



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 08:14 PM UTC

PDB ID : 6Z2K / pdb_00006z2k
EMDB ID : EMD-11042
Title : The structure of the tetrameric HDAC1/MIDEAS/DNTTIP1 MiDAC deacetylase complex
Authors : Fairall, L.; Saleh, A.; Ragan, T.J.; Millard, C.J.; Savva, C.G.; Schwabe, J.W.R.
Deposited on : 2020-05-16
Resolution : 4.50 Å (reported)
Based on initial models : 5ICN, 4D6K

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:

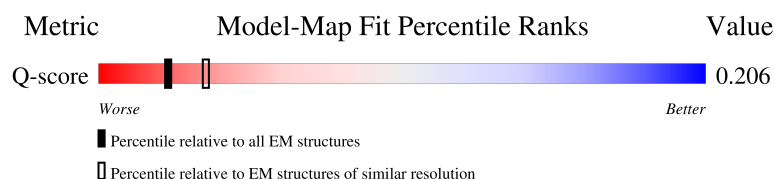
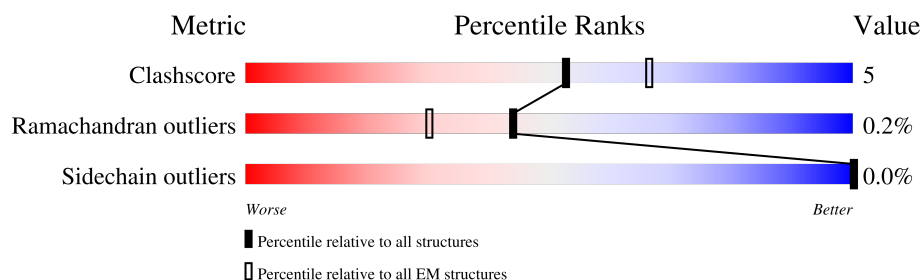
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 4.50 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.



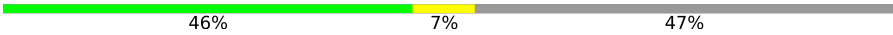
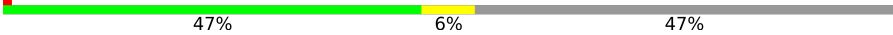
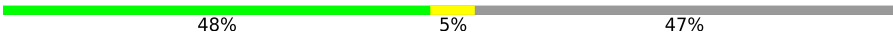
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2937 (4.00 - 5.00)

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	C	482	
1	E	482	
1	I	482	
1	K	482	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	A	130	
2	B	130	
2	G	130	
2	H	130	
3	D	173	
3	F	173	
3	J	173	
3	L	173	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 37867 atoms, of which 18601 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 Histone deacetylase 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	C	369	Total	C	H	N	O	S	0	0
			5802	1892	2838	501	549	22		
1	E	368	Total	C	H	N	O	S	0	0
			5787	1888	2831	499	547	22		
1	K	369	Total	C	H	N	O	S	0	0
			5802	1892	2838	501	549	22		
1	I	369	Total	C	H	N	O	S	0	0
			5802	1892	2838	501	549	22		

- Molecule 2 is a protein called Deoxynucleotidyltransferase terminal-interacting protein 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	A	69	Total	C	H	N	O	S	4	0
			1139	356	573	96	109	5		
2	B	69	Total	C	H	N	O	S	0	0
			1098	346	549	92	107	4		
2	G	69	Total	C	H	N	O	S	4	0
			1139	356	573	96	109	5		
2	H	69	Total	C	H	N	O	S	0	0
			1098	346	549	92	107	4		

- Molecule 3 is a protein called Mitotic deacetylase-associated SANT domain protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	F	157	Total	C	H	N	O	S	0	0
			2505	808	1247	221	225	4		
3	D	157	Total	C	H	N	O	S	0	0
			2505	808	1247	221	225	4		
3	L	157	Total	C	H	N	O	S	0	0
			2505	808	1247	221	225	4		
3	J	157	Total	C	H	N	O	S	0	0
			2505	808	1247	221	225	4		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	715	GLY	-	expression tag	UNP Q6PJG2
F	716	ALA	-	expression tag	UNP Q6PJG2
D	715	GLY	-	expression tag	UNP Q6PJG2
D	716	ALA	-	expression tag	UNP Q6PJG2
L	715	GLY	-	expression tag	UNP Q6PJG2
L	716	ALA	-	expression tag	UNP Q6PJG2
J	715	GLY	-	expression tag	UNP Q6PJG2
J	716	ALA	-	expression tag	UNP Q6PJG2

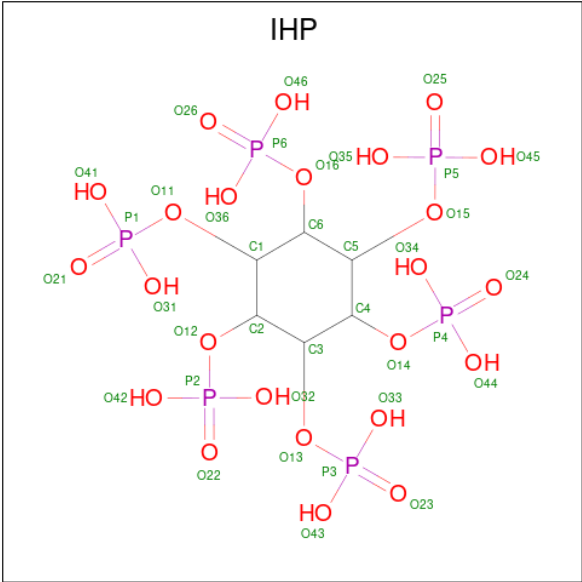
- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	AltConf
4	C	1	Total Zn 1 1	0
4	E	1	Total Zn 1 1	0
4	K	1	Total Zn 1 1	0
4	I	1	Total Zn 1 1	0

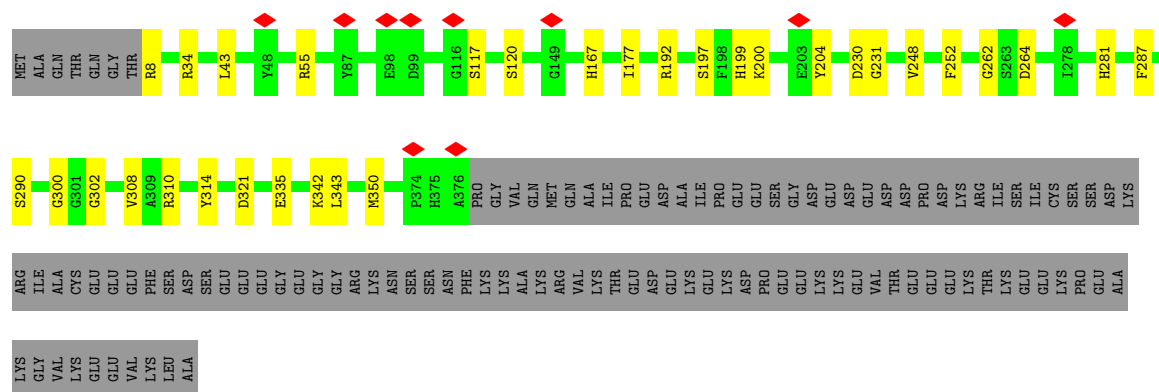
- Molecule 5 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	AltConf
5	C	2	Total K 2 2	0
5	E	2	Total K 2 2	0
5	K	2	Total K 2 2	0
5	I	2	Total K 2 2	0

- Molecule 6 is INOSITOL HEXAKISPHOSPHATE (CCD ID: IHP) (formula: C₆H₁₈O₂₄P₆).

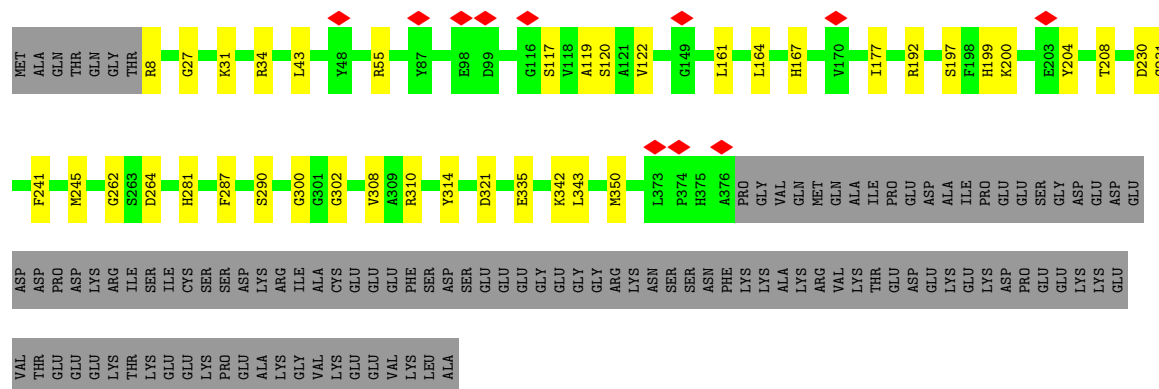


Mol	Chain	Residues	Atoms					AltConf
6	E	1	Total	C	H	O	P	0
			42	6	6	24	6	
6	F	1	Total	C	H	O	P	0
			42	6	6	24	6	
6	D	1	Total	C	H	O	P	0
			42	6	6	24	6	
6	J	1	Total	C	H	O	P	0
			42	6	6	24	6	



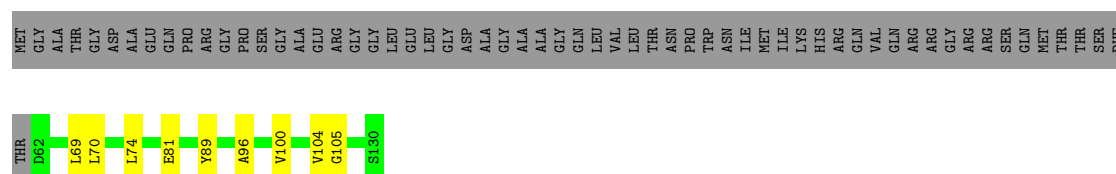
• Molecule 1: Histone deacetylase 1

Chain I: 68% 8% 23%



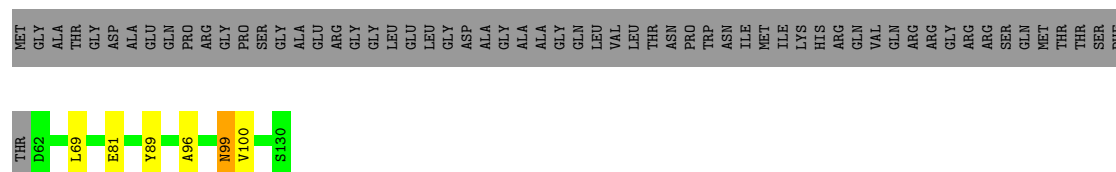
• Molecule 2: Deoxynucleotidyltransferase terminal-interacting protein 1

Chain A: 46% 7% 47%



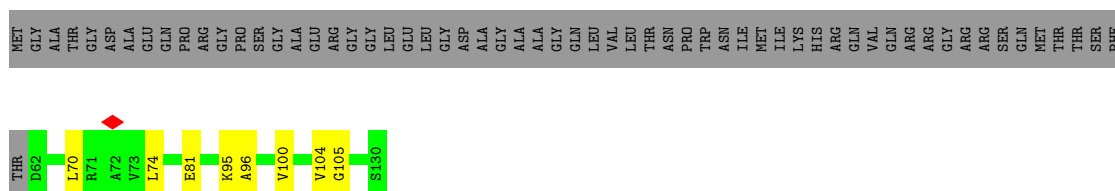
• Molecule 2: Deoxynucleotidyltransferase terminal-interacting protein 1

Chain B: 48% 0% 47%



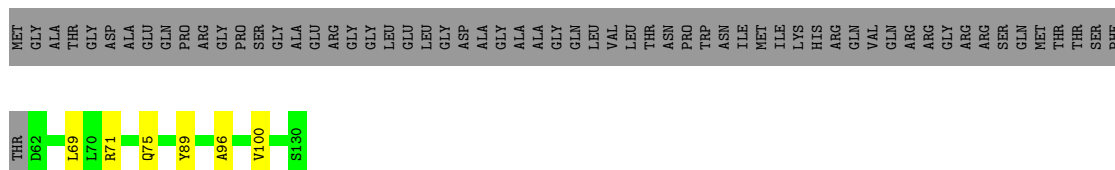
• Molecule 2: Deoxynucleotidyltransferase terminal-interacting protein 1

Chain G: 47% 6% 47%



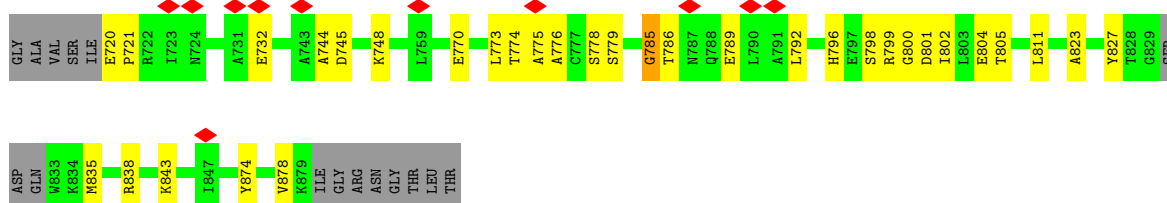
- Molecule 2: Deoxynucleotidyltransferase terminal-interacting protein 1

Chain H: 48% 5% 47%



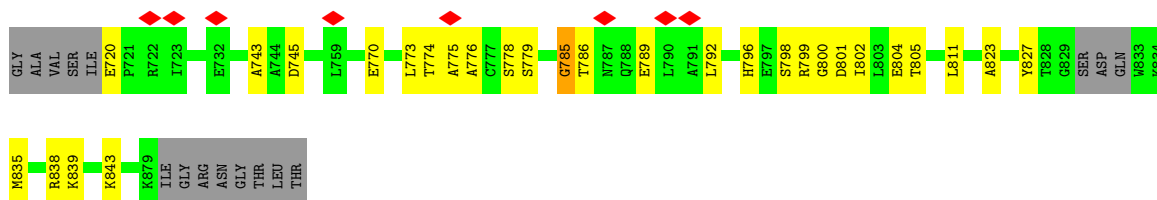
- Molecule 3: Mitotic deacetylase-associated SANT domain protein

Chain F: 6% 72% 18% 9%



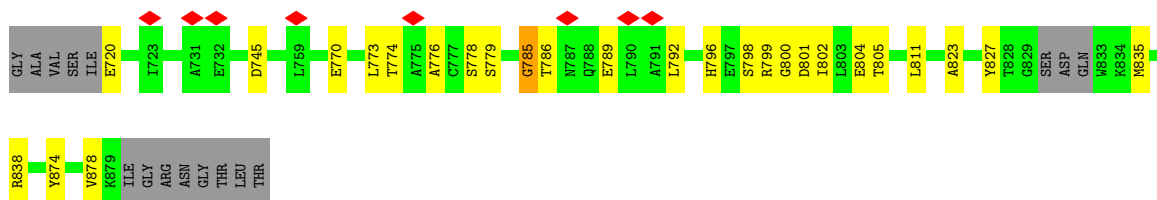
- Molecule 3: Mitotic deacetylase-associated SANT domain protein

Chain D: 5% 74% 16% 9%

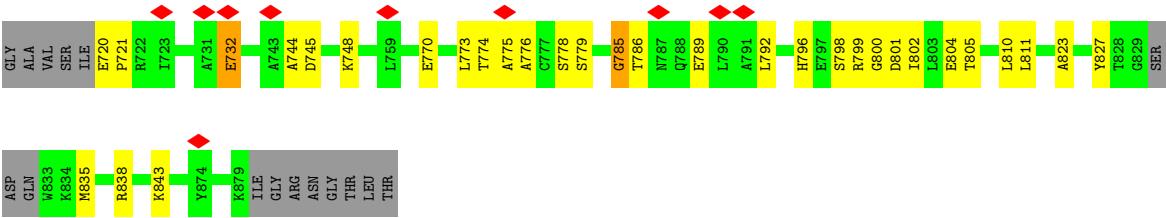


- Molecule 3: Mitotic deacetylase-associated SANT domain protein

Chain L: 5% 75% 15% 9%



- Molecule 3: Mitotic deacetylase-associated SANT domain protein



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D2	Depositor
Number of particles used	63222	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 ($e^-/\text{\AA}^2$)	36	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	500	Depositor
Magnification	129629	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.061	Depositor
Minimum map value	-0.037	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	432.00003, 432.00003, 432.00003	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	C	0.25	0/3042	0.49	0/4112
1	E	0.25	0/3033	0.49	0/4098
1	I	0.25	0/3042	0.49	0/4112
1	K	0.25	0/3042	0.49	0/4112
2	A	0.26	0/586	0.49	0/788
2	B	0.25	0/555	0.42	0/747
2	G	0.27	0/586	0.48	0/788
2	H	0.25	0/555	0.43	0/747
3	D	0.30	0/1288	0.74	2/1743 (0.1%)
3	F	0.30	0/1288	0.74	2/1743 (0.1%)
3	J	0.30	0/1288	0.74	2/1743 (0.1%)
3	L	0.30	0/1288	0.74	2/1743 (0.1%)
All	All	0.27	0/19593	0.56	8/26476 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	D	0	2
3	F	0	2
3	J	0	2
3	L	0	2
All	All	0	8

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	785	GLY	CA-C-N	5.86	129.53	121.33
3	D	785	GLY	C-N-CA	5.86	129.53	121.33
3	J	785	GLY	CA-C-N	5.83	129.50	121.33
3	J	785	GLY	C-N-CA	5.83	129.50	121.33
3	F	785	GLY	CA-C-N	5.83	129.49	121.33
3	F	785	GLY	C-N-CA	5.83	129.49	121.33
3	L	785	GLY	CA-C-N	5.82	129.47	121.33
3	L	785	GLY	C-N-CA	5.82	129.47	121.33

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	745	ASP	Peptide
3	D	811	LEU	Peptide
3	F	745	ASP	Peptide
3	F	811	LEU	Peptide
3	J	745	ASP	Peptide
3	J	811	LEU	Peptide
3	L	745	ASP	Peptide
3	L	811	LEU	Peptide

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	C	2964	2838	2836	23	0
1	E	2956	2831	2829	27	0
1	I	2964	2838	2836	30	0
1	K	2964	2838	2836	21	0
2	A	566	573	561	9	0
2	B	549	549	549	9	0
2	G	566	573	561	9	0
2	H	549	549	549	6	0
3	D	1258	1247	1247	26	0
3	F	1258	1247	1247	33	0
3	J	1258	1247	1247	34	0
3	L	1258	1247	1247	20	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	I	1	0	0	0	0
4	K	1	0	0	0	0
5	C	2	0	0	0	0
5	E	2	0	0	0	0
5	I	2	0	0	0	0
5	K	2	0	0	0	0
6	D	36	6	6	6	0
6	E	36	6	6	0	0
6	F	36	6	6	3	0
6	J	36	6	6	5	0
All	All	19266	18601	18569	199	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 5.

All (199) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:839:LYS:HE3	3:D:843:LYS:HE2	1.47	0.97
3:F:843:LYS:NZ	6:F:901:IHP:O32	1.99	0.94
3:J:843:LYS:NZ	6:J:901:IHP:O32	2.01	0.93
3:F:773:LEU:HD11	3:F:792:LEU:HD12	1.53	0.90
3:J:773:LEU:HD11	3:J:792:LEU:HD12	1.53	0.90
3:D:773:LEU:HD11	3:D:792:LEU:HD12	1.53	0.88
3:L:773:LEU:HD11	3:L:792:LEU:HD12	1.53	0.88
1:I:27:GLY:HA2	6:D:901:IHP:O41	1.81	0.79
3:D:796:HIS:O	3:D:799:ARG:NH2	2.19	0.76
3:L:796:HIS:O	3:L:799:ARG:NH2	2.19	0.76
3:D:839:LYS:NZ	6:D:901:IHP:O34	2.18	0.76
3:L:801:ASP:O	3:L:805:THR:OG1	2.04	0.76
3:F:796:HIS:O	3:F:799:ARG:NH2	2.19	0.76
3:J:796:HIS:O	3:J:799:ARG:NH2	2.19	0.76
1:K:34:ARG:NH1	1:K:302:GLY:O	2.20	0.75
3:D:801:ASP:O	3:D:805:THR:OG1	2.04	0.75
1:C:34:ARG:NH1	1:C:302:GLY:O	2.20	0.75
1:C:167:HIS:O	1:C:192:ARG:NH2	2.20	0.75
1:E:43:LEU:HD21	3:F:792:LEU:CD2	2.17	0.75
1:K:167:HIS:O	1:K:192:ARG:NH2	2.20	0.75
1:E:167:HIS:O	1:E:192:ARG:NH2	2.20	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:167:HIS:O	1:I:192:ARG:NH2	2.20	0.75
3:J:801:ASP:O	3:J:805:THR:OG1	2.04	0.74
3:F:801:ASP:O	3:F:805:THR:OG1	2.04	0.74
1:E:34:ARG:NH1	1:E:302:GLY:O	2.20	0.74
1:I:34:ARG:NH1	1:I:302:GLY:O	2.20	0.73
2:G:95:LYS:HG2	3:J:810:LEU:HD11	1.70	0.72
2:B:69:LEU:HD13	3:D:775:ALA:HA	1.71	0.71
2:G:81:GLU:OE2	2:H:89:TYR:OH	2.09	0.71
1:K:43:LEU:HD21	3:L:792:LEU:CD2	2.23	0.69
1:K:34:ARG:NH2	1:K:300:GLY:O	2.26	0.69
3:J:843:LYS:CE	6:J:901:IHP:O32	2.40	0.69
1:C:34:ARG:NH2	1:C:300:GLY:O	2.26	0.69
1:E:34:ARG:NH2	1:E:300:GLY:O	2.26	0.69
1:I:34:ARG:NH2	1:I:300:GLY:O	2.26	0.69
2:A:81:GLU:OE2	2:B:89:TYR:OH	2.11	0.69
1:C:8:ARG:NH2	1:C:321:ASP:OD1	2.26	0.69
1:K:8:ARG:NH2	1:K:321:ASP:OD1	2.26	0.69
1:I:8:ARG:NH2	1:I:321:ASP:OD1	2.26	0.68
1:E:8:ARG:NH2	1:E:321:ASP:OD1	2.26	0.68
1:I:55:ARG:HD3	3:J:823:ALA:HB2	1.76	0.68
3:F:843:LYS:HZ1	6:F:901:IHP:P2	2.20	0.65
2:H:71:ARG:NH1	2:H:75:GLN:OE1	2.31	0.63
1:E:55:ARG:HD3	3:F:823:ALA:HB2	1.81	0.62
3:L:798:SER:OG	3:L:804:GLU:OE1	2.16	0.62
1:I:287:PHE:O	1:I:290:SER:OG	2.14	0.62
1:E:287:PHE:O	1:E:290:SER:OG	2.14	0.62
1:I:43:LEU:HD21	3:J:792:LEU:CD2	2.30	0.61
3:F:773:LEU:HD11	3:F:792:LEU:CD1	2.28	0.60
1:K:287:PHE:O	1:K:290:SER:OG	2.14	0.60
3:D:773:LEU:HD11	3:D:792:LEU:CD1	2.28	0.60
3:J:773:LEU:HD11	3:J:792:LEU:CD1	2.28	0.60
1:C:287:PHE:O	1:C:290:SER:OG	2.14	0.60
3:L:773:LEU:HD11	3:L:792:LEU:CD1	2.28	0.60
1:C:55:ARG:HD3	3:D:823:ALA:HB2	1.84	0.60
3:F:843:LYS:CE	6:F:901:IHP:O32	2.50	0.59
1:E:122:VAL:HG11	3:F:744:ALA:O	2.03	0.58
1:I:122:VAL:HG11	3:J:744:ALA:O	2.02	0.58
2:G:100:VAL:HG22	3:J:779:SER:CB	2.33	0.58
1:C:43:LEU:HD21	3:D:792:LEU:CD2	2.33	0.58
3:F:874:TYR:OH	3:L:878:VAL:HG22	2.05	0.57
2:B:69:LEU:HB3	3:D:775:ALA:HA	1.88	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:843:LYS:HZ1	6:J:901:IHP:P2	2.28	0.56
1:I:164:LEU:HD13	3:J:732:GLU:CB	2.36	0.56
2:A:96:ALA:O	2:A:100:VAL:HG23	2.06	0.56
2:H:69:LEU:HD13	3:J:775:ALA:HA	1.87	0.55
1:K:231:GLY:N	1:K:350:MET:HE2	2.22	0.55
1:C:231:GLY:N	1:C:350:MET:HE2	2.22	0.55
1:E:119:ALA:HB2	3:F:748:LYS:HB3	1.88	0.55
1:E:231:GLY:N	1:E:350:MET:HE2	2.22	0.55
1:I:231:GLY:N	1:I:350:MET:HE2	2.22	0.55
1:E:335:GLU:N	1:E:335:GLU:OE1	2.40	0.54
1:I:335:GLU:N	1:I:335:GLU:OE1	2.40	0.54
2:G:100:VAL:HG22	3:J:779:SER:HB2	1.90	0.54
3:D:839:LYS:HZ3	6:D:901:IHP:P4	2.31	0.54
1:C:335:GLU:OE1	1:C:335:GLU:N	2.40	0.53
1:K:55:ARG:HD3	3:L:823:ALA:HB2	1.89	0.53
2:H:100:VAL:HG22	3:L:779:SER:CB	2.38	0.53
2:B:100:VAL:HG22	3:F:779:SER:CB	2.38	0.53
1:K:335:GLU:OE1	1:K:335:GLU:N	2.40	0.53
3:F:798:SER:OG	3:F:804:GLU:OE1	2.16	0.53
3:J:835:MET:SD	3:J:838:ARG:NH2	2.82	0.53
3:F:835:MET:SD	3:F:838:ARG:NH2	2.82	0.53
3:D:801:ASP:OD1	3:D:802:ILE:N	2.41	0.53
3:D:835:MET:SD	3:D:838:ARG:NH2	2.82	0.53
3:L:801:ASP:OD1	3:L:802:ILE:N	2.41	0.53
3:L:835:MET:SD	3:L:838:ARG:NH2	2.82	0.52
3:J:798:SER:OG	3:J:804:GLU:OE1	2.16	0.52
3:J:801:ASP:OD1	3:J:802:ILE:N	2.41	0.52
3:F:785:GLY:O	3:F:786:THR:HG23	2.09	0.52
3:D:785:GLY:O	3:D:786:THR:HG23	2.09	0.52
3:J:785:GLY:O	3:J:786:THR:HG23	2.09	0.52
3:L:785:GLY:O	3:L:786:THR:HG23	2.09	0.52
3:F:801:ASP:OD1	3:F:802:ILE:N	2.41	0.52
1:I:200:LYS:NZ	1:I:230:ASP:OD1	2.43	0.52
3:J:843:LYS:HE3	6:J:901:IHP:O32	2.09	0.51
1:E:34:ARG:HD2	1:E:308:VAL:HG11	1.93	0.51
1:I:34:ARG:HD2	1:I:308:VAL:HG11	1.93	0.51
3:D:839:LYS:NZ	6:D:901:IHP:P4	2.84	0.51
1:I:231:GLY:CA	1:I:350:MET:HE2	2.40	0.51
1:C:34:ARG:HD2	1:C:308:VAL:HG11	1.93	0.50
1:C:177:ILE:HG22	1:C:264:ASP:CG	2.36	0.50
1:E:231:GLY:CA	1:E:350:MET:HE2	2.41	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:34:ARG:HD2	1:K:308:VAL:HG11	1.93	0.50
1:K:177:ILE:HG22	1:K:264:ASP:CG	2.36	0.50
1:I:310:ARG:NE	1:I:342:LYS:O	2.39	0.50
1:K:231:GLY:CA	1:K:350:MET:HE2	2.41	0.50
1:C:231:GLY:CA	1:C:350:MET:HE2	2.41	0.50
1:E:310:ARG:NE	1:E:342:LYS:O	2.39	0.50
1:I:177:ILE:HG22	1:I:264:ASP:CG	2.36	0.50
2:A:100:VAL:HG22	3:D:779:SER:CB	2.42	0.50
1:E:177:ILE:HG22	1:E:264:ASP:CG	2.36	0.49
1:C:117:SER:O	1:C:120:SER:OG	2.23	0.49
1:K:117:SER:O	1:K:120:SER:OG	2.23	0.49
3:D:798:SER:OG	3:D:804:GLU:OE1	2.15	0.49
3:F:789:GLU:HA	3:F:792:LEU:HD13	1.95	0.48
3:F:878:VAL:HG22	3:L:874:TYR:OH	2.13	0.48
3:D:789:GLU:HA	3:D:792:LEU:HD13	1.95	0.48
3:L:789:GLU:HA	3:L:792:LEU:HD13	1.95	0.48
3:J:789:GLU:HA	3:J:792:LEU:HD13	1.96	0.48
1:K:43:LEU:HD21	3:L:792:LEU:HD21	1.96	0.48
1:E:164:LEU:HD13	3:F:732:GLU:CB	2.44	0.48
2:A:89:TYR:OH	2:B:81:GLU:OE2	2.31	0.47
2:B:100:VAL:HG22	3:F:779:SER:HB3	1.95	0.47
1:E:262:GLY:O	1:E:281:HIS:NE2	2.47	0.47
1:K:200:LYS:NZ	1:K:230:ASP:OD1	2.43	0.47
1:C:200:LYS:NZ	1:C:230:ASP:OD1	2.43	0.47
1:I:262:GLY:O	1:I:281:HIS:NE2	2.47	0.47
2:H:100:VAL:HG22	3:L:779:SER:HB2	1.95	0.47
1:E:43:LEU:HD21	3:F:792:LEU:HD21	1.96	0.47
1:K:262:GLY:O	1:K:281:HIS:NE2	2.47	0.47
1:C:262:GLY:O	1:C:281:HIS:NE2	2.47	0.47
1:E:197:SER:OG	1:E:199:HIS:ND1	2.48	0.47
1:I:197:SER:OG	1:I:199:HIS:ND1	2.48	0.47
1:I:31:LYS:HE3	6:D:901:IHP:O26	2.15	0.46
1:I:164:LEU:HD13	3:J:732:GLU:HB2	1.97	0.46
1:E:200:LYS:NZ	1:E:230:ASP:OD1	2.43	0.46
2:G:100:VAL:HG22	3:J:779:SER:HB3	1.97	0.46
1:E:314:TYR:CD1	1:E:343:LEU:HD21	2.51	0.46
3:D:720:GLU:N	3:D:720:GLU:OE1	2.49	0.46
3:L:720:GLU:N	3:L:720:GLU:OE1	2.49	0.46
1:I:314:TYR:CD1	1:I:343:LEU:HD21	2.51	0.45
3:J:720:GLU:N	3:J:720:GLU:OE1	2.49	0.45
2:A:100:VAL:HG22	3:D:779:SER:HB2	1.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:720:GLU:N	3:F:720:GLU:OE1	2.49	0.45
3:D:770:GLU:O	3:D:774:THR:HG22	2.16	0.45
3:J:770:GLU:O	3:J:774:THR:HG22	2.16	0.45
1:K:197:SER:OG	1:K:199:HIS:ND1	2.48	0.45
3:F:770:GLU:O	3:F:774:THR:HG22	2.16	0.45
3:J:843:LYS:NZ	6:J:901:IHP:P2	2.87	0.45
3:L:770:GLU:O	3:L:774:THR:HG22	2.16	0.45
1:C:314:TYR:CD1	1:C:343:LEU:HD21	2.51	0.45
1:C:197:SER:OG	1:C:199:HIS:ND1	2.48	0.45
1:C:43:LEU:HD21	3:D:792:LEU:HD21	1.97	0.45
1:K:314:TYR:CD1	1:K:343:LEU:HD21	2.51	0.45
1:I:119:ALA:HB2	3:J:748:LYS:HB3	1.97	0.44
1:I:43:LEU:HD21	3:J:792:LEU:HD21	1.99	0.44
2:H:96:ALA:O	2:H:100:VAL:HG23	2.17	0.44
3:D:776:ALA:O	3:D:778:SER:N	2.50	0.44
3:L:776:ALA:O	3:L:778:SER:N	2.50	0.44
1:C:310:ARG:NE	1:C:342:LYS:O	2.39	0.44
1:K:310:ARG:NE	1:K:342:LYS:O	2.39	0.44
1:C:126:LYS:HE2	3:D:743:ALA:HB3	1.99	0.44
1:I:27:GLY:CA	6:D:901:IHP:O41	2.58	0.44
1:E:161:LEU:HD22	3:F:732:GLU:HA	1.99	0.43
2:B:99:ASN:C	2:B:99:ASN:HD22	2.27	0.43
2:B:100:VAL:HG22	3:F:779:SER:HB2	2.01	0.42
2:A:69:LEU:CB	3:F:775:ALA:HA	2.50	0.42
1:E:57:HIS:N	3:F:748:LYS:O	2.48	0.42
2:B:96:ALA:O	2:B:100:VAL:HG23	2.18	0.42
1:E:177:ILE:HG23	1:E:204:TYR:OH	2.20	0.42
1:E:208:THR:HG22	3:F:721:PRO:HD3	2.01	0.42
1:I:177:ILE:HG23	1:I:204:TYR:OH	2.20	0.42
2:A:104:VAL:HG12	2:A:105:GLY:H	1.84	0.42
2:G:95:LYS:CG	3:J:810:LEU:HD11	2.45	0.41
1:C:248:VAL:O	1:C:252:PHE:N	2.50	0.41
2:A:70:LEU:HD11	2:A:74:LEU:HD11	2.03	0.41
1:C:177:ILE:HG23	1:C:204:TYR:OH	2.20	0.41
1:K:248:VAL:O	1:K:252:PHE:N	2.50	0.41
2:G:70:LEU:HD11	2:G:74:LEU:HD11	2.02	0.41
1:K:177:ILE:HG23	1:K:204:TYR:OH	2.20	0.41
1:I:117:SER:O	1:I:120:SER:OG	2.23	0.41
3:D:823:ALA:HB1	3:D:827:TYR:CZ	2.56	0.41
3:L:823:ALA:HB1	3:L:827:TYR:CZ	2.56	0.41
1:E:117:SER:O	1:E:120:SER:OG	2.23	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:96:ALA:O	2:G:100:VAL:HG23	2.20	0.41
3:F:776:ALA:O	3:F:778:SER:N	2.50	0.41
2:G:104:VAL:HG12	2:G:105:GLY:H	1.85	0.41
3:J:776:ALA:O	3:J:778:SER:N	2.50	0.41
3:J:823:ALA:HB1	3:J:827:TYR:CZ	2.56	0.40
1:C:43:LEU:HD11	3:D:792:LEU:HD23	2.03	0.40
1:I:161:LEU:HD22	3:J:732:GLU:HA	2.03	0.40
1:I:208:THR:HG22	3:J:721:PRO:HD3	2.02	0.40
3:F:823:ALA:HB1	3:F:827:TYR:CZ	2.56	0.40
3:F:744:ALA:HB3	3:F:748:LYS:NZ	2.37	0.40
3:J:744:ALA:HB3	3:J:748:LYS:NZ	2.37	0.40
1:E:241:PHE:CE1	1:E:245:MET:HE3	2.57	0.40
1:I:241:PHE:CE1	1:I:245:MET:HE3	2.57	0.40
2:A:69:LEU:HD12	3:F:775:ALA:HA	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	
1	C	367/482 (76%)	351 (96%)	16 (4%)	0	100	100
1	E	364/482 (76%)	349 (96%)	15 (4%)	0	100	100
1	I	367/482 (76%)	351 (96%)	16 (4%)	0	100	100
1	K	367/482 (76%)	351 (96%)	16 (4%)	0	100	100
2	A	71/130 (55%)	71 (100%)	0	0	100	100
2	B	67/130 (52%)	67 (100%)	0	0	100	100
2	G	71/130 (55%)	71 (100%)	0	0	100	100
2	H	67/130 (52%)	67 (100%)	0	0	100	100
3	D	153/173 (88%)	115 (75%)	37 (24%)	1 (1%)	18	55

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	F	153/173 (88%)	115 (75%)	37 (24%)	1 (1%)	18	55
3	J	153/173 (88%)	115 (75%)	36 (24%)	2 (1%)	9	41
3	L	153/173 (88%)	115 (75%)	37 (24%)	1 (1%)	18	55
All	All	2353/3140 (75%)	2138 (91%)	210 (9%)	5 (0%)	44	77

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	800	GLY
3	D	800	GLY
3	L	800	GLY
3	J	732	GLU
3	J	800	GLY

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	C	315/419 (75%)	315 (100%)	0	100	100
1	E	314/419 (75%)	314 (100%)	0	100	100
1	I	315/419 (75%)	315 (100%)	0	100	100
1	K	315/419 (75%)	315 (100%)	0	100	100
2	A	65/106 (61%)	65 (100%)	0	100	100
2	B	61/106 (58%)	60 (98%)	1 (2%)	55	69
2	G	65/106 (61%)	65 (100%)	0	100	100
2	H	61/106 (58%)	61 (100%)	0	100	100
3	D	131/149 (88%)	131 (100%)	0	100	100
3	F	131/149 (88%)	131 (100%)	0	100	100
3	J	131/149 (88%)	131 (100%)	0	100	100
3	L	131/149 (88%)	131 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	2035/2696 (76%)	2034 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	99	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	90	GLN
1	C	111	GLN
1	C	140	HIS
1	C	292	ASN
1	E	90	GLN
1	E	111	GLN
1	E	292	ASN
1	K	90	GLN
1	K	111	GLN
1	K	292	ASN
1	I	90	GLN
1	I	111	GLN
1	I	292	ASN
2	B	115	GLN
2	H	99	ASN
2	H	115	GLN
3	F	766	GLN
3	D	766	GLN
3	L	766	GLN
3	J	766	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	IHP	D	901	-	36,36,36	0.84	1 (2%)	60,60,60	1.11	4 (6%)
6	IHP	E	501	-	36,36,36	0.84	1 (2%)	60,60,60	1.11	5 (8%)
6	IHP	F	901	-	36,36,36	0.83	1 (2%)	60,60,60	1.11	5 (8%)
6	IHP	J	901	-	36,36,36	0.83	1 (2%)	60,60,60	1.11	4 (6%)

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
6	IHP	D	901	-	-	1/30/54/54	0/1/1/1
6	IHP	E	501	-	-	1/30/54/54	0/1/1/1
6	IHP	F	901	-	-	1/30/54/54	0/1/1/1
6	IHP	J	901	-	-	1/30/54/54	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	901	IHP	P3-O13	2.31	1.63	1.59
6	E	501	IHP	P3-O13	2.31	1.63	1.59
6	J	901	IHP	P3-O13	2.27	1.63	1.59
6	F	901	IHP	P3-O13	2.18	1.63	1.59

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	901	IHP	O11-C1-C6	2.26	113.56	108.76
6	E	501	IHP	O11-C1-C6	2.25	113.55	108.76
6	J	901	IHP	O11-C1-C6	2.25	113.54	108.76
6	D	901	IHP	O11-C1-C6	2.24	113.52	108.76
6	E	501	IHP	O43-P3-O33	2.22	116.12	107.80
6	D	901	IHP	O43-P3-O33	2.21	116.10	107.80
6	J	901	IHP	O43-P3-O33	2.21	116.09	107.80
6	F	901	IHP	O43-P3-O33	2.21	116.08	107.80
6	D	901	IHP	O41-P1-O31	2.20	116.06	107.80
6	E	501	IHP	O41-P1-O31	2.20	116.06	107.80
6	J	901	IHP	O41-P1-O31	2.20	116.05	107.80
6	F	901	IHP	O41-P1-O31	2.19	116.02	107.80
6	F	901	IHP	O13-P3-O23	-2.03	102.10	109.33
6	E	501	IHP	O13-P3-O23	-2.02	102.13	109.33
6	D	901	IHP	O13-P3-O23	-2.02	102.14	109.33
6	J	901	IHP	O13-P3-O23	-2.01	102.17	109.33
6	E	501	IHP	O14-P4-O24	-2.00	102.19	109.33
6	F	901	IHP	O14-P4-O24	-2.00	102.20	109.33

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	E	501	IHP	C2-O12-P2-O42
6	F	901	IHP	C2-O12-P2-O42
6	D	901	IHP	C2-O12-P2-O42
6	J	901	IHP	C2-O12-P2-O42

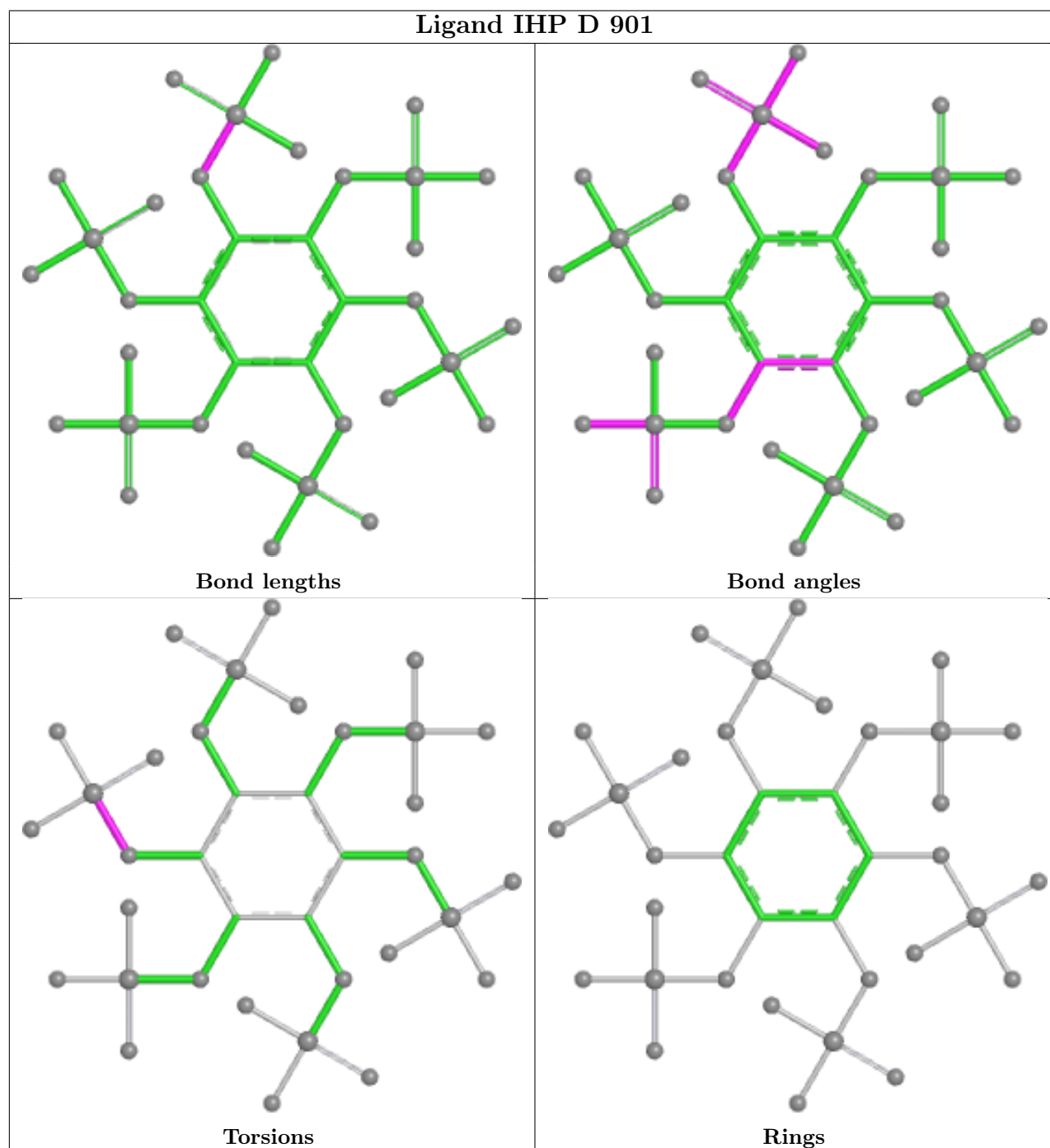
There are no ring outliers.

3 monomers are involved in 14 short contacts:

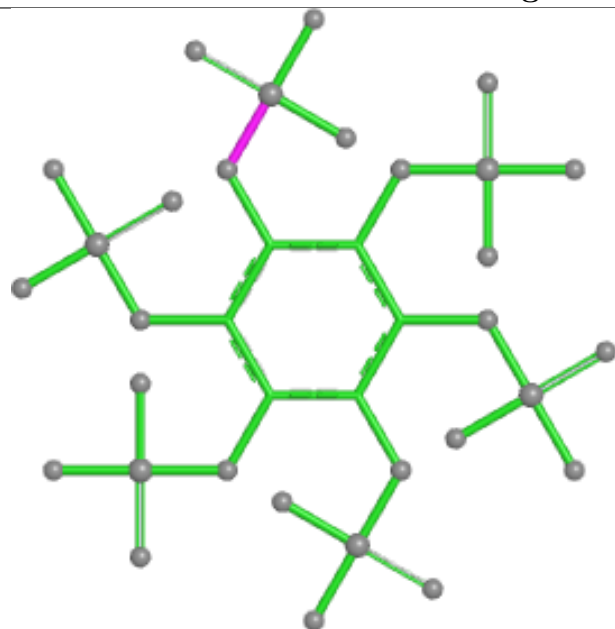
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	901	IHP	6	0
6	F	901	IHP	3	0
6	J	901	IHP	5	0

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

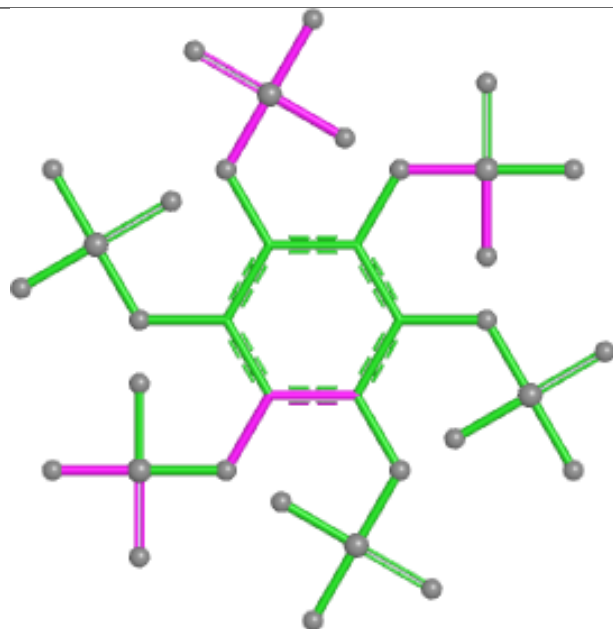
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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.



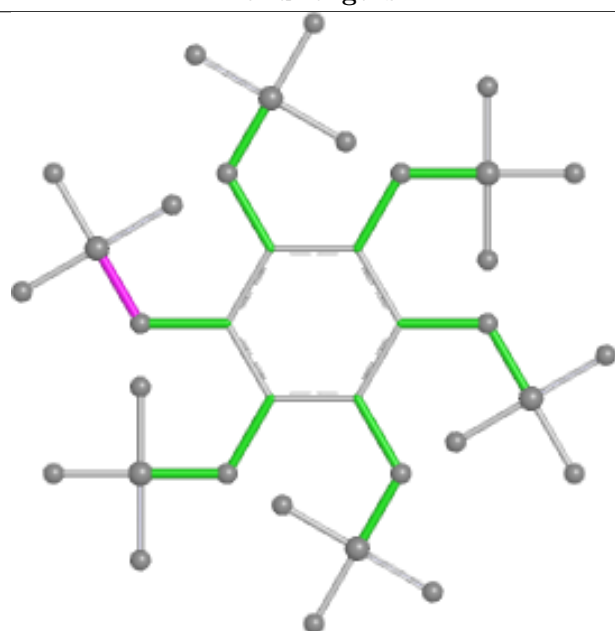
Ligand IHP E 501



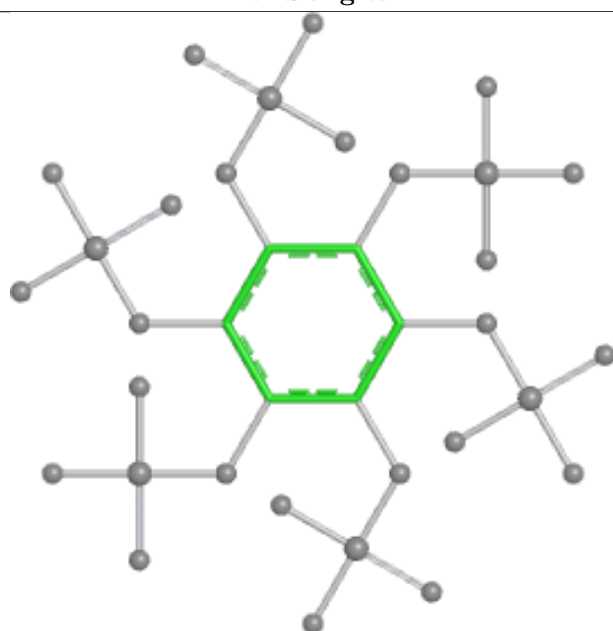
Bond lengths



Bond angles

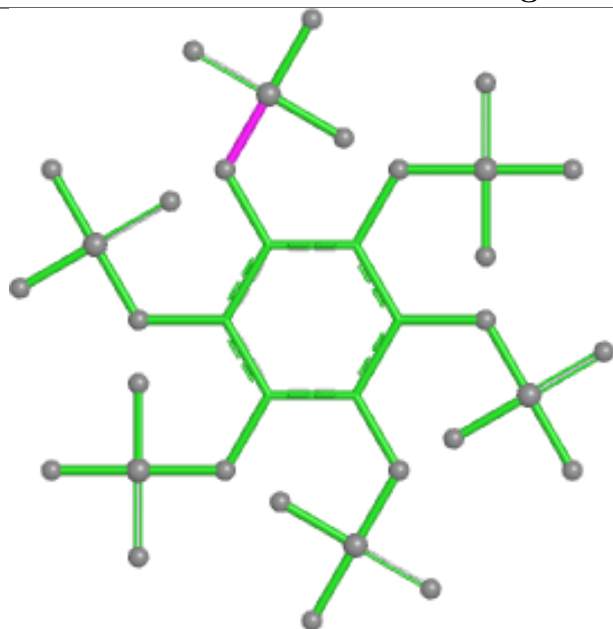


Torsions

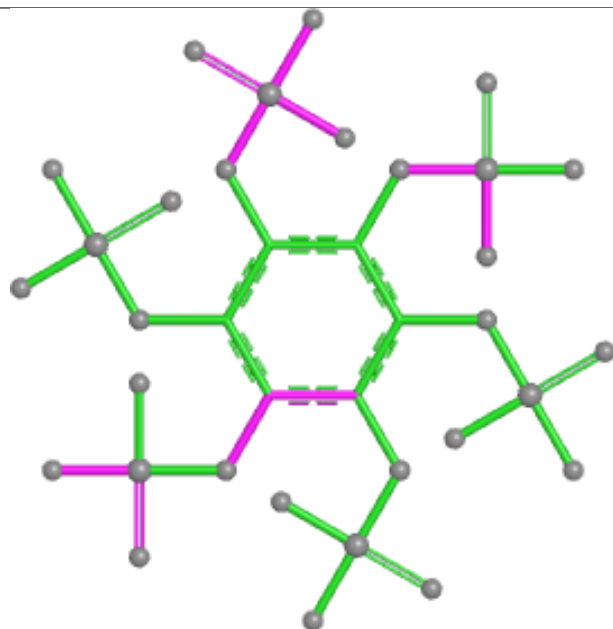


Rings

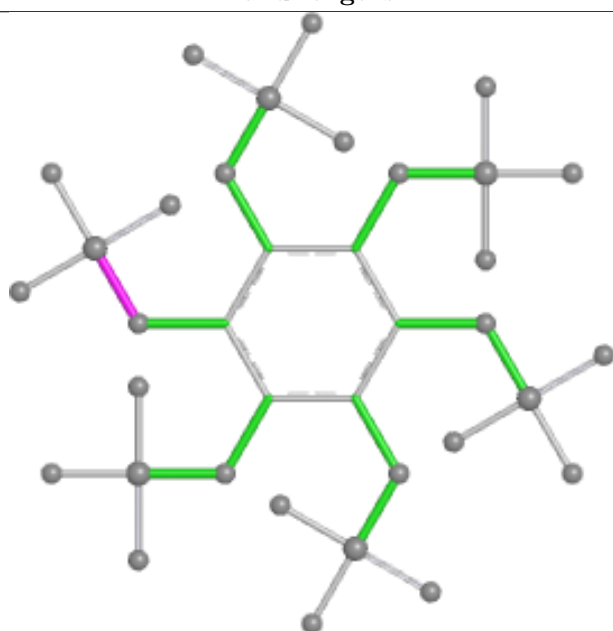
Ligand IHP F 901



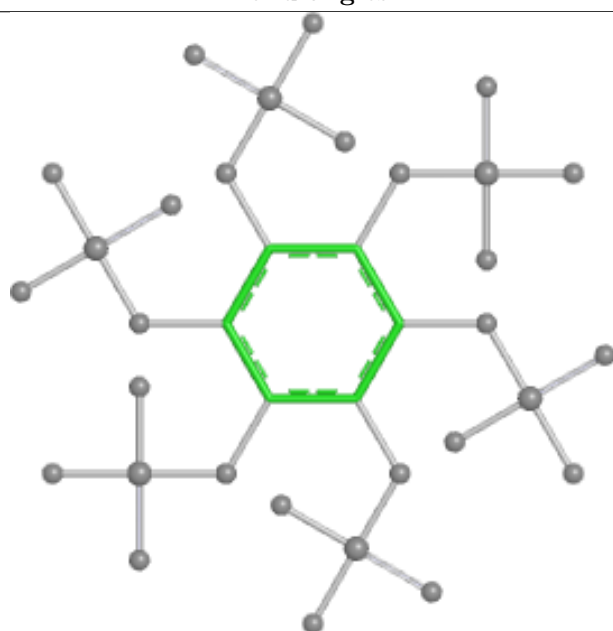
Bond lengths



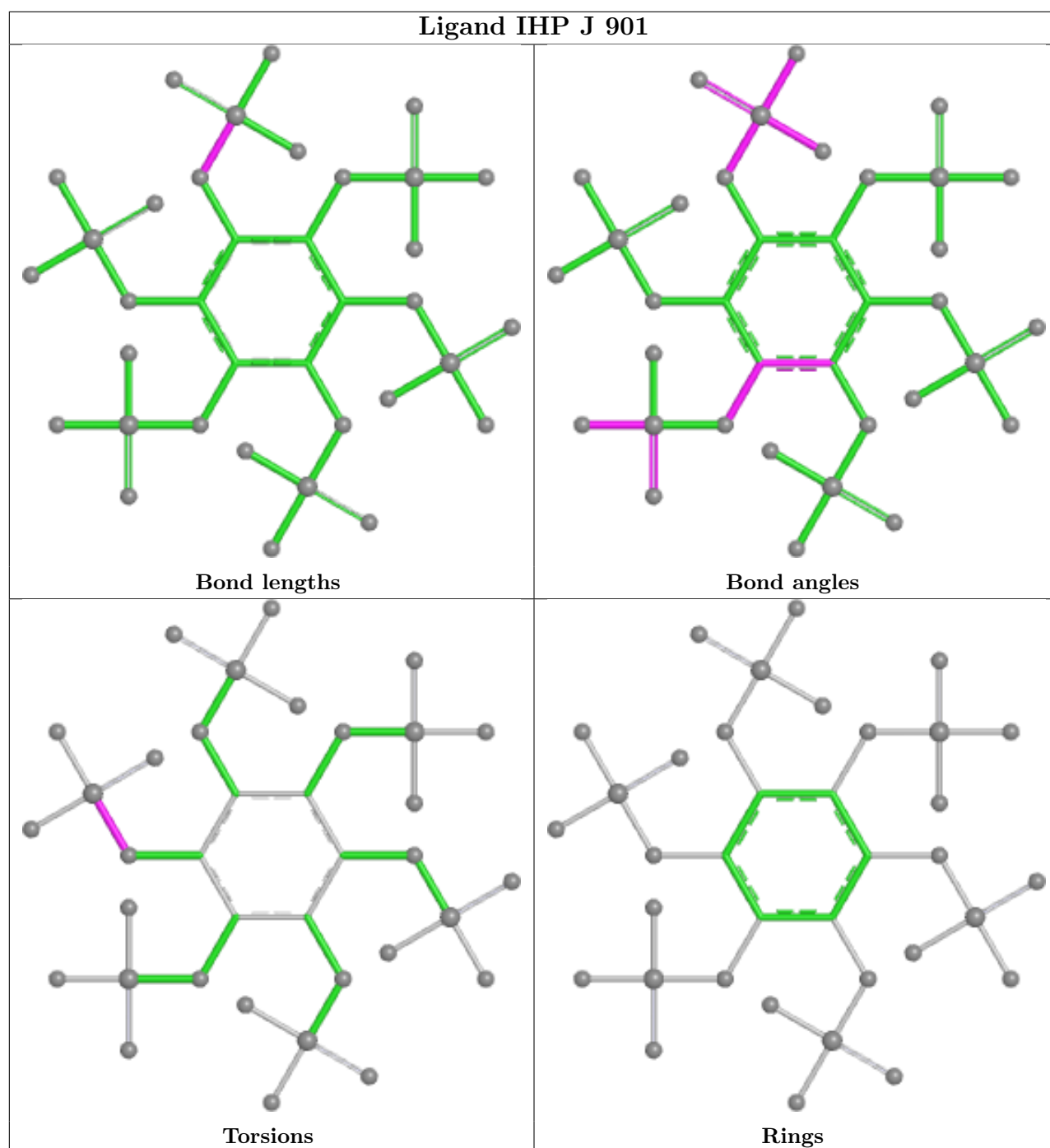
Bond angles



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.

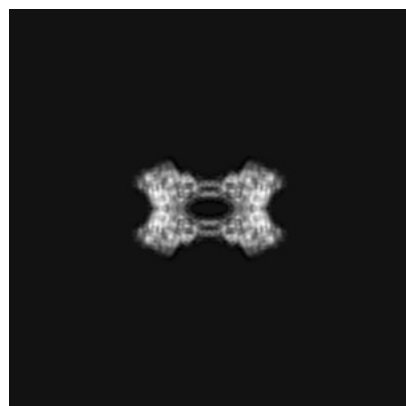
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11042. These allow visual inspection of the internal detail of the map and identification of artifacts.

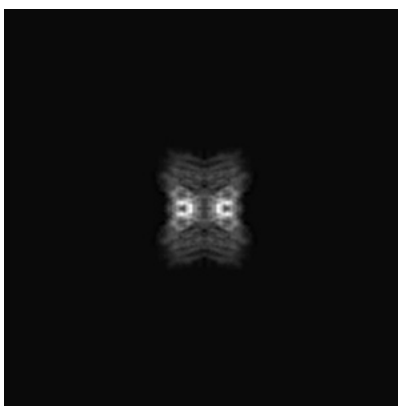
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



X

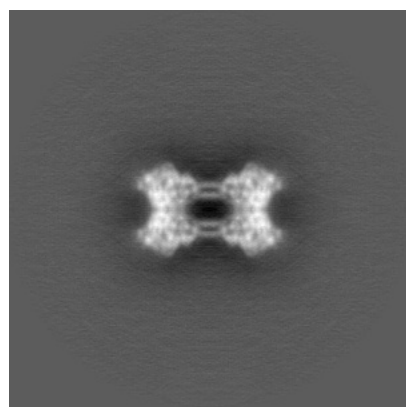


Y

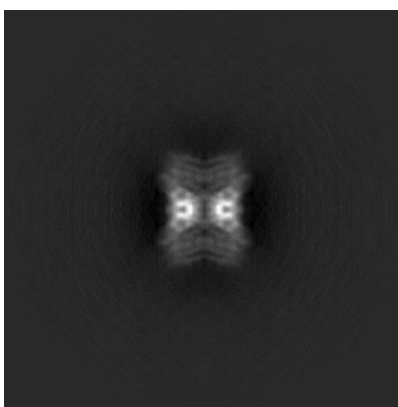


Z

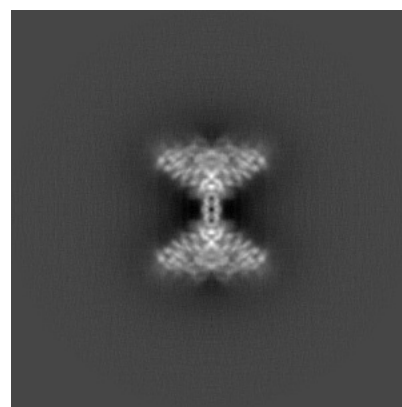
6.1.2 Raw map



X



Y



Z

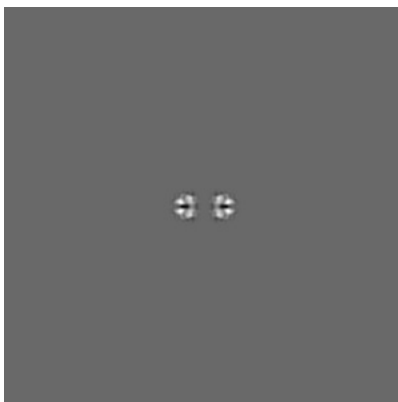
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 200

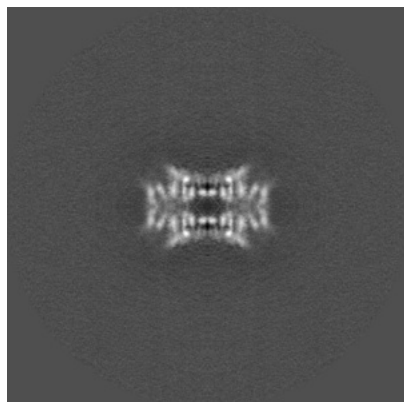


Y Index: 200

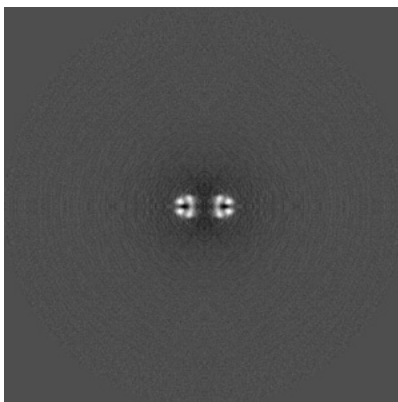


Z Index: 200

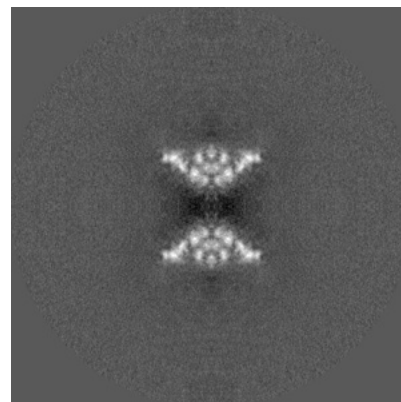
6.2.2 Raw map



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: 205

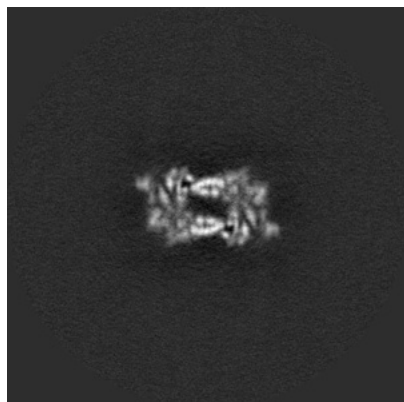


Y Index: 150

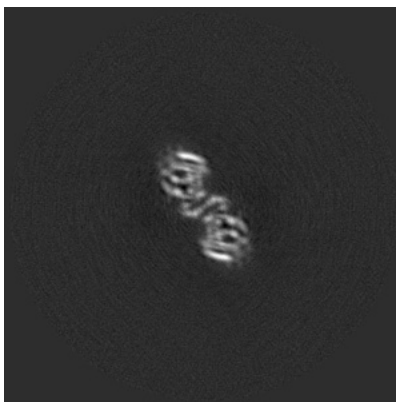


Z Index: 217

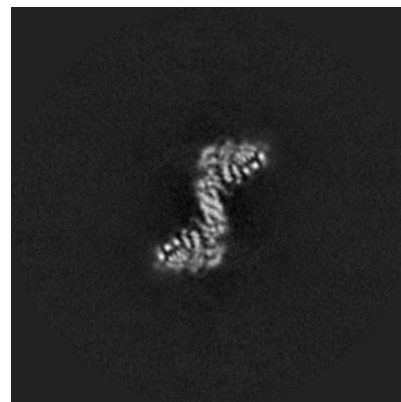
6.3.2 Raw map



X Index: 195



Y Index: 150

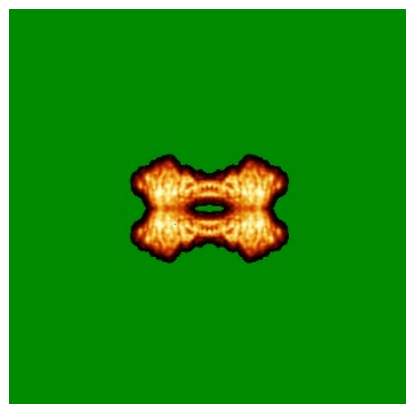


Z Index: 215

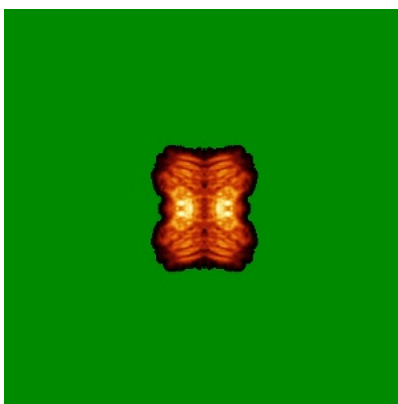
The images above show the largest variance slices of the map in three orthogonal directions.

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

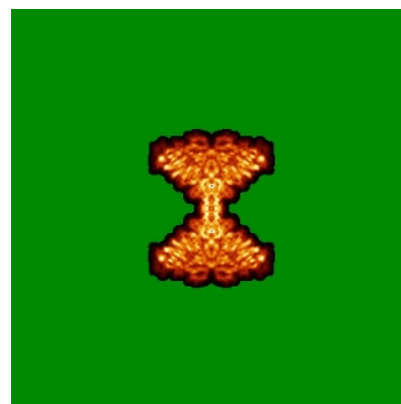
6.4.1 Primary map



X

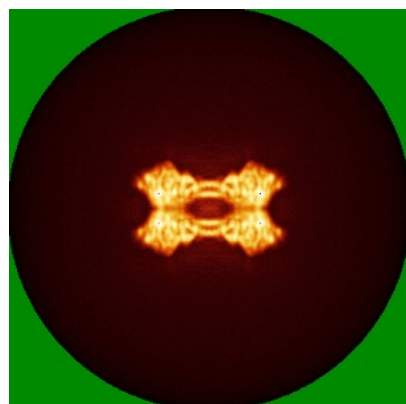


Y

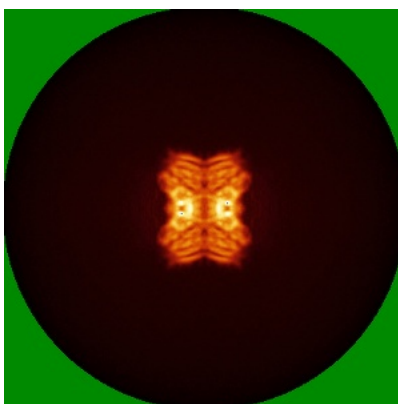


Z

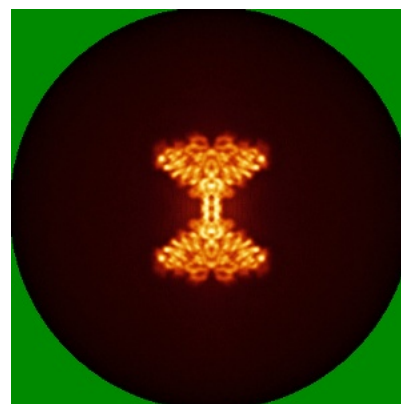
6.4.2 Raw map



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 0.01. 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.

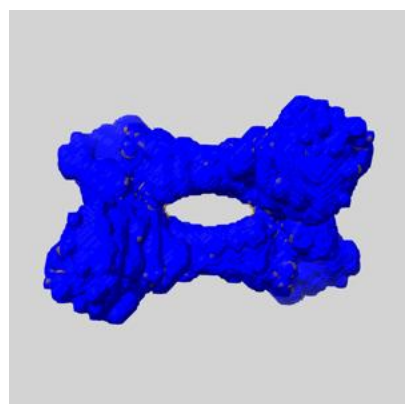
6.6 Mask visualisation [i](#)

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

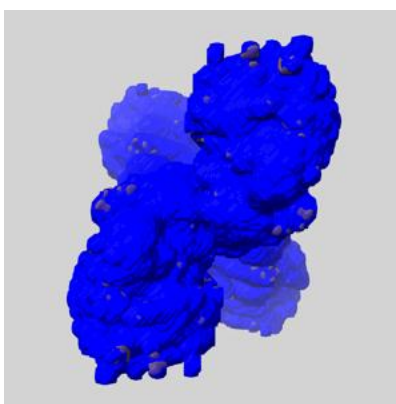
A mask typically either:

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

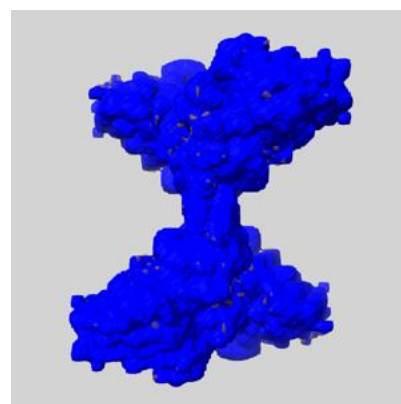
6.6.1 emd_11042_msk_1.map [i](#)



X



Y

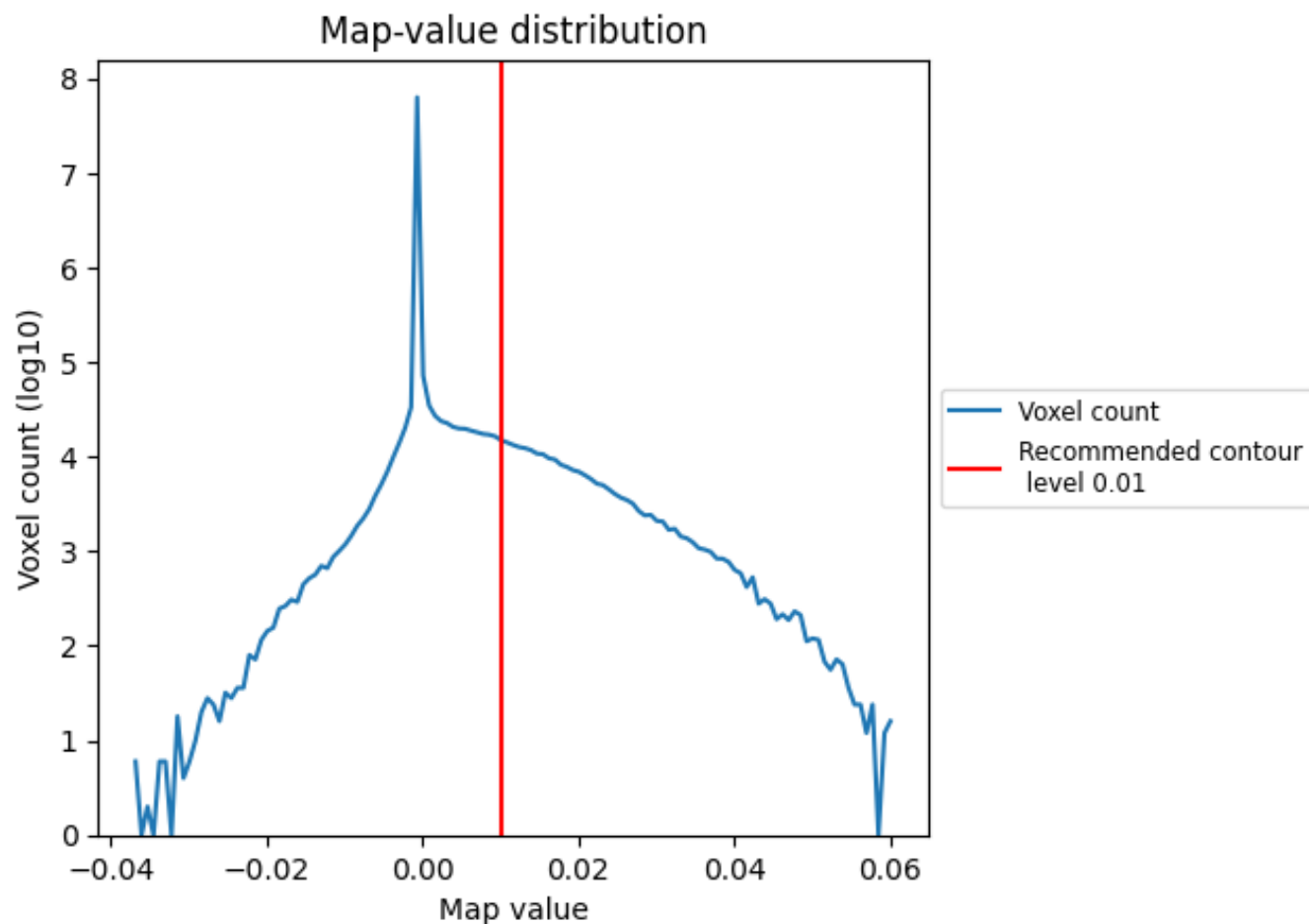


Z

7 Map analysis [i](#)

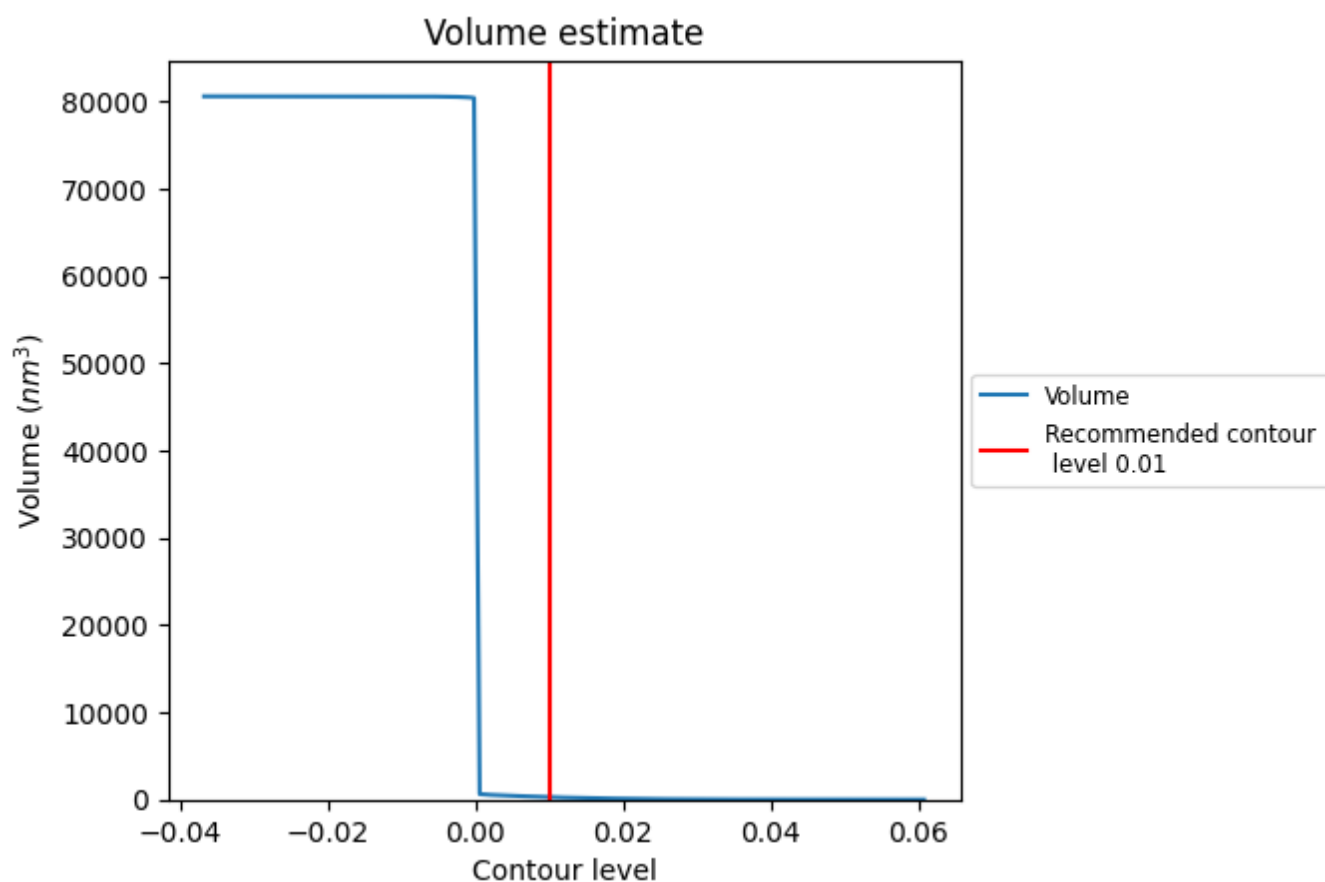
This section contains the results of statistical analysis of the map.

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.

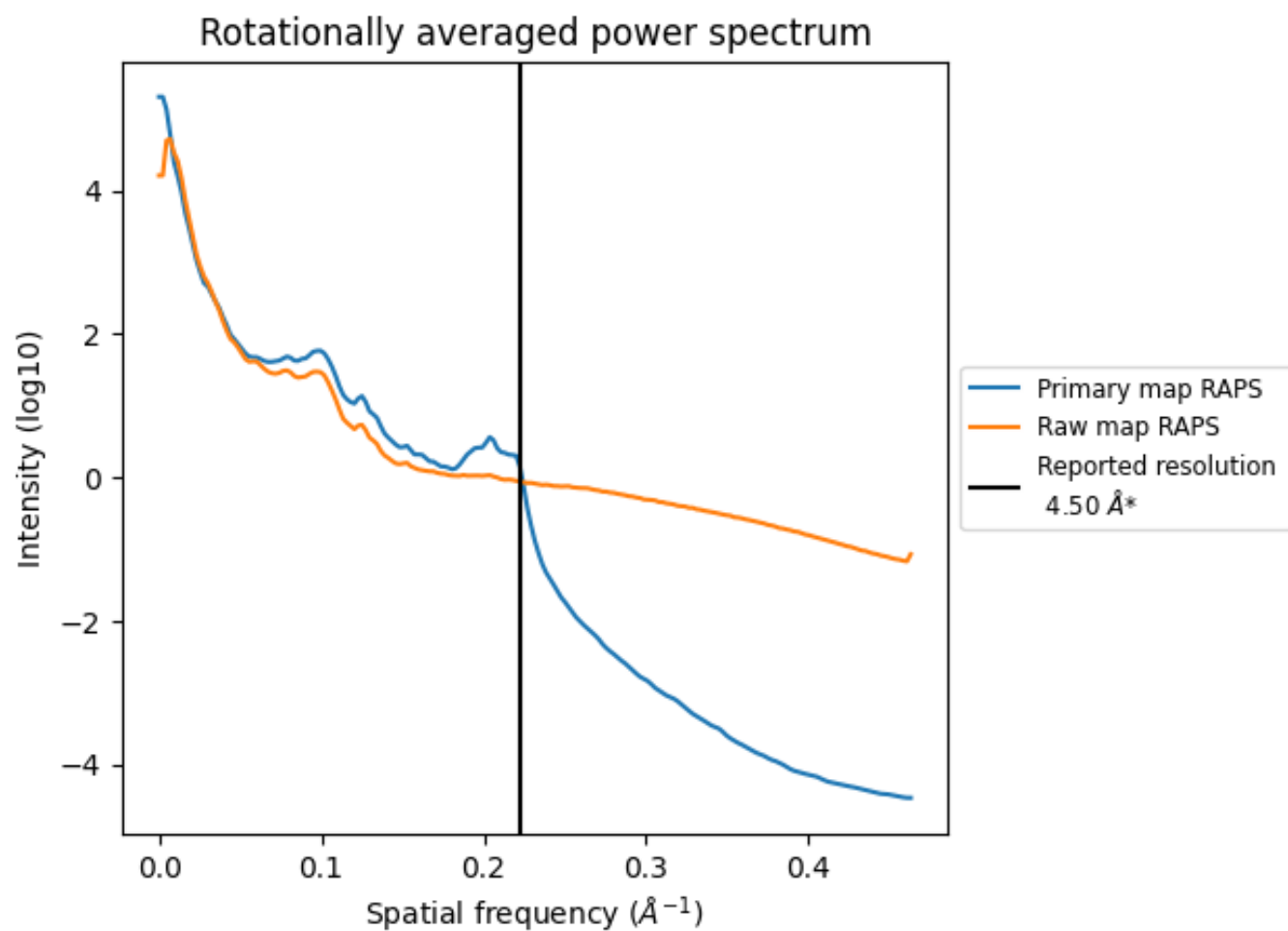
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 280 nm³; this corresponds to an approximate mass of 253 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 ⓘ

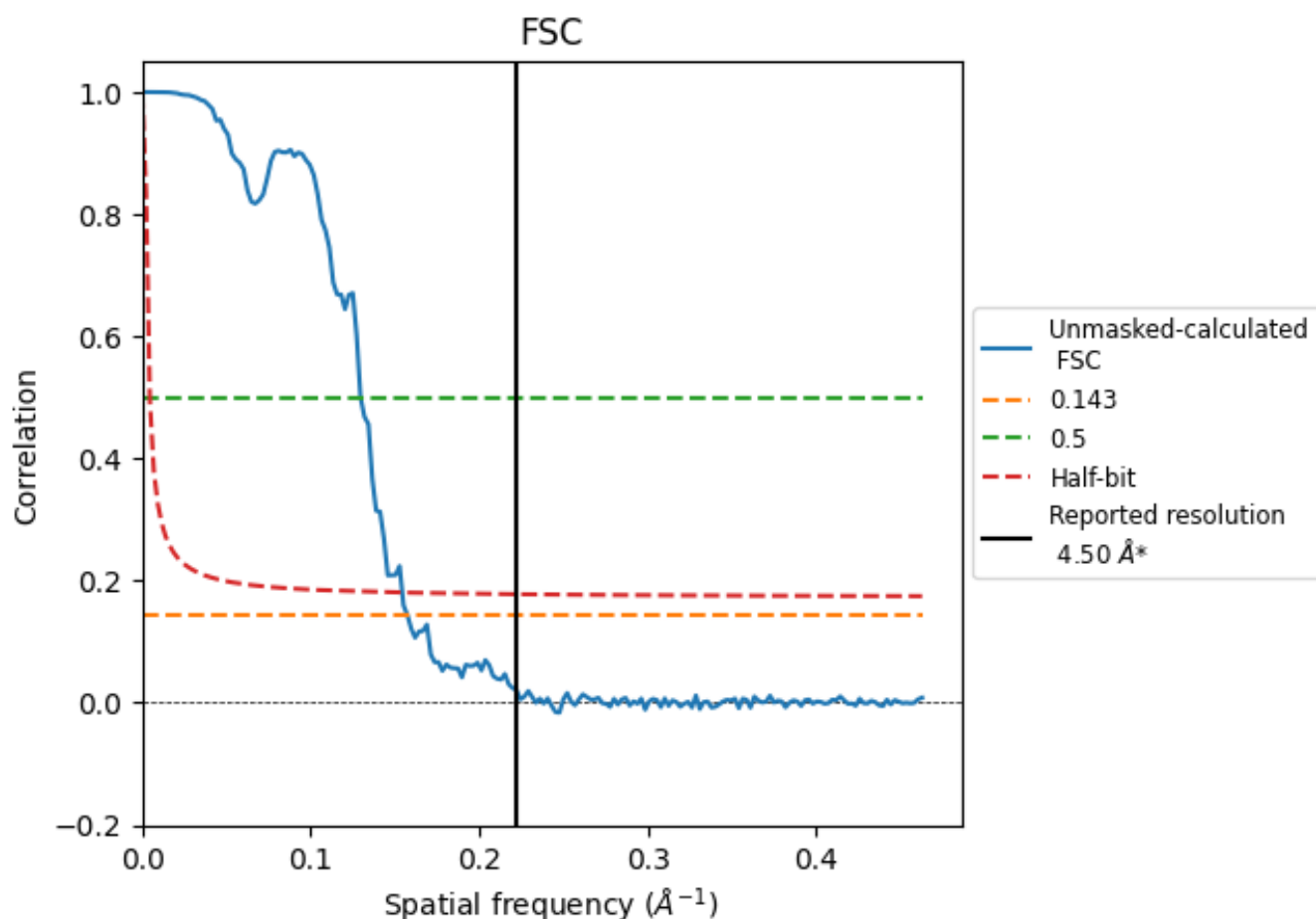


*Reported resolution corresponds to spatial frequency of 0.222 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.222 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

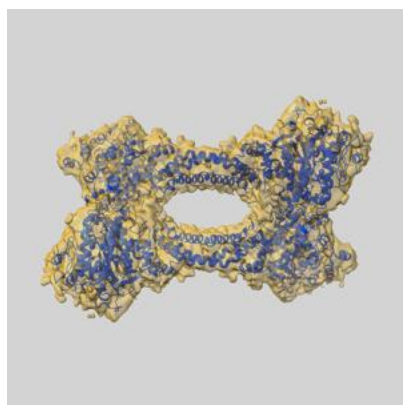
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	6.35	7.72	6.48

*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 6.35 differs from the reported value 4.5 by more than 10 %

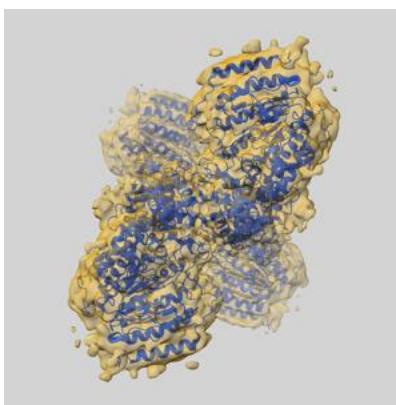
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-11042 and PDB model 6Z2K. Per-residue inclusion information can be found in section [3](#) on page [7](#).

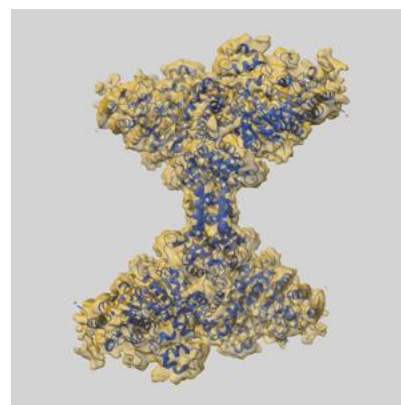
9.1 Map-model overlay [i](#)



X



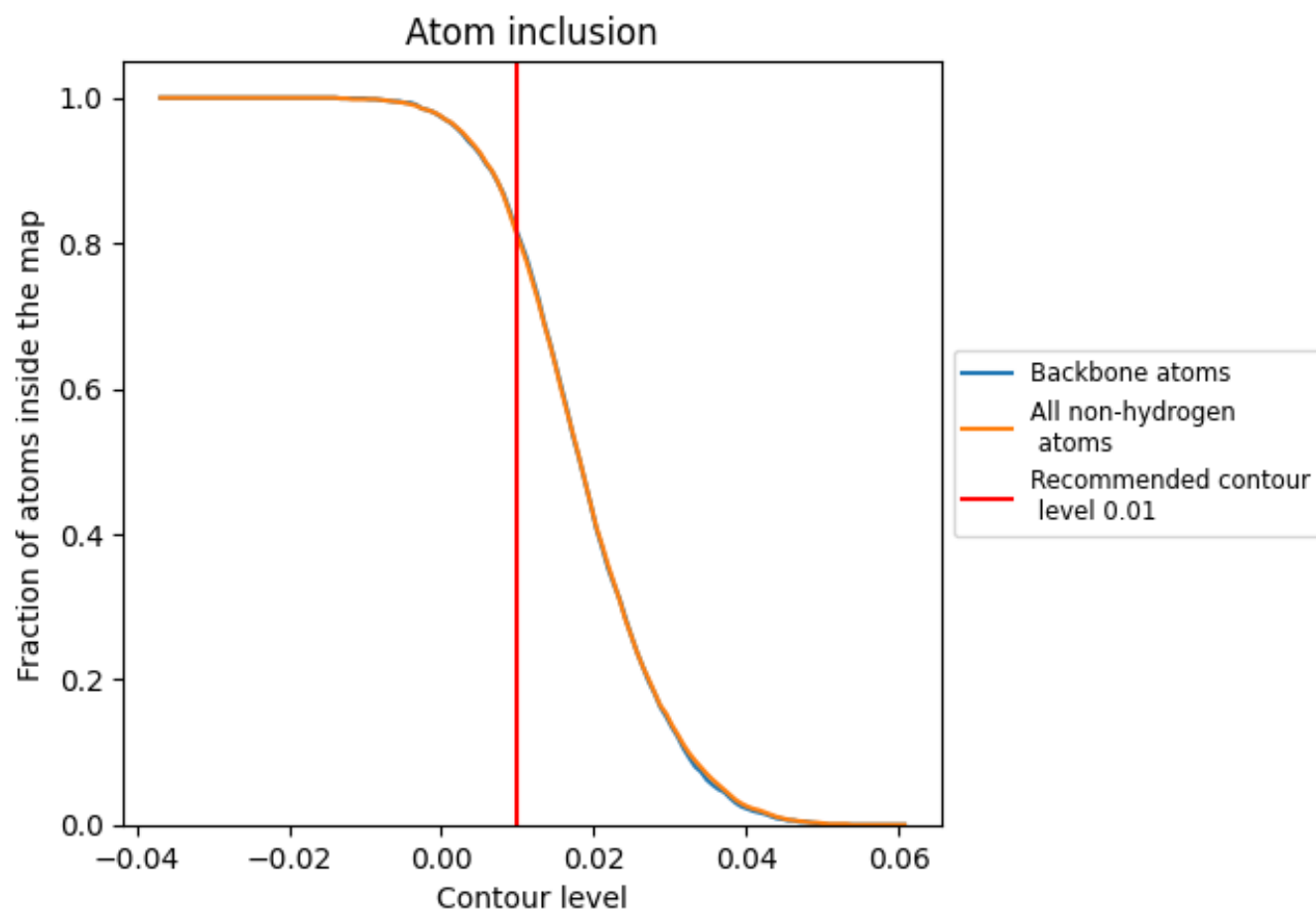
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.01 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.

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 81% 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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8090	0.2060
A	0.8600	0.2300
B	0.8620	0.2490
C	0.8180	0.2100
D	0.7990	0.1910
E	0.8180	0.2090
F	0.7910	0.1800
G	0.8670	0.2460
H	0.8640	0.2440
I	0.8170	0.2050
J	0.8040	0.1860
K	0.8160	0.2110
L	0.7930	0.1820

1.0

0.0

<0.0