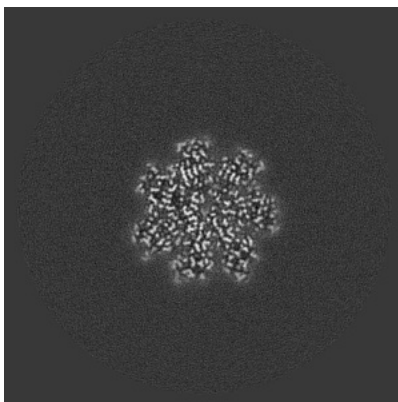


Z Index: 165

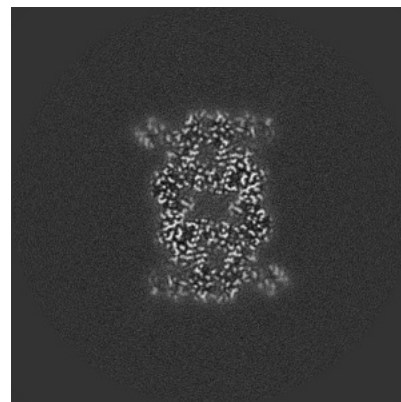
6.3.2 Raw map



X Index: 161



Y Index: 263

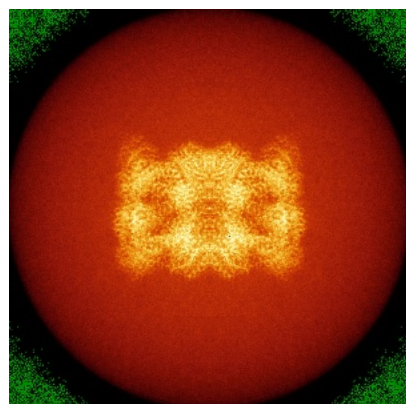


Z Index: 165

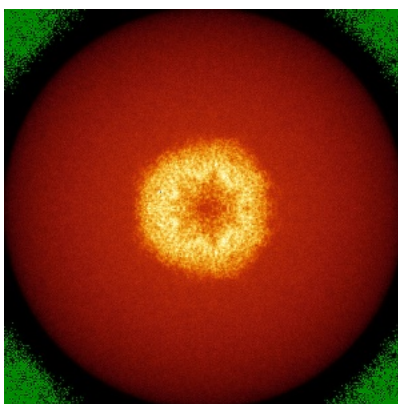
The images above show the largest variance slices of the map in three orthogonal directions.

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

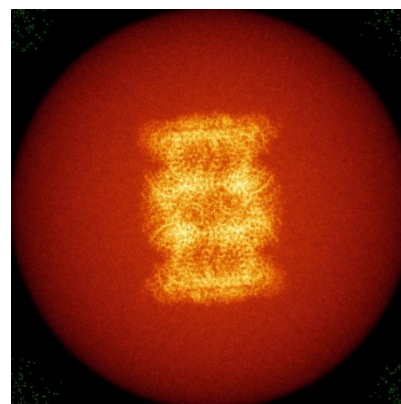
6.4.1 Primary map



X

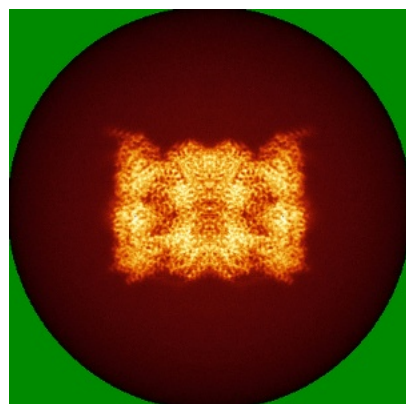


Y

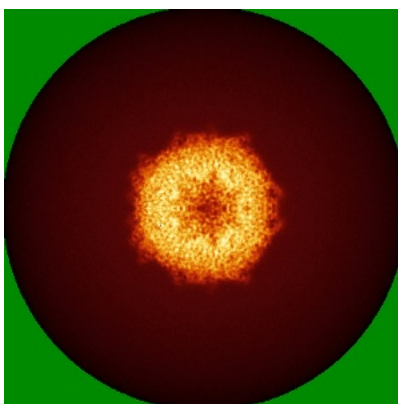


Z

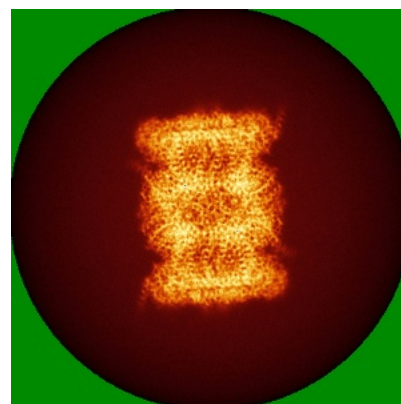
6.4.2 Raw map



X



Y

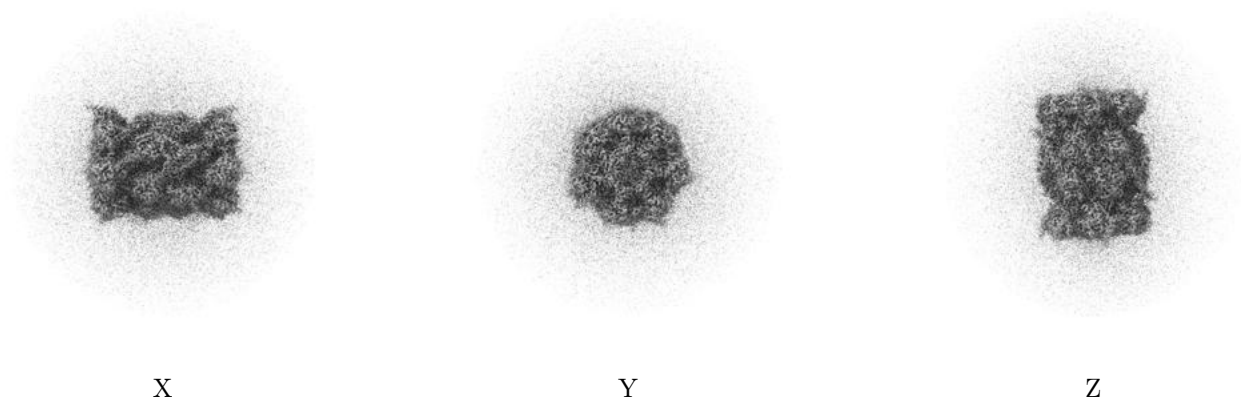


Z

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.

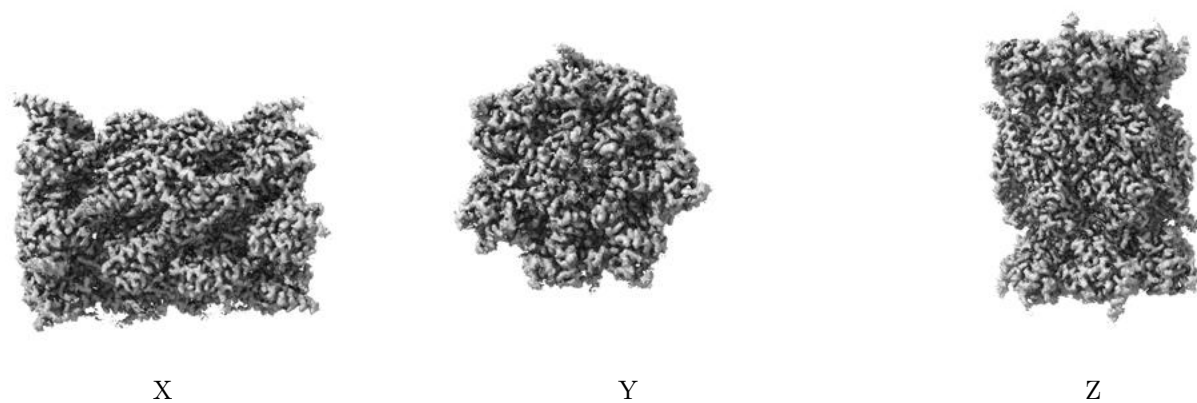
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.025. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

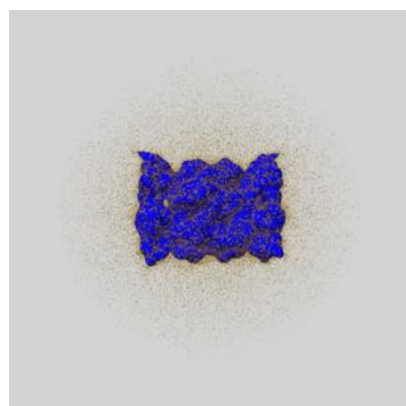
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

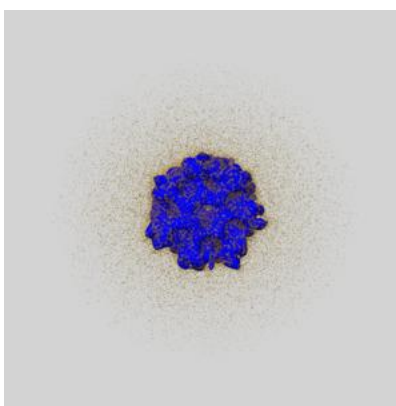
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

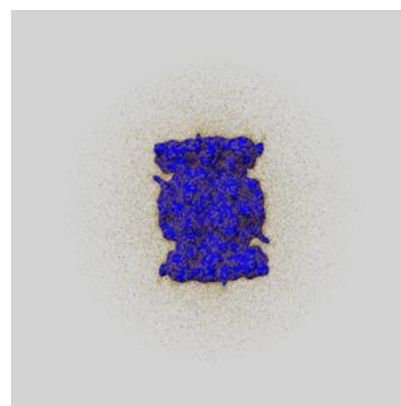
6.6.1 emd_29764_msk_1.map [i](#)



X



Y

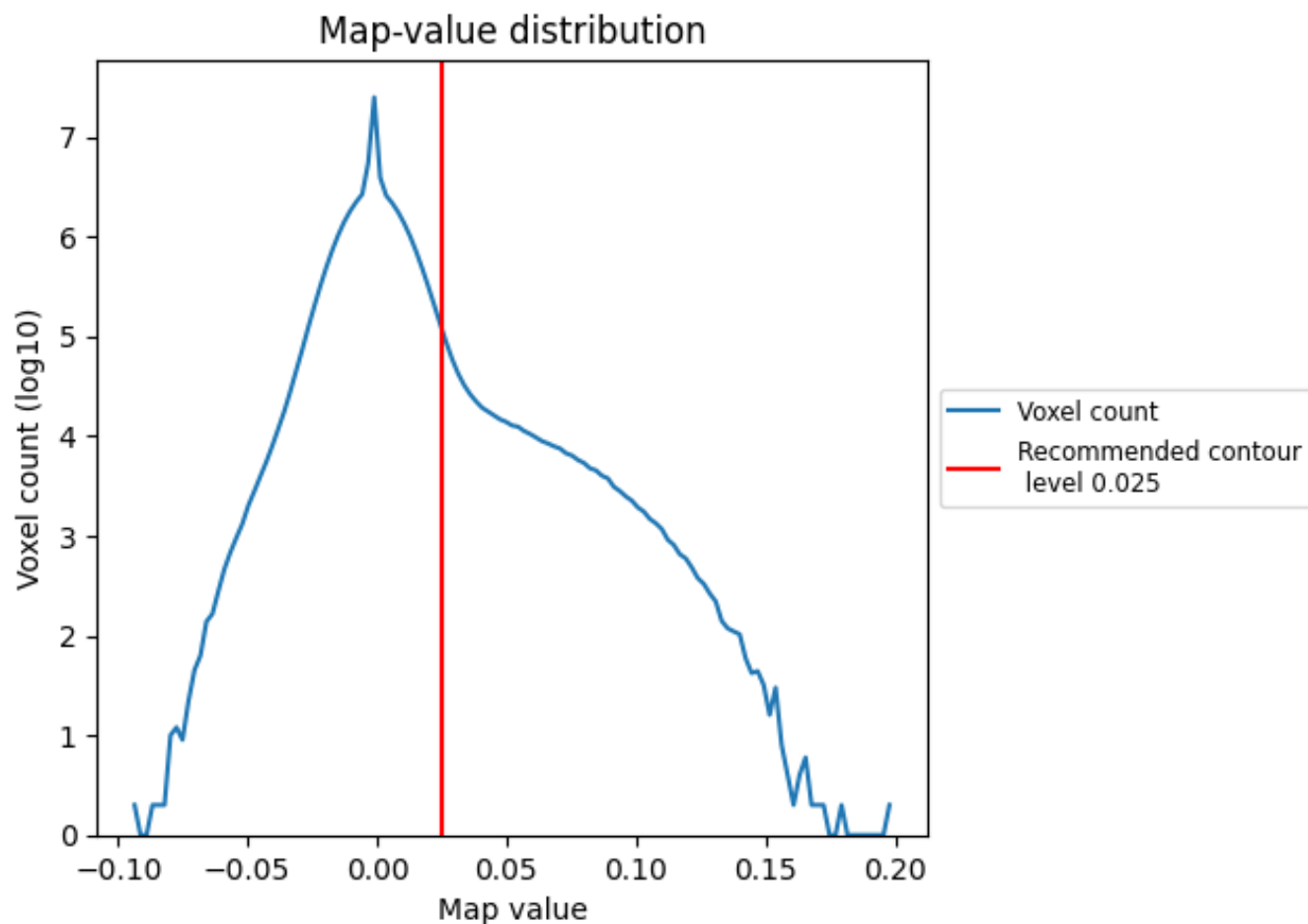


Z

7 Map analysis [i](#)

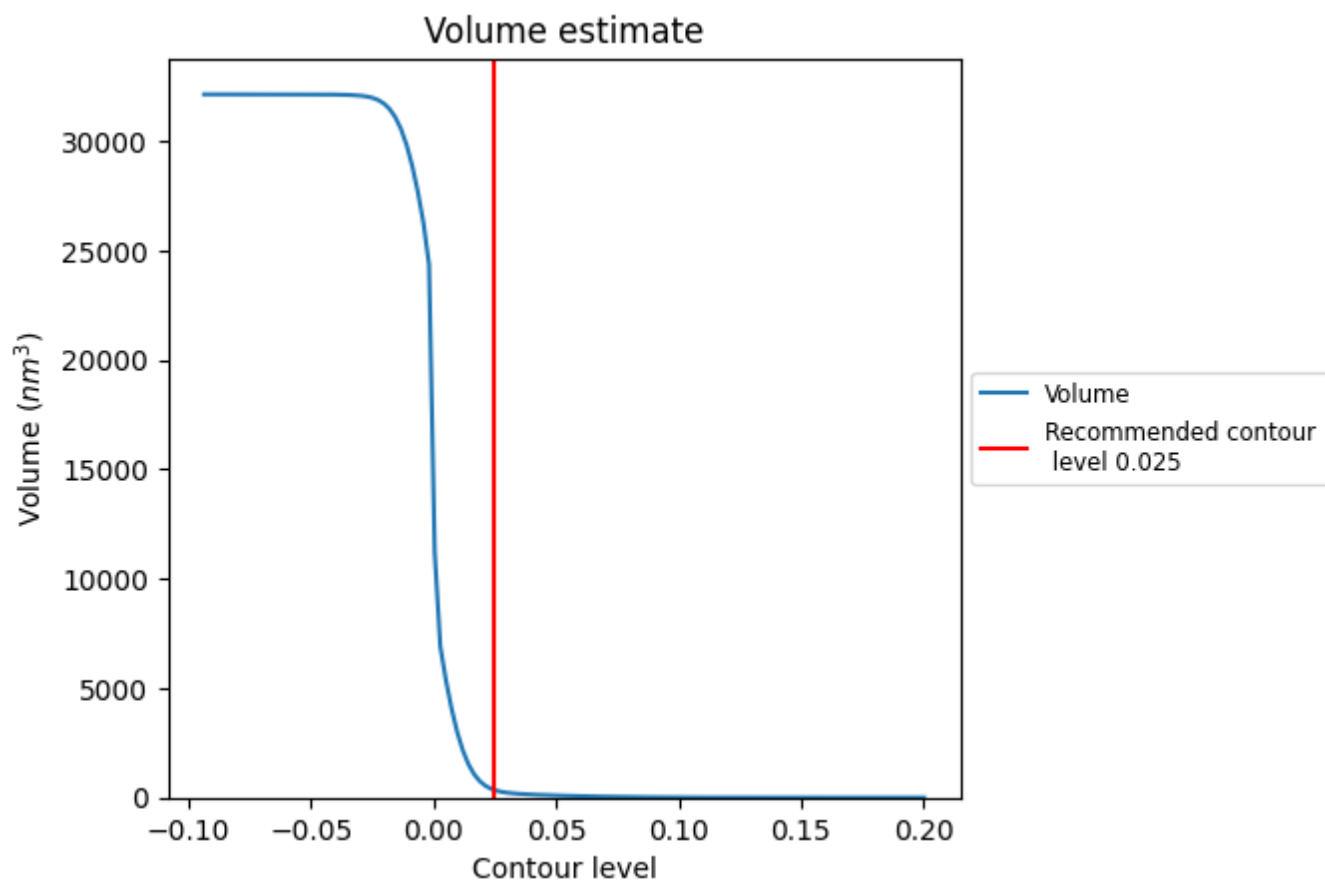
This section contains the results of statistical analysis of the map.

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.

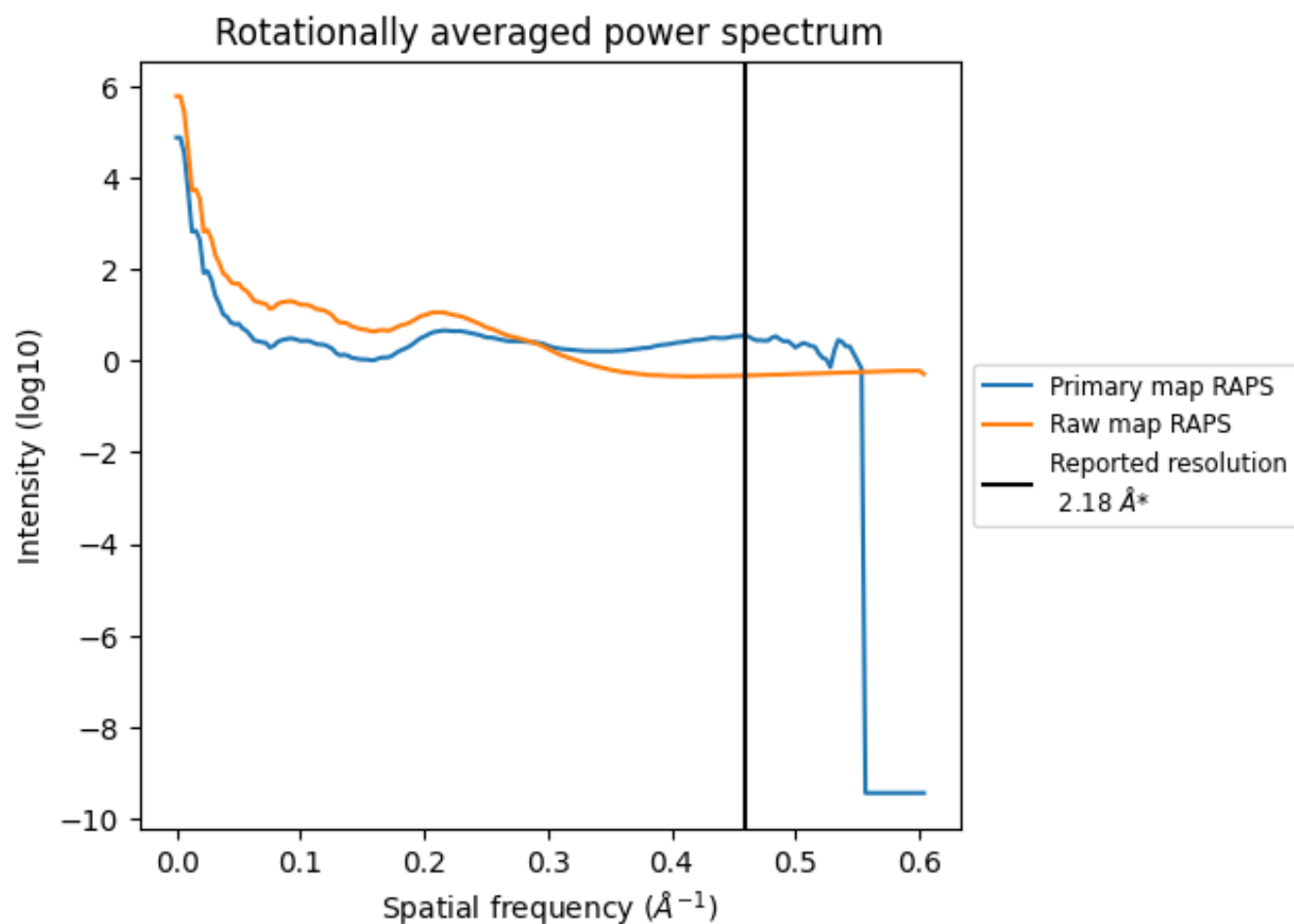
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 343 nm³; this corresponds to an approximate mass of 310 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 ⓘ

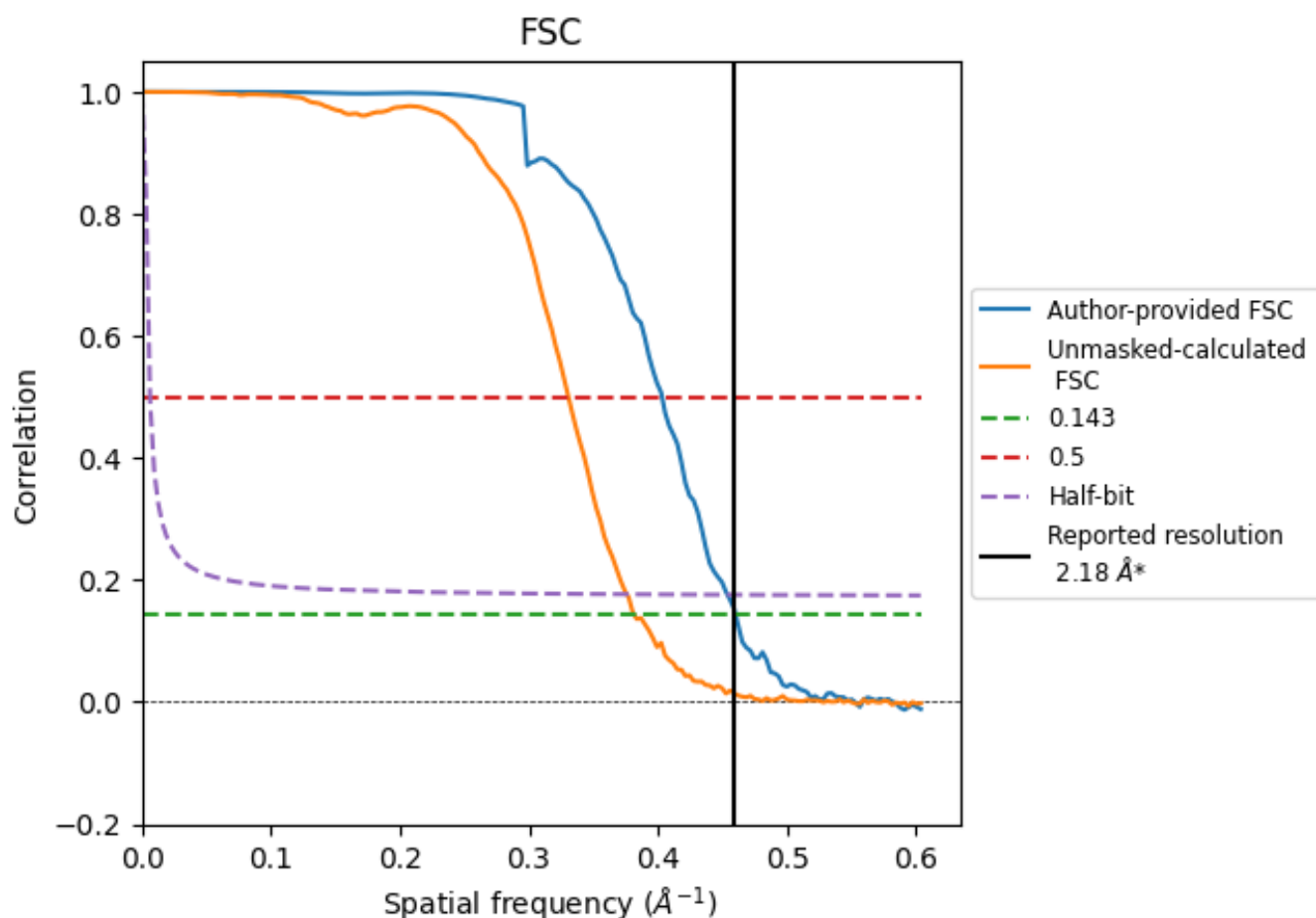


*Reported resolution corresponds to spatial frequency of 0.459 Å⁻¹

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 0.459 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

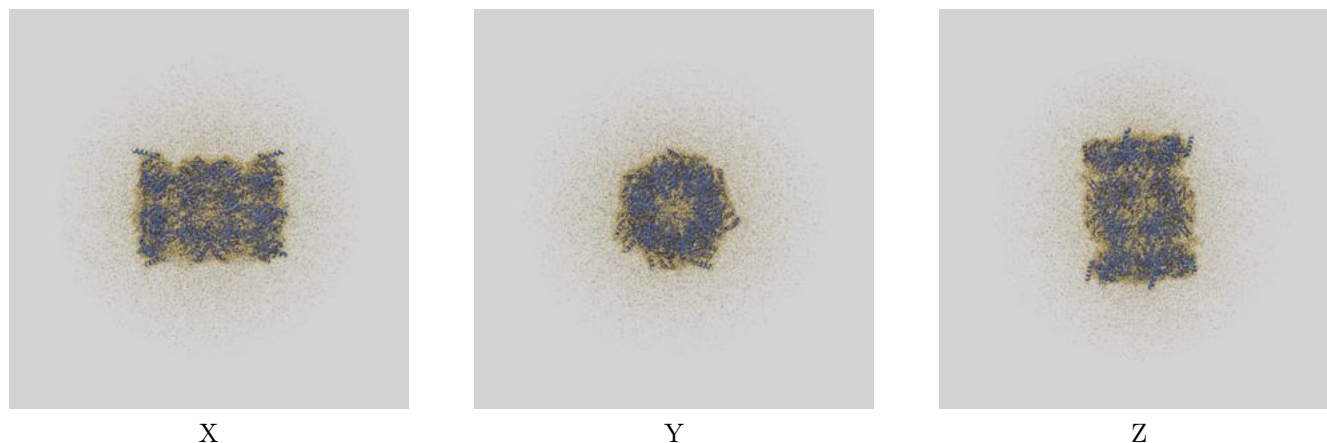
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.18	-	-
Author-provided FSC curve	2.17	2.48	2.20
Unmasked-calculated*	2.62	3.02	2.66

*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 2.62 differs from the reported value 2.18 by more than 10 %

9 Map-model fit [i](#)

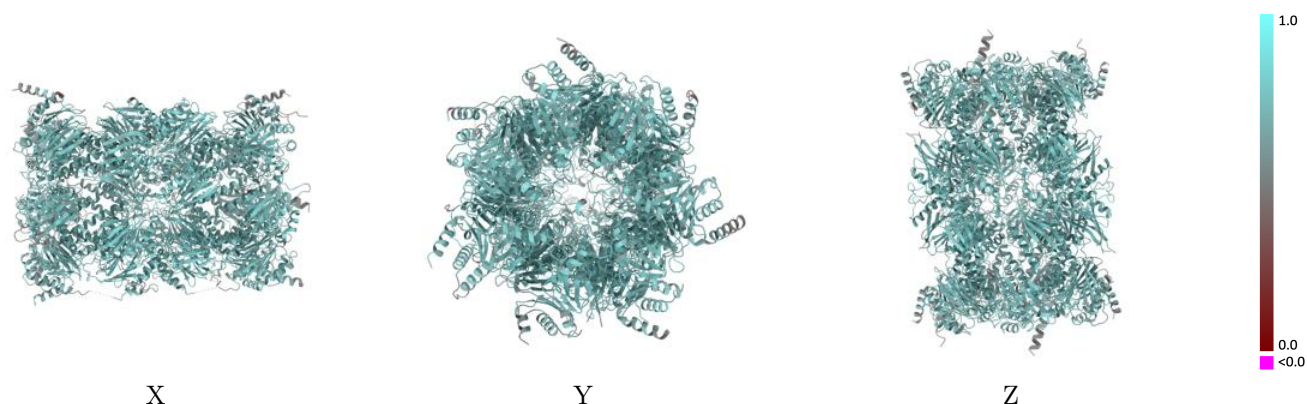
This section contains information regarding the fit between EMDB map EMD-29764 and PDB model 8G6E. Per-residue inclusion information can be found in [section 3](#) on [page 10](#).

9.1 Map-model overlay [i](#)



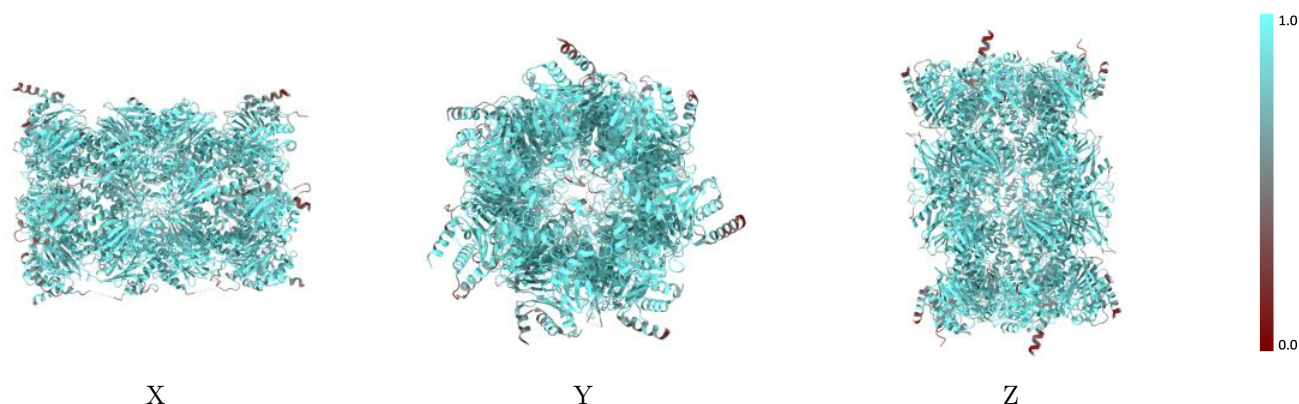
The images above show the 3D surface view of the map at the recommended contour level 0.025 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



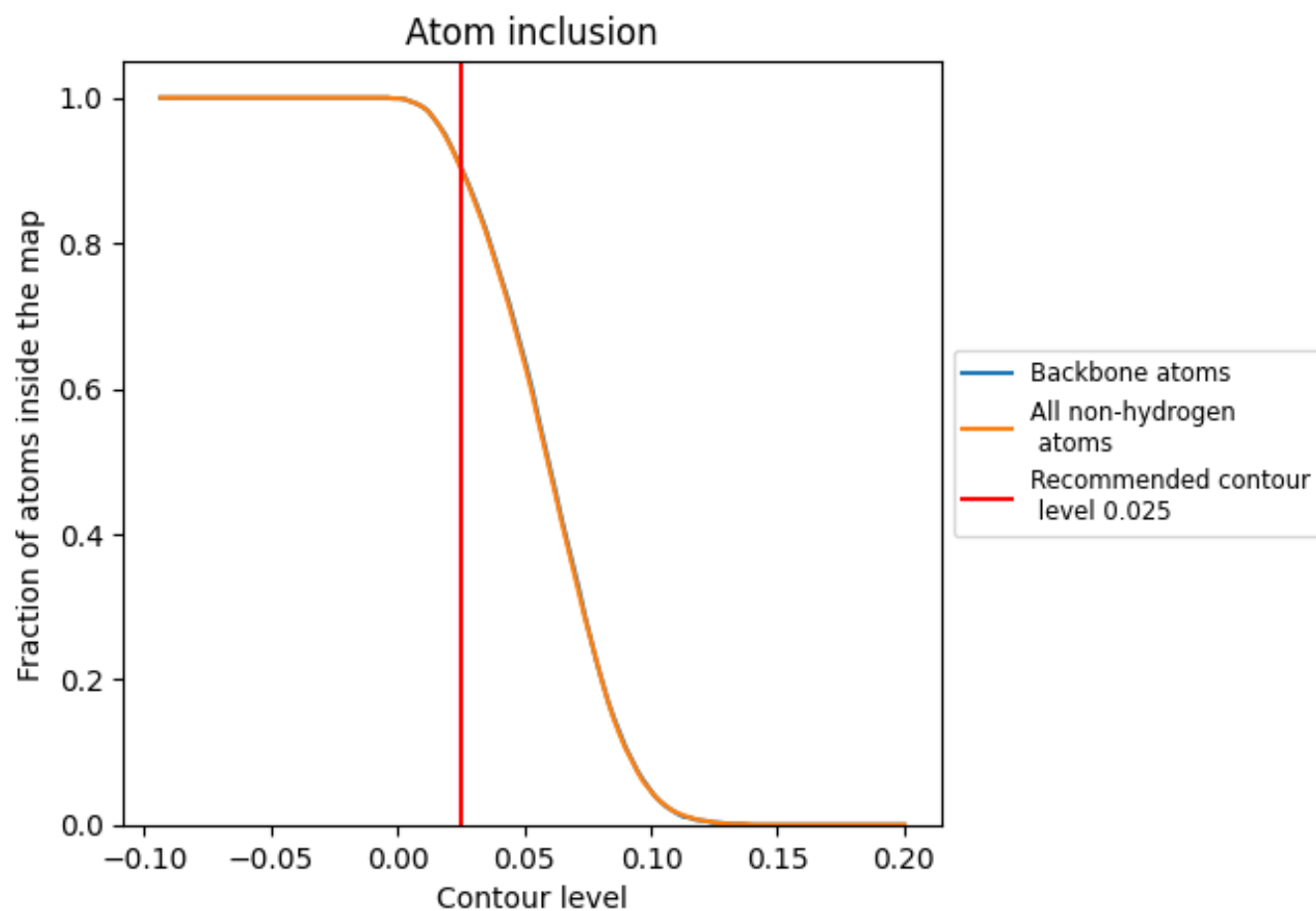
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 90% 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 (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9030	 0.7150
A	 0.8820	 0.7020
B	 0.8420	 0.6740
C	 0.8790	 0.6960
D	 0.8610	 0.6880
E	 0.8490	 0.6760
F	 0.8760	 0.6960
G	 0.9140	 0.7130
H	 0.9170	 0.7300
I	 0.9500	 0.7390
J	 0.9460	 0.7470
K	 0.9460	 0.7360
L	 0.9620	 0.7480
M	 0.9380	 0.7330
N	 0.9350	 0.7380
O	 0.8800	 0.7020
P	 0.8430	 0.6740
Q	 0.8830	 0.6990
R	 0.8580	 0.6890
S	 0.8470	 0.6780
T	 0.8780	 0.6970
U	 0.9150	 0.7140
V	 0.9150	 0.7290
W	 0.9480	 0.7370
X	 0.9450	 0.7450
Y	 0.9480	 0.7360
Z	 0.9610	 0.7480
a	 0.9360	 0.7350
b	 0.9350	 0.7360

