



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 01:16 PM UTC

PDB ID : 9F44 / pdb_00009f44
EMDB ID : EMD-50183
Title : cryo-EM structure of LST2 TOS peptide bound to human mTOR complex 1
Authors : Craigie, L.M.; Maier, T.
Deposited on : 2024-04-26
Resolution : 3.68 Å(reported)
Based on initial model : 7PEB

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

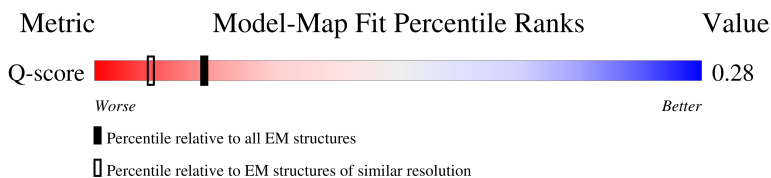
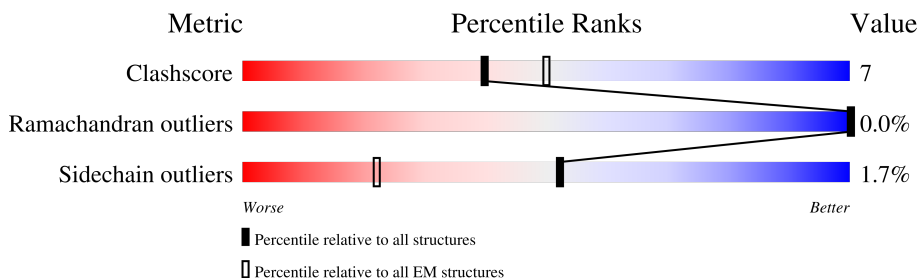
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.68 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.



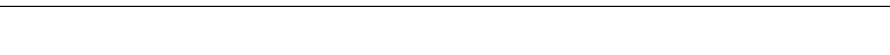
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11376 (3.18 - 4.18)

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	2549	11% 70% 16% 14%
1	B	2549	11% 70% 16% 14%
2	C	326	5% 81% 16% .
2	D	326	5% 82% 15% .

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Mol	Chain	Length	Quality of chain
3	E	1363	 5% 62% 15% 23%
3	F	1363	 5% 61% 16% 23%
4	G	13	 8% 54% 15% 31%
4	H	13	 8% 54% 15% 31%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 56848 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 Serine/threonine-protein kinase mTOR.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	2192	Total	C	N	O	S	0	0
			17467	11158	3060	3137	112		
1	B	2192	Total	C	N	O	S	0	0
			17467	11158	3060	3137	112		

- Molecule 2 is a protein called Target of rapamycin complex subunit LST8.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	317	Total	C	N	O	S	0	0
			2456	1526	436	476	18		
2	D	317	Total	C	N	O	S	0	0
			2456	1526	436	476	18		

- Molecule 3 is a protein called Regulatory-associated protein of mTOR.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	1052	Total	C	N	O	S	0	0
			8385	5361	1450	1518	56		
3	F	1052	Total	C	N	O	S	0	0
			8385	5361	1450	1518	56		

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-27	HIS	-	expression tag	UNP Q8N122
E	-26	HIS	-	expression tag	UNP Q8N122
E	-25	HIS	-	expression tag	UNP Q8N122
E	-24	HIS	-	expression tag	UNP Q8N122
E	-23	HIS	-	expression tag	UNP Q8N122
E	-22	HIS	-	expression tag	UNP Q8N122
E	-21	HIS	-	expression tag	UNP Q8N122
E	-20	HIS	-	expression tag	UNP Q8N122
E	-19	HIS	-	expression tag	UNP Q8N122

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-18	HIS	-	expression tag	UNP Q8N122
E	-17	GLU	-	expression tag	UNP Q8N122
E	-16	GLN	-	expression tag	UNP Q8N122
E	-15	LYS	-	expression tag	UNP Q8N122
E	-14	LEU	-	expression tag	UNP Q8N122
E	-13	ILE	-	expression tag	UNP Q8N122
E	-12	SER	-	expression tag	UNP Q8N122
E	-11	GLU	-	expression tag	UNP Q8N122
E	-10	GLU	-	expression tag	UNP Q8N122
E	-9	ASP	-	expression tag	UNP Q8N122
E	-8	LEU	-	expression tag	UNP Q8N122
E	-7	ASP	-	expression tag	UNP Q8N122
E	-6	TYR	-	expression tag	UNP Q8N122
E	-5	LYS	-	expression tag	UNP Q8N122
E	-4	ASP	-	expression tag	UNP Q8N122
E	-3	ASP	-	expression tag	UNP Q8N122
E	-2	ASP	-	expression tag	UNP Q8N122
E	-1	ASP	-	expression tag	UNP Q8N122
E	0	LYS	-	expression tag	UNP Q8N122
F	-27	HIS	-	expression tag	UNP Q8N122
F	-26	HIS	-	expression tag	UNP Q8N122
F	-25	HIS	-	expression tag	UNP Q8N122
F	-24	HIS	-	expression tag	UNP Q8N122
F	-23	HIS	-	expression tag	UNP Q8N122
F	-22	HIS	-	expression tag	UNP Q8N122
F	-21	HIS	-	expression tag	UNP Q8N122
F	-20	HIS	-	expression tag	UNP Q8N122
F	-19	HIS	-	expression tag	UNP Q8N122
F	-18	HIS	-	expression tag	UNP Q8N122
F	-17	GLU	-	expression tag	UNP Q8N122
F	-16	GLN	-	expression tag	UNP Q8N122
F	-15	LYS	-	expression tag	UNP Q8N122
F	-14	LEU	-	expression tag	UNP Q8N122
F	-13	ILE	-	expression tag	UNP Q8N122
F	-12	SER	-	expression tag	UNP Q8N122
F	-11	GLU	-	expression tag	UNP Q8N122
F	-10	GLU	-	expression tag	UNP Q8N122
F	-9	ASP	-	expression tag	UNP Q8N122
F	-8	LEU	-	expression tag	UNP Q8N122
F	-7	ASP	-	expression tag	UNP Q8N122
F	-6	TYR	-	expression tag	UNP Q8N122
F	-5	LYS	-	expression tag	UNP Q8N122

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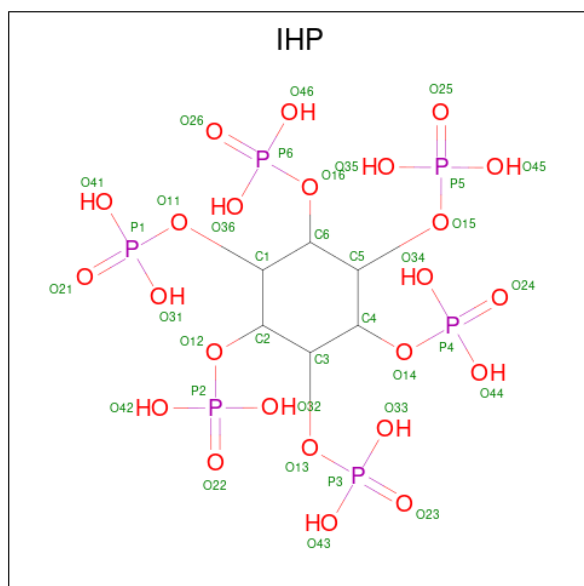
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Chain	Residue	Modelled	Actual	Comment	Reference
F	-4	ASP	-	expression tag	UNP Q8N122
F	-3	ASP	-	expression tag	UNP Q8N122
F	-2	ASP	-	expression tag	UNP Q8N122
F	-1	ASP	-	expression tag	UNP Q8N122
F	0	LYS	-	expression tag	UNP Q8N122

- Molecule 4 is a protein called Lateral signaling target protein 2 homolog.

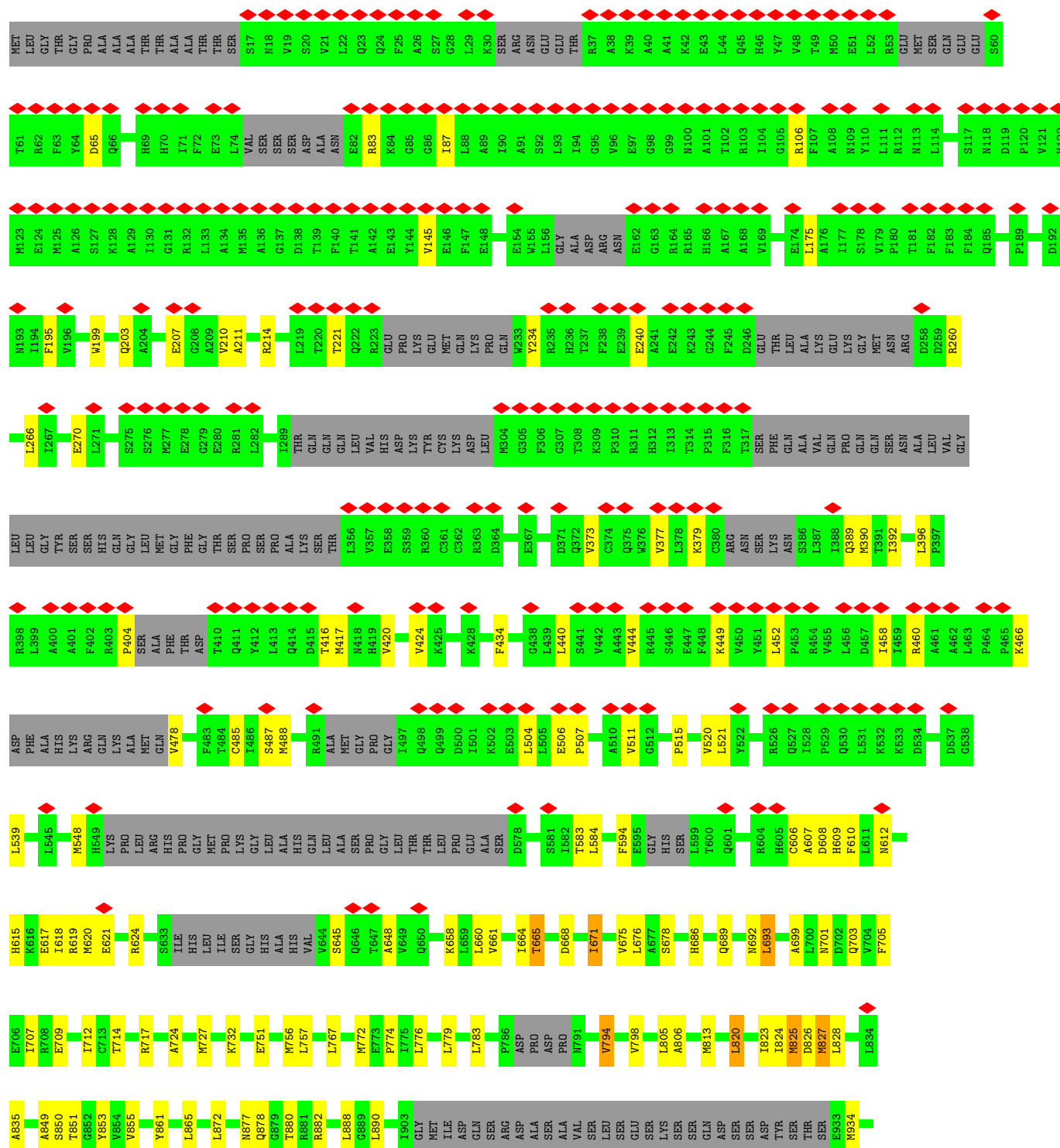
Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	9	Total	C	N	O	S	0	0
			80	52	12	15	1		
4	H	9	Total	C	N	O	S	0	0
			80	52	12	15	1		

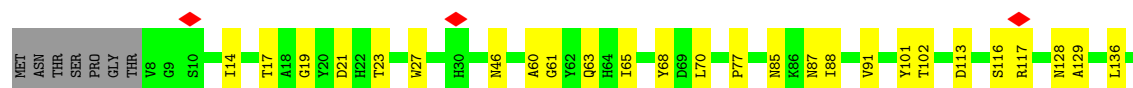
- Molecule 5 is INOSITOL HEXAKISPHOSPHATE (CCD ID: IHP) (formula: $C_6H_{18}O_{24}P_6$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	O	P	0
			36	6	24	6	
5	B	1	Total	C	O	P	0
			36	6	24	6	









4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	782845	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	48.43	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.959	Depositor
Minimum map value	-0.449	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.014	Depositor
Recommended contour level	0.13	Depositor
Map size (Å)	670.72, 670.72, 670.72	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.31, 1.31, 1.31	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.12	0/17814	0.28	0/24116
1	B	0.12	0/17814	0.28	0/24116
2	C	0.11	0/2514	0.29	0/3426
2	D	0.11	0/2514	0.29	0/3426
3	E	0.10	0/8585	0.25	0/11680
3	F	0.10	0/8585	0.25	0/11680
4	G	0.11	0/81	0.27	0/107
4	H	0.11	0/81	0.27	0/107
All	All	0.11	0/57988	0.27	0/78658

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	17467	0	17553	267	0
1	B	17467	0	17553	264	0
2	C	2456	0	2341	30	0
2	D	2456	0	2341	28	0
3	E	8385	0	8375	125	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	8385	0	8375	134	0
4	G	80	0	71	2	0
4	H	80	0	71	2	0
5	A	36	0	6	2	0
5	B	36	0	6	2	0
All	All	56848	0	56692	808	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 7.

All (808) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:506:GLU:HG3	1:B:507:PRO:HD3	1.65	0.79
3:F:660:ARG:NH2	3:F:1111:ALA:O	2.16	0.78
1:A:506:GLU:HG3	1:A:507:PRO:HD3	1.65	0.77
3:E:660:ARG:NH2	3:E:1111:ALA:O	2.16	0.77
1:A:1943:ILE:HA	1:A:1946:ILE:HD13	1.66	0.76
1:B:1943:ILE:HA	1:B:1946:ILE:HD13	1.66	0.76
1:B:1745:LYS:HE3	1:B:1749:ARG:HE	1.52	0.74
3:E:1252:ILE:HG22	3:E:1253:VAL:HG23	1.71	0.73
3:F:1252:ILE:HG22	3:F:1253:VAL:HG23	1.71	0.73
1:A:1745:LYS:HE3	1:A:1749:ARG:HE	1.52	0.71
1:A:606:CYS:HA	1:A:609:HIS:HB3	1.73	0.70
1:B:606:CYS:HA	1:B:609:HIS:HB3	1.73	0.69
1:A:878:GLN:NE2	1:A:1572:ASP:OD2	2.25	0.69
2:D:166:ILE:HD11	2:D:169:PRO:HA	1.74	0.68
1:B:487:SER:HB3	1:B:520:VAL:HG13	1.76	0.68
2:C:166:ILE:HD11	2:C:169:PRO:HA	1.74	0.68
1:A:2222:ILE:HD11	1:A:2330:VAL:HG21	1.76	0.68
2:D:101:TYR:HE2	2:D:136:LEU:HD13	1.59	0.68
1:B:2222:ILE:HD11	1:B:2330:VAL:HG21	1.76	0.68
1:A:2277:HIS:O	2:C:87:ASN:ND2	2.27	0.67
2:C:101:TYR:HE2	2:C:136:LEU:HD13	1.59	0.67
1:B:878:GLN:NE2	1:B:1572:ASP:OD2	2.25	0.67
1:A:1987:ARG:HH22	1:A:2426:LEU:HD11	1.60	0.67
1:B:2277:HIS:O	2:D:87:ASN:ND2	2.26	0.67
1:A:487:SER:HB3	1:A:520:VAL:HG13	1.76	0.67
1:A:1150:ASP:OD1	1:B:701:ASN:ND2	2.27	0.67
1:B:1987:ARG:HH22	1:B:2426:LEU:HD11	1.60	0.67
1:A:2271:MET:HE3	1:A:2286:VAL:HG13	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:701:ASN:ND2	1:B:1150:ASP:OD1	2.27	0.66
1:B:2271:MET:HE3	1:B:2286:VAL:HG13	1.77	0.65
2:C:19:GLY:HA2	2:C:316:VAL:HG22	1.79	0.65
1:B:2408:ARG:NH2	1:B:2524:GLN:OE1	2.30	0.64
1:A:1174:MET:HE3	1:A:1204:ILE:HG21	1.80	0.64
1:A:2408:ARG:NH2	1:A:2524:GLN:OE1	2.30	0.64
1:B:2018:ARG:NH1	1:B:2067:GLU:OE2	2.30	0.64
1:B:877:ASN:HB3	1:B:880:THR:HG22	1.80	0.64
1:A:2018:ARG:NH1	1:A:2067:GLU:OE2	2.30	0.64
1:B:1174:MET:HE3	1:B:1204:ILE:HG21	1.80	0.63
2:D:19:GLY:HA2	2:D:316:VAL:HG22	1.79	0.63
1:A:779:LEU:HD22	1:A:798:VAL:HG23	1.80	0.63
1:A:877:ASN:HB3	1:A:880:THR:HG22	1.80	0.63
1:A:938:MET:HE1	1:A:1366:HIS:HB3	1.80	0.63
1:A:2288:GLU:OE1	2:C:221:ARG:NH2	2.32	0.63
1:B:938:MET:HE1	1:B:1366:HIS:HB3	1.80	0.62
3:E:413:LEU:HD21	3:E:439:VAL:HG11	1.80	0.62
1:A:65:ASP:OD2	1:A:106:ARG:NH1	2.32	0.62
1:A:608:ASP:OD1	1:A:612:ASN:ND2	2.32	0.62
1:A:1552:LEU:O	1:A:1606:LYS:NZ	2.33	0.62
1:B:2288:GLU:OE1	2:D:221:ARG:NH2	2.32	0.62
3:F:413:LEU:HD21	3:F:439:VAL:HG11	1.80	0.62
1:B:65:ASP:OD2	1:B:106:ARG:NH1	2.32	0.62
1:B:779:LEU:HD22	1:B:798:VAL:HG23	1.80	0.62
1:B:1552:LEU:O	1:B:1606:LYS:NZ	2.33	0.62
1:B:2016:LEU:HD12	1:B:2186:LEU:HD21	1.82	0.61
3:E:1139:MET:SD	3:E:1139:MET:N	2.73	0.61
1:A:645:SER:HB3	1:A:648:ALA:HB3	1.82	0.61
3:F:1139:MET:SD	3:F:1139:MET:N	2.73	0.61
1:A:583:THR:HG23	1:A:621:GLU:HG3	1.82	0.61
1:B:645:SER:HB3	1:B:648:ALA:HB3	1.82	0.61
3:E:681:VAL:HG21	3:E:808:TYR:HB2	1.82	0.61
1:B:608:ASP:OD1	1:B:612:ASN:ND2	2.32	0.61
2:D:229:SER:OG	2:D:231:ASP:OD1	2.18	0.61
3:F:184:THR:OG1	3:F:213:ARG:NH2	2.34	0.61
1:A:751:GLU:HG3	1:A:794:VAL:HG22	1.83	0.61
1:A:2016:LEU:HD12	1:A:2186:LEU:HD21	1.82	0.61
1:B:2170:ARG:HB2	1:B:2186:LEU:HB2	1.83	0.61
3:E:558:LEU:HD22	3:E:584:ILE:HG13	1.83	0.61
3:E:184:THR:OG1	3:E:213:ARG:NH2	2.34	0.60
1:B:583:THR:HG23	1:B:621:GLU:HG3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:888:LEU:HD22	1:B:1590:MET:HE1	1.83	0.60
3:F:681:VAL:HG21	3:F:808:TYR:HB2	1.82	0.60
2:C:117:ARG:O	2:C:117:ARG:NE	2.33	0.60
1:A:1514:ARG:NH1	1:A:1542:ASP:OD2	2.35	0.59
1:B:1514:ARG:NH1	1:B:1542:ASP:OD2	2.35	0.59
2:D:46:ASN:N	2:D:60:ALA:O	2.31	0.59
3:E:592:ARG:NH1	3:E:626:PHE:O	2.36	0.59
3:F:1281:ASN:OD1	3:F:1282:SER:N	2.35	0.59
2:C:46:ASN:N	2:C:60:ALA:O	2.31	0.59
3:F:558:LEU:HD22	3:F:584:ILE:HG13	1.83	0.59
1:A:2170:ARG:HB2	1:A:2186:LEU:HB2	1.83	0.59
3:F:440:LEU:HD12	3:F:472:ILE:HD11	1.85	0.59
3:E:275:LEU:HD21	3:E:312:LEU:HD13	1.85	0.59
3:E:440:LEU:HD12	3:E:472:ILE:HD11	1.85	0.59
1:B:2368:ARG:NH1	1:B:2370:LYS:O	2.36	0.59
1:A:888:LEU:HD22	1:A:1590:MET:HE1	1.83	0.59
1:A:1188:TYR:HH	1:A:1209:TYR:HH	1.48	0.59
3:F:592:ARG:NH1	3:F:626:PHE:O	2.36	0.59
1:B:751:GLU:HG3	1:B:794:VAL:HG22	1.83	0.58
3:F:1083:LEU:HD23	3:F:1097:LYS:HD3	1.85	0.58
1:B:619:ARG:NH2	1:B:668:ASP:OD2	2.36	0.58
3:F:559:ILE:HD11	3:F:584:ILE:HG21	1.85	0.58
1:B:1444:GLU:OE1	1:B:1444:GLU:N	2.36	0.58
3:E:1083:LEU:HD23	3:E:1097:LYS:HD3	1.85	0.58
1:A:686:HIS:O	1:A:692:ASN:ND2	2.36	0.58
1:A:2332:TYR:OH	1:A:2512:ASP:OD2	2.20	0.58
1:A:2368:ARG:NH1	1:A:2370:LYS:O	2.36	0.58
1:B:1164:ASP:OD1	1:B:1202:HIS:NE2	2.28	0.58
1:A:1444:GLU:N	1:A:1444:GLU:OE1	2.36	0.58
1:A:1684:GLN:HB3	1:A:1687:HIS:HB3	1.85	0.58
3:E:1077:ASN:OD1	3:E:1133:GLN:NE2	2.34	0.58
1:A:211:ALA:HA	1:A:214:ARG:HD3	1.86	0.58
2:D:91:VAL:HG12	2:D:102:THR:HG22	1.86	0.58
2:D:152:ILE:N	2:D:165:LEU:O	2.36	0.58
1:A:619:ARG:NH2	1:A:668:ASP:OD2	2.36	0.57
1:A:850:SER:O	1:A:1611:ARG:NH2	2.37	0.57
1:B:686:HIS:O	1:B:692:ASN:ND2	2.36	0.57
3:F:1062:ASN:O	3:F:1094:ARG:NH1	2.37	0.57
3:F:275:LEU:HD21	3:F:312:LEU:HD13	1.85	0.57
1:B:676:LEU:HD22	1:B:707:ILE:HG23	1.86	0.57
1:B:1684:GLN:HB3	1:B:1687:HIS:HB3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:1077:ASN:OD1	3:F:1133:GLN:NE2	2.34	0.57
1:B:2022:LEU:HD11	1:B:2121:LEU:HD21	1.87	0.57
1:B:2332:TYR:OH	1:B:2512:ASP:OD2	2.20	0.57
3:E:1062:ASN:O	3:E:1094:ARG:NH1	2.37	0.57
1:B:1697:THR:HG22	1:B:1701:MET:HE2	1.87	0.57
3:E:266:CYS:HA	3:E:273:ILE:HG21	1.87	0.57
3:E:559:ILE:HD11	3:E:584:ILE:HG21	1.85	0.57
1:A:1164:ASP:OD1	1:A:1202:HIS:NE2	2.28	0.57
1:B:211:ALA:HA	1:B:214:ARG:HD3	1.86	0.57
2:C:91:VAL:HG12	2:C:102:THR:HG22	1.86	0.57
3:E:1281:ASN:OD1	3:E:1282:SER:N	2.35	0.57
3:F:266:CYS:HA	3:F:273:ILE:HG21	1.87	0.57
1:A:676:LEU:HD22	1:A:707:ILE:HG23	1.86	0.56
1:B:850:SER:O	1:B:1611:ARG:NH2	2.37	0.56
3:E:1138:LEU:O	3:E:1150:TRP:N	2.38	0.56
1:A:1631:GLU:OE1	1:A:1631:GLU:N	2.37	0.56
1:A:2022:LEU:HD11	1:A:2121:LEU:HD21	1.87	0.56
2:D:117:ARG:O	2:D:117:ARG:NE	2.33	0.56
3:F:494:ILE:O	3:F:498:ILE:HG12	2.06	0.56
1:B:1631:GLU:N	1:B:1631:GLU:OE1	2.37	0.56
3:F:1138:LEU:O	3:F:1150:TRP:N	2.38	0.56
3:E:494:ILE:O	3:E:498:ILE:HG12	2.05	0.56
1:A:1697:THR:HG22	1:A:1701:MET:HE2	1.87	0.56
1:B:1119:VAL:HA	1:B:1122:PHE:CE2	2.41	0.56
1:A:701:ASN:ND2	1:B:1150:ASP:O	2.39	0.55
1:A:1119:VAL:HA	1:A:1122:PHE:CE2	2.41	0.55
1:A:1598:GLU:OE2	1:A:1640:ARG:NE	2.37	0.55
1:B:2339:ARG:NH2	1:B:2343:ASN:HB3	2.21	0.55
2:D:63:GLN:HG2	2:D:85:ASN:HA	1.89	0.55
2:C:152:ILE:N	2:C:165:LEU:O	2.36	0.55
1:A:377:VAL:O	1:A:389:GLN:NE2	2.39	0.55
1:A:2339:ARG:NH2	1:A:2343:ASN:HB3	2.21	0.55
1:B:1331:GLU:HA	1:B:1334:GLN:HB2	1.89	0.55
1:B:377:VAL:O	1:B:389:GLN:NE2	2.39	0.54
2:C:229:SER:OG	2:C:231:ASP:OD1	2.18	0.54
1:A:515:PRO:HD3	1:A:584:LEU:HD11	1.90	0.54
1:B:1724:MET:HE2	1:B:1747:MET:HB2	1.89	0.54
1:A:727:MET:HG3	1:A:767:LEU:HD23	1.90	0.54
1:B:2337:GLY:O	1:B:2339:ARG:NH1	2.40	0.54
2:C:63:GLN:HG2	2:C:85:ASN:HA	1.89	0.54
1:A:952:MET:HE2	1:A:993:VAL:HG22	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1277:LYS:HA	1:A:1350:ILE:HD13	1.90	0.54
1:A:1527:ASP:N	1:A:1527:ASP:OD1	2.40	0.54
1:B:417:MET:HE1	1:B:440:LEU:HD13	1.89	0.54
1:B:952:MET:HE2	1:B:993:VAL:HG22	1.90	0.54
1:B:1277:LYS:HA	1:B:1350:ILE:HD13	1.90	0.54
1:A:1150:ASP:O	1:B:701:ASN:ND2	2.40	0.54
1:A:2337:GLY:O	1:A:2339:ARG:NH1	2.40	0.54
1:B:515:PRO:HD3	1:B:584:LEU:HD11	1.90	0.54
1:B:1598:GLU:OE2	1:B:1640:ARG:NE	2.37	0.54
3:E:1032:HIS:CG	3:E:1035:THR:HG22	2.43	0.54
1:A:1331:GLU:HA	1:A:1334:GLN:HB2	1.89	0.54
1:B:890:LEU:HD12	1:B:1597:SER:HB2	1.90	0.54
2:C:173:ILE:HD11	2:C:187:ALA:HB1	1.90	0.54
3:E:151:ASN:OD1	3:E:153:HIS:ND1	2.41	0.54
3:F:1032:HIS:CG	3:F:1035:THR:HG22	2.43	0.54
1:A:1342:GLU:OE2	1:A:1375:ARG:N	2.41	0.53
1:B:945:GLU:OE1	1:B:945:GLU:N	2.40	0.53
1:B:727:MET:HG3	1:B:767:LEU:HD23	1.90	0.53
1:A:620:MET:HE2	1:A:624:ARG:HG3	1.89	0.53
1:B:620:MET:HE2	1:B:624:ARG:HG3	1.89	0.53
2:D:173:ILE:HD11	2:D:187:ALA:HB1	1.90	0.53
1:A:417:MET:HE1	1:A:440:LEU:HD13	1.89	0.53
1:A:1109:ASP:CG	3:E:282:LYS:HZ3	2.16	0.53
1:A:1724:MET:HE2	1:A:1747:MET:HB2	1.90	0.53
1:A:701:ASN:O	1:B:1153:SER:OG	2.26	0.53
1:B:958:ILE:HD12	1:B:1306:LEU:HD21	1.91	0.53
1:A:945:GLU:N	1:A:945:GLU:OE1	2.40	0.53
1:B:772:MET:HG2	1:B:805:LEU:HD11	1.90	0.53
1:B:1527:ASP:OD1	1:B:1527:ASP:N	2.40	0.53
3:E:530:GLU:OE2	3:E:570:HIS:NE2	2.42	0.53
3:F:1088:THR:OG1	3:F:1090:ASP:OD1	2.21	0.53
1:A:689:GLN:HE22	3:F:419:TRP:HA	1.75	0.53
1:A:1629:ILE:HG22	1:A:1632:ASP:H	1.74	0.52
1:A:890:LEU:HD12	1:A:1597:SER:HB2	1.90	0.52
1:A:2194:GLN:HE21	1:A:2198:VAL:HG23	1.74	0.52
1:B:990:LEU:HD21	1:B:1023:VAL:HG21	1.92	0.52
1:B:2084:TRP:HB2	1:B:2100:ALA:HB2	1.92	0.52
1:A:806:ALA:HA	1:A:813:MET:SD	2.50	0.52
1:A:1153:SER:OG	1:B:701:ASN:O	2.26	0.52
1:A:1892:ILE:HG21	1:A:1930:ILE:HD11	1.92	0.52
1:B:2194:GLN:HE21	1:B:2198:VAL:HG23	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:147:LEU:HD12	3:E:191:ILE:HD11	1.91	0.52
3:F:151:ASN:OD1	3:F:153:HIS:ND1	2.41	0.52
3:F:509:LEU:HD13	3:F:542:ILE:HD11	1.92	0.52
1:B:221:THR:OG1	1:B:234:TYR:OH	2.26	0.52
1:B:806:ALA:HA	1:B:813:MET:SD	2.50	0.52
3:F:446:ARG:HD2	3:F:479:LEU:HD11	1.92	0.52
3:E:446:ARG:HD2	3:E:479:LEU:HD11	1.92	0.52
1:B:1329:LEU:HB2	1:B:1334:GLN:HG3	1.92	0.52
3:E:1088:THR:OG1	3:E:1090:ASP:OD1	2.21	0.52
1:A:1745:LYS:HE3	1:A:1749:ARG:NE	2.24	0.52
1:B:1892:ILE:HG21	1:B:1930:ILE:HD11	1.92	0.52
3:F:147:LEU:HD12	3:F:191:ILE:HD11	1.92	0.52
3:F:530:GLU:OE2	3:F:570:HIS:NE2	2.42	0.52
1:A:210:VAL:HG11	1:A:266:LEU:HD13	1.92	0.52
1:A:1066:GLU:OE1	1:A:1066:GLU:N	2.42	0.52
1:A:1161:ARG:HH22	1:B:703:GLN:NE2	2.08	0.52
1:B:1066:GLU:N	1:B:1066:GLU:OE1	2.42	0.51
1:B:1167:PRO:HA	1:B:1170:ARG:HG3	1.92	0.51
3:F:979:ASP:OD2	3:F:982:SER:HB3	2.10	0.51
1:A:1598:GLU:O	1:A:1602:VAL:HG23	2.10	0.51
1:A:772:MET:HG2	1:A:805:LEU:HD11	1.91	0.51
1:B:1533:THR:HA	1:B:1536:ILE:HG13	1.92	0.51
3:E:1262:HIS:HD2	3:E:1263:PRO:HD2	1.75	0.51
1:B:266:LEU:O	1:B:270:GLU:HG2	2.10	0.51
3:E:244:ALA:HB3	3:E:365:VAL:HB	1.93	0.51
1:A:1167:PRO:HA	1:A:1170:ARG:HG3	1.92	0.51
1:A:1533:THR:HA	1:A:1536:ILE:HG13	1.91	0.51
1:B:210:VAL:HG11	1:B:266:LEU:HD13	1.92	0.51
1:B:1342:GLU:OE2	1:B:1375:ARG:N	2.41	0.51
3:E:294:LEU:HD11	3:E:393:LEU:HB3	1.93	0.51
1:A:958:ILE:HD12	1:A:1306:LEU:HD21	1.91	0.51
1:A:990:LEU:HD21	1:A:1023:VAL:HG21	1.92	0.51
1:A:1329:LEU:HB2	1:A:1334:GLN:HG3	1.92	0.51
3:E:509:LEU:HD13	3:E:542:ILE:HD11	1.92	0.51
3:E:979:ASP:OD2	3:E:982:SER:HB3	2.10	0.51
1:B:689:GLN:HE22	3:E:419:TRP:HA	1.75	0.51
1:A:2084:TRP:HB2	1:A:2100:ALA:HB2	1.92	0.51
1:B:2271:MET:HE1	1:B:2289:HIS:HB3	1.92	0.51
3:F:566:LEU:HD23	3:F:603:LYS:HG2	1.93	0.51
1:A:266:LEU:O	1:A:270:GLU:HG2	2.10	0.51
1:A:703:GLN:NE2	1:B:1161:ARG:HH22	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2077:ASP:HB3	1:A:2103:LEU:HD22	1.92	0.51
1:A:732:LYS:NZ	1:B:1150:ASP:OD2	2.34	0.50
1:A:2364:VAL:HG22	1:A:2434:THR:HG21	1.93	0.50
1:B:1629:ILE:HG22	1:B:1632:ASP:H	1.74	0.50
1:A:2188:GLY:O	1:A:2189:HIS:ND1	2.44	0.50
3:F:294:LEU:HD11	3:F:393:LEU:HB3	1.93	0.50
1:B:2188:GLY:O	1:B:2189:HIS:ND1	2.44	0.50
3:E:413:LEU:HD12	3:E:452:LEU:HD12	1.94	0.50
1:A:1199:LEU:O	1:A:1203:ARG:N	2.45	0.50
1:B:615:HIS:HB3	1:B:618:ILE:HB	1.94	0.50
1:B:1658:SER:OG	5:B:2601:IHP:O32	2.29	0.50
1:B:2275:TYR:HA	1:B:2278:LEU:HG	1.93	0.50
3:F:499:LEU:HB3	3:F:541:VAL:HG21	1.94	0.50
1:A:2275:TYR:HA	1:A:2278:LEU:HG	1.93	0.50
1:B:2387:MET:HE1	1:B:2392:LEU:HA	1.93	0.50
1:A:615:HIS:HB3	1:A:618:ILE:HB	1.94	0.50
1:B:1109:ASP:CG	3:F:282:LYS:HZ3	2.19	0.50
3:F:1189:SER:OG	3:F:1191:ARG:NH1	2.45	0.50
3:F:1262:HIS:HD2	3:F:1263:PRO:HD2	1.75	0.50
1:A:2291:VAL:HG13	1:A:2381:ARG:HE	1.77	0.50
1:B:1598:GLU:O	1:B:1602:VAL:HG23	2.10	0.50
1:B:2077:ASP:HB3	1:B:2103:LEU:HD22	1.92	0.50
3:F:244:ALA:HB3	3:F:365:VAL:HB	1.93	0.50
1:A:2271:MET:HE1	1:A:2289:HIS:HB3	1.92	0.50
1:B:1064:GLY:O	1:B:1106:ASN:ND2	2.44	0.50
1:B:1199:LEU:O	1:B:1203:ARG:N	2.45	0.50
1:B:1318:PHE:CE2	1:B:1360:LEU:HD12	2.47	0.50
3:E:131:LYS:HZ3	3:E:185:TRP:CD1	2.28	0.50
2:D:223:ALA:HA	2:D:239:SER:HA	1.94	0.50
3:E:464:VAL:HB	3:E:501:VAL:HG21	1.94	0.50
3:F:413:LEU:HD12	3:F:452:LEU:HD12	1.94	0.50
1:A:1728:ALA:O	1:A:1732:ILE:HG13	2.12	0.49
1:A:2374:LYS:O	1:A:2537:ASN:ND2	2.45	0.49
3:E:499:LEU:HB3	3:E:541:VAL:HG21	1.94	0.49
3:E:1060:PHE:CZ	3:E:1106:PRO:HB3	2.47	0.49
3:E:1189:SER:OG	3:E:1191:ARG:NH1	2.45	0.49
1:A:1318:PHE:CE2	1:A:1360:LEU:HD12	2.46	0.49
1:A:2010:MET:O	1:A:2014:GLU:HG2	2.12	0.49
1:B:994:MET:HE2	1:B:1027:ILE:HD11	1.95	0.49
1:B:1084:HIS:O	1:B:1084:HIS:ND1	2.46	0.49
1:B:1323:VAL:HG21	1:B:1363:PHE:HE2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1655:LYS:NZ	5:B:2601:IHP:O43	2.46	0.49
3:F:131:LYS:HZ3	3:F:185:TRP:CD1	2.31	0.49
2:C:223:ALA:HA	2:C:239:SER:HA	1.95	0.49
2:D:147:ASP:OD1	2:D:151:ALA:N	2.46	0.49
1:A:776:LEU:HD11	1:A:820:LEU:HD21	1.94	0.49
1:A:2022:LEU:O	1:A:2026:MET:HG3	2.13	0.49
1:A:2387:MET:HE1	1:A:2392:LEU:HA	1.93	0.49
1:B:757:LEU:HD23	1:B:779:LEU:HD21	1.95	0.49
1:A:2192:LEU:HD12	1:A:2235:GLY:HA3	1.95	0.49
1:B:466:LYS:HZ1	1:B:478:VAL:N	2.11	0.49
1:B:2364:VAL:HG22	1:B:2434:THR:HG21	1.93	0.49
3:E:53:ASP:N	3:E:53:ASP:OD1	2.46	0.49
3:F:53:ASP:OD1	3:F:53:ASP:N	2.46	0.49
3:F:395:THR:O	3:F:400:GLY:N	2.40	0.49
1:A:671:ILE:O	1:A:675:VAL:HG12	2.13	0.49
1:B:2022:LEU:O	1:B:2026:MET:HG3	2.12	0.49
2:C:300:GLU:HG2	2:C:301:THR:HG23	1.95	0.49
1:A:1084:HIS:O	1:A:1084:HIS:ND1	2.46	0.49
1:B:671:ILE:O	1:B:675:VAL:HG12	2.13	0.49
1:B:2374:LYS:O	1:B:2537:ASN:ND2	2.45	0.49
3:E:637:SER:HA	3:E:640:ILE:HG22	1.95	0.49
3:F:1060:PHE:CZ	3:F:1106:PRO:HB3	2.47	0.49
1:A:221:THR:OG1	1:A:234:TYR:OH	2.26	0.49
1:B:2088:TYR:O	1:B:2092:GLY:N	2.45	0.49
2:D:65:ILE:HD11	2:D:88:ILE:HG21	1.95	0.49
3:E:180:TYR:O	3:E:184:THR:OG1	2.26	0.49
3:F:476:VAL:HG13	3:F:495:TRP:CZ2	2.48	0.49
1:A:2088:TYR:O	1:A:2092:GLY:N	2.45	0.48
1:B:2291:VAL:HG13	1:B:2381:ARG:HE	1.77	0.48
2:D:300:GLU:HG2	2:D:301:THR:HG23	1.95	0.48
1:B:661:VAL:O	1:B:665:THR:OG1	2.24	0.48
1:B:2010:MET:O	1:B:2014:GLU:HG2	2.13	0.48
3:E:566:LEU:HD23	3:E:603:LYS:HG2	1.93	0.48
1:A:872:LEU:HD13	1:A:1575:LEU:HD22	1.96	0.48
1:B:1199:LEU:HD22	1:B:1204:ILE:HB	1.95	0.48
3:F:292:LEU:HD12	3:F:292:LEU:H	1.78	0.48
1:A:757:LEU:HD23	1:A:779:LEU:HD21	1.95	0.48
1:A:994:MET:HE2	1:A:1027:ILE:HD11	1.95	0.48
1:A:1323:VAL:HG21	1:A:1363:PHE:HE2	1.78	0.48
1:B:1745:LYS:HE3	1:B:1749:ARG:NE	2.24	0.48
2:C:147:ASP:OD1	2:C:151:ALA:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:292:LEU:HD12	3:E:292:LEU:H	1.78	0.48
3:E:1226:HIS:ND1	3:E:1240:ASP:OD1	2.46	0.48
3:F:166:VAL:HG23	3:F:175:ILE:HG13	1.96	0.48
3:F:1226:HIS:ND1	3:F:1240:ASP:OD1	2.46	0.48
1:B:872:LEU:HD13	1:B:1575:LEU:HD22	1.96	0.48
3:E:128:ASP:OD1	3:E:129:GLU:N	2.47	0.48
3:E:305:ARG:NH2	4:G:402:PHE:O	2.46	0.48
3:F:464:VAL:HB	3:F:501:VAL:HG21	1.95	0.48
1:A:424:VAL:HG12	1:A:458:ILE:HG21	1.96	0.48
1:A:607:ALA:HA	1:A:610:PHE:CE2	2.49	0.48
1:B:776:LEU:HD11	1:B:820:LEU:HD21	1.94	0.48
1:B:1724:MET:SD	1:B:1743:LEU:HG	2.54	0.48
3:E:126:THR:OG1	3:E:128:ASP:OD1	2.20	0.48
1:B:607:ALA:HA	1:B:610:PHE:CE2	2.49	0.48
3:E:277:TRP:HA	3:E:280:MET:HE2	1.96	0.48
3:F:305:ARG:NH2	4:H:402:PHE:O	2.46	0.48
3:F:611:PRO:HB3	3:F:1154:ARG:HE	1.79	0.48
3:F:637:SER:HA	3:F:640:ILE:HG22	1.95	0.48
1:A:1724:MET:SD	1:A:1743:LEU:HG	2.54	0.48
1:B:1728:ALA:O	1:B:1732:ILE:HG13	2.12	0.48
3:E:476:VAL:HG13	3:E:495:TRP:CZ2	2.48	0.48
3:F:49:TRP:CD2	3:F:468:LEU:HD12	2.49	0.48
1:A:1064:GLY:O	1:A:1106:ASN:ND2	2.44	0.48
1:A:1428:ALA:O	1:A:1432:VAL:HG23	2.14	0.48
1:A:1658:SER:OG	5:A:2601:IHP:O32	2.29	0.48
1:B:615:HIS:CE1	1:B:617:GLU:HB2	2.49	0.48
1:B:825:MET:HE1	1:B:861:TYR:CG	2.49	0.48
3:E:36:CYS:SG	3:E:37:GLU:N	2.87	0.48
1:B:882:ARG:NH2	1:B:1569:ASP:OD2	2.47	0.47
3:E:447:LEU:HD23	3:E:447:LEU:H	1.79	0.47
3:F:36:CYS:SG	3:F:37:GLU:N	2.87	0.47
1:B:1428:ALA:O	1:B:1432:VAL:HG23	2.14	0.47
3:E:611:PRO:HB3	3:E:1154:ARG:HE	1.79	0.47
1:A:615:HIS:CE1	1:A:617:GLU:HB2	2.49	0.47
1:A:882:ARG:NH2	1:A:1569:ASP:OD2	2.47	0.47
1:B:2192:LEU:HD12	1:B:2235:GLY:HA3	1.95	0.47
1:B:2281:MET:HA	1:B:2284:VAL:HG12	1.96	0.47
3:E:23:ASP:OD2	3:E:999:ARG:NH2	2.45	0.47
3:E:318:ALA:HB1	3:E:435:ILE:HD13	1.96	0.47
3:F:1212:ALA:HB3	3:F:1231:SER:HB2	1.96	0.47
1:A:2133:CYS:O	1:A:2134:ARG:NE	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:65:ILE:HD11	2:C:88:ILE:HG21	1.95	0.47
3:E:49:TRP:CD2	3:E:468:LEU:HD12	2.48	0.47
1:A:825:MET:HE1	1:A:861:TYR:CG	2.49	0.47
1:A:1655:LYS:NZ	5:A:2601:IHP:O43	2.46	0.47
1:B:424:VAL:HG12	1:B:458:ILE:HG21	1.96	0.47
3:F:128:ASP:OD1	3:F:129:GLU:N	2.47	0.47
1:B:539:LEU:HD22	1:B:594:PHE:HE1	1.80	0.47
1:B:1505:ASN:O	1:B:1509:GLN:HB2	2.15	0.47
1:B:1979:ALA:HB1	1:B:1987:ARG:HD2	1.97	0.47
1:B:2133:CYS:O	1:B:2134:ARG:NE	2.48	0.47
1:B:2194:GLN:OE1	1:B:2430:ARG:NH1	2.48	0.47
3:E:166:VAL:HG23	3:E:175:ILE:HG13	1.96	0.47
3:E:318:ALA:HA	3:E:438:GLN:HG3	1.97	0.47
1:A:539:LEU:HD22	1:A:594:PHE:HE1	1.80	0.47
3:F:277:TRP:HA	3:F:280:MET:HE2	1.96	0.47
3:F:602:GLU:HA	3:F:605:TYR:CD2	2.50	0.47
1:A:1199:LEU:HD22	1:A:1204:ILE:HB	1.96	0.47
1:A:2281:MET:HA	1:A:2284:VAL:HG12	1.96	0.47
3:E:602:GLU:HA	3:E:605:TYR:CD2	2.50	0.47
3:E:1212:ALA:HB3	3:E:1231:SER:HB2	1.96	0.47
3:E:1283:SER:OG	3:E:1285:GLU:OE1	2.33	0.47
3:F:562:CYS:HB3	3:F:577:VAL:HG13	1.97	0.47
1:A:396:LEU:HB3	1:A:416:THR:HG23	1.97	0.47
1:A:705:PHE:O	1:A:709:GLU:HG3	2.15	0.47
1:A:964:LEU:HG	1:A:968:HIS:CE1	2.50	0.47
1:A:1505:ASN:O	1:A:1509:GLN:HB2	2.15	0.47
1:B:2546:CYS:HB3	1:B:2549:TRP:HB2	1.97	0.47
1:A:689:GLN:OE1	3:F:427:ARG:NH1	2.48	0.46
1:A:1150:ASP:OD2	1:B:732:LYS:NZ	2.33	0.46
1:B:145:VAL:HG13	1:B:175:LEU:HD13	1.97	0.46
1:B:783:LEU:HD13	1:B:823:ILE:HD12	1.97	0.46
3:F:447:LEU:H	3:F:447:LEU:HD23	1.79	0.46
1:A:665:THR:HG21	1:B:1191:PHE:HA	1.97	0.46
1:B:1626:CYS:SG	1:B:1627:GLN:N	2.88	0.46
3:E:191:ILE:HG22	3:E:241:ILE:HB	1.96	0.46
3:F:191:ILE:HG22	3:F:241:ILE:HB	1.96	0.46
1:A:1600:GLU:O	1:A:1603:ILE:HG12	2.16	0.46
1:A:2194:GLN:OE1	1:A:2430:ARG:NH1	2.48	0.46
3:E:562:CYS:HB3	3:E:577:VAL:HG13	1.97	0.46
1:A:1626:CYS:SG	1:A:1627:GLN:N	2.88	0.46
1:A:1979:ALA:HB1	1:A:1987:ARG:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2412:ASP:OD1	1:A:2413:SER:N	2.48	0.46
1:B:705:PHE:O	1:B:709:GLU:HG3	2.15	0.46
1:B:964:LEU:HG	1:B:968:HIS:CE1	2.50	0.46
3:F:318:ALA:HA	3:F:438:GLN:HG3	1.97	0.46
3:F:318:ALA:HB1	3:F:435:ILE:HD13	1.96	0.46
1:B:1981:LYS:NZ	1:B:2014:GLU:OE2	2.48	0.46
1:A:714:THR:HG22	1:A:717:ARG:HH21	1.81	0.46
1:A:2326:VAL:HG22	1:A:2403:VAL:HG21	1.98	0.46
3:E:544:ASN:OD1	3:E:583:ARG:NH1	2.49	0.46
1:A:1284:LEU:HD12	1:A:1353:VAL:HG22	1.98	0.46
3:F:1262:HIS:CD2	3:F:1263:PRO:HD2	2.51	0.46
1:B:1122:PHE:HB3	1:B:1135:ALA:HB3	1.98	0.46
3:E:419:TRP:CZ2	3:E:430:PRO:HD3	2.51	0.46
1:A:1024:LYS:HA	1:A:1062:ALA:HB1	1.98	0.46
1:A:1111:LEU:HD21	1:A:1148:PHE:CE2	2.51	0.46
1:A:2057:MET:HB3	1:A:2057:MET:HE3	1.72	0.46
1:B:1284:LEU:HD12	1:B:1353:VAL:HG22	1.98	0.46
3:E:833:VAL:O	3:E:836:SER:OG	2.31	0.46
3:F:544:ASN:OD1	3:F:583:ARG:NH1	2.49	0.46
1:A:145:VAL:HG13	1:A:175:LEU:HD13	1.97	0.46
1:A:1161:ARG:HH22	1:B:703:GLN:HE22	1.64	0.46
1:A:2265:HIS:O	1:A:2268:MET:HG2	2.16	0.46
1:A:2265:HIS:HA	1:A:2268:MET:SD	2.56	0.46
1:A:2387:MET:HE1	1:A:2392:LEU:HD23	1.98	0.46
1:B:955:LEU:HD22	1:B:971:VAL:HG23	1.98	0.46
1:B:2055:HIS:CE1	1:B:2082:GLN:HB2	2.51	0.46
1:B:2265:HIS:O	1:B:2268:MET:HG2	2.16	0.46
3:F:419:TRP:CZ2	3:F:430:PRO:HD3	2.51	0.46
1:A:2546:CYS:HB3	1:A:2549:TRP:HB2	1.98	0.45
1:B:714:THR:HG22	1:B:717:ARG:HH21	1.81	0.45
1:B:2015:GLU:OE1	1:B:2127:SER:OG	2.32	0.45
3:E:1028:VAL:HG13	3:E:1073:MET:HE2	1.98	0.45
3:F:1283:SER:OG	3:F:1285:GLU:OE1	2.33	0.45
1:A:466:LYS:HZ1	1:A:478:VAL:N	2.13	0.45
1:A:1332:ASP:OD1	1:A:1332:ASP:N	2.44	0.45
1:B:1111:LEU:HD21	1:B:1148:PHE:CE2	2.51	0.45
1:A:783:LEU:HD13	1:A:823:ILE:HD12	1.97	0.45
1:A:955:LEU:HD22	1:A:971:VAL:HG23	1.98	0.45
1:A:2418:LEU:HD13	1:A:2504:VAL:HG11	1.98	0.45
1:B:732:LYS:HD3	3:E:380:GLN:NE2	2.32	0.45
1:B:2412:ASP:OD1	1:B:2413:SER:N	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2418:LEU:HD23	1:B:2421:PHE:HE2	1.81	0.45
1:A:1191:PHE:HA	1:B:665:THR:HG21	1.98	0.45
1:B:1325:CYS:O	1:B:1329:LEU:HG	2.16	0.45
3:E:1220:GLN:OE1	3:E:1222:ARG:NH1	2.49	0.45
1:A:1325:CYS:O	1:A:1329:LEU:HG	2.16	0.45
1:A:2418:LEU:HD23	1:A:2421:PHE:HE2	1.81	0.45
3:E:468:LEU:HD21	3:E:498:ILE:HD12	1.98	0.45
3:F:23:ASP:OD2	3:F:999:ARG:NH2	2.45	0.45
3:F:442:SER:OG	3:F:443:GLN:N	2.50	0.45
1:A:661:VAL:O	1:A:665:THR:OG1	2.24	0.45
1:B:460:ARG:HH21	1:B:504:LEU:HD21	1.82	0.45
1:B:1024:LYS:HA	1:B:1062:ALA:HB1	1.98	0.45
1:B:2265:HIS:HA	1:B:2268:MET:SD	2.56	0.45
3:E:60:VAL:HG13	3:E:147:LEU:HD22	1.99	0.45
1:A:1122:PHE:HB3	1:A:1135:ALA:HB3	1.98	0.45
1:B:2387:MET:HE1	1:B:2392:LEU:HD23	1.98	0.45
3:E:79:THR:HB	3:E:80:PRO:HD3	1.99	0.45
1:A:2425:PRO:HA	1:A:2428:ASN:ND2	2.32	0.45
1:B:934:MET:HE2	1:B:934:MET:HB2	1.87	0.45
1:B:1060:VAL:HG11	1:B:1103:PHE:CE1	2.52	0.45
3:F:180:TYR:O	3:F:184:THR:OG1	2.26	0.45
3:F:1220:GLN:OE1	3:F:1222:ARG:NH1	2.49	0.45
1:A:460:ARG:HH21	1:A:504:LEU:HD21	1.82	0.45
1:A:947:TYR:HB2	1:A:948:PRO:HD3	1.99	0.45
1:A:1060:VAL:HG11	1:A:1103:PHE:CE1	2.52	0.45
1:B:396:LEU:HB3	1:B:416:THR:HG23	1.97	0.45
1:B:947:TYR:HB2	1:B:948:PRO:HD3	1.99	0.45
1:B:979:PHE:CD2	1:B:986:CYS:HB2	2.52	0.45
1:B:1600:GLU:O	1:B:1603:ILE:HG12	2.16	0.45
1:A:849:ALA:HB2	1:A:890:LEU:HD21	1.99	0.45
1:A:1530:GLU:HA	1:A:1550:LEU:HD11	1.98	0.45
1:A:2055:HIS:CE1	1:A:2082:GLN:HB2	2.52	0.45
1:B:2425:PRO:HA	1:B:2428:ASN:ND2	2.32	0.45
3:E:442:SER:OG	3:E:443:GLN:N	2.50	0.45
1:A:703:GLN:HE22	1:B:1161:ARG:HH22	1.65	0.44
1:A:724:ALA:HB3	3:F:415:ALA:HB2	1.99	0.44
1:A:2154:GLN:HB3	1:A:2175:MET:HG3	1.98	0.44
1:B:705:PHE:CE1	1:B:756:MET:HE2	2.52	0.44
1:B:1530:GLU:HA	1:B:1550:LEU:HD11	1.98	0.44
1:B:2326:VAL:HG22	1:B:2403:VAL:HG21	1.98	0.44
2:D:274:TRP:CZ2	2:D:316:VAL:HG12	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:468:LEU:HD21	3:F:498:ILE:HD12	1.98	0.44
2:C:287:THR:HB	2:C:297:TRP:HE1	1.83	0.44
1:A:732:LYS:HD3	3:F:380:GLN:NE2	2.32	0.44
1:A:824:ILE:O	1:A:828:LEU:HB2	2.18	0.44
1:B:404:PRO:HB3	1:B:444:VAL:HG22	2.00	0.44
1:B:693:LEU:HD12	1:B:693:LEU:HA	1.73	0.44
1:B:849:ALA:HB2	1:B:890:LEU:HD21	1.99	0.44
1:B:2418:LEU:HD13	1:B:2504:VAL:HG11	1.98	0.44
2:C:274:TRP:CZ2	2:C:316:VAL:HG12	2.52	0.44
3:F:1028:VAL:HG13	3:F:1073:MET:HE2	1.98	0.44
1:A:2011:MET:HE2	1:A:2011:MET:HB2	1.88	0.44
1:A:2339:ARG:HH21	1:A:2343:ASN:HB3	1.82	0.44
1:B:709:GLU:O	1:B:712:ILE:HG22	2.17	0.44
1:B:1278:ASP:OD1	1:B:1279:ASP:N	2.50	0.44
1:B:2285:GLU:HB2	2:D:272:TRP:CE3	2.53	0.44
3:E:1262:HIS:CD2	3:E:1263:PRO:HD2	2.51	0.44
3:F:293:ASP:OD1	3:F:294:LEU:N	2.51	0.44
3:F:1013:LEU:HD23	3:F:1013:LEU:H	1.82	0.44
1:A:404:PRO:HB3	1:A:444:VAL:HG22	2.00	0.44
1:A:979:PHE:CD2	1:A:986:CYS:HB2	2.52	0.44
1:A:1278:ASP:OD1	1:A:1279:ASP:N	2.50	0.44
3:E:369:ARG:HA	3:E:369:ARG:HH11	1.83	0.44
3:E:1013:LEU:HD23	3:E:1013:LEU:H	1.82	0.44
1:A:1432:VAL:HG22	1:A:2390:THR:HG21	2.00	0.44
1:B:724:ALA:HB3	3:E:415:ALA:HB2	1.98	0.44
1:B:2154:GLN:HB3	1:B:2175:MET:HG3	1.98	0.44
2:D:113:ASP:HB3	2:D:116:SER:OG	2.18	0.44
3:F:60:VAL:HG13	3:F:147:LEU:HD22	1.99	0.44
1:A:434:PHE:HB3	1:A:485:CYS:SG	2.58	0.44
1:A:1493:LEU:HD23	1:A:1519:ALA:HB2	2.00	0.44
1:A:1910:TRP:CD1	1:A:1923:LEU:HD13	2.53	0.44
1:A:2266:ARG:HB2	1:A:2266:ARG:NH1	2.33	0.44
1:B:689:GLN:OE1	3:E:427:ARG:NH1	2.50	0.44
2:D:287:THR:HB	2:D:297:TRP:HE1	1.83	0.44
3:F:980:LEU:HG	3:F:981:GLU:HG3	2.00	0.44
1:B:507:PRO:O	1:B:511:VAL:HG23	2.18	0.44
3:E:390:LEU:HA	3:E:393:LEU:HD13	2.00	0.44
3:F:126:THR:OG1	3:F:128:ASP:OD1	2.20	0.44
3:F:521:VAL:HG13	3:F:527:MET:SD	2.58	0.44
1:A:2421:PHE:HA	1:A:2424:ASP:HB2	2.00	0.44
1:B:434:PHE:HB3	1:B:485:CYS:SG	2.58	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:658:LYS:HA	1:B:661:VAL:HG22	2.00	0.44
3:F:79:THR:HB	3:F:80:PRO:HD3	1.99	0.44
3:F:390:LEU:HA	3:F:393:LEU:HD13	2.00	0.44
3:F:582:GLY:HA2	3:F:585:TRP:NE1	2.33	0.44
1:A:507:PRO:O	1:A:511:VAL:HG23	2.18	0.43
1:A:665:THR:HG23	1:B:1191:PHE:CD2	2.53	0.43
1:B:1338:ILE:HD11	1:B:1372:LEU:HB3	2.00	0.43
3:E:521:VAL:HG13	3:E:527:MET:SD	2.58	0.43
3:F:1074:GLU:OE1	3:F:1075:TYR:N	2.51	0.43
1:A:210:VAL:O	1:A:214:ARG:HG3	2.18	0.43
1:A:615:HIS:ND1	1:A:618:ILE:HG13	2.34	0.43
1:B:824:ILE:O	1:B:828:LEU:HB2	2.17	0.43
1:B:2002:SER:O	1:B:2003:ASN:HB3	2.18	0.43
1:B:2197:ARG:NH2	1:B:2424:ASP:OD2	2.51	0.43
1:A:709:GLU:O	1:A:712:ILE:HG22	2.17	0.43
1:A:865:LEU:HD21	1:A:1583:TYR:CE1	2.54	0.43
1:B:1432:VAL:HG22	1:B:2390:THR:HG21	2.00	0.43
1:B:1493:LEU:HD23	1:B:1519:ALA:HB2	1.99	0.43
1:B:1629:ILE:HD13	1:B:1629:ILE:HA	1.92	0.43
2:D:21:ASP:OD2	2:D:23:THR:OG1	2.33	0.43
3:E:293:ASP:OD1	3:E:294:LEU:N	2.51	0.43
3:E:563:LEU:HG	3:E:603:LYS:HD2	2.00	0.43
3:E:993:LEU:O	3:E:997:ARG:HG2	2.18	0.43
3:E:1089:ASP:N	3:E:1089:ASP:OD1	2.51	0.43
3:F:563:LEU:HG	3:F:603:LYS:HD2	2.00	0.43
1:A:2285:GLU:HB2	2:C:272:TRP:CE3	2.54	0.43
1:B:1564:ILE:HD11	1:B:1599:LEU:HD23	2.00	0.43
3:E:582:GLY:HA2	3:E:585:TRP:NE1	2.33	0.43
3:E:1074:GLU:OE1	3:E:1075:TYR:N	2.51	0.43
1:A:679:LEU:HD23	1:A:679:LEU:HA	1.86	0.43
1:A:705:PHE:CE1	1:A:756:MET:HE2	2.52	0.43
1:B:2266:ARG:NH1	1:B:2266:ARG:HB2	2.33	0.43
3:E:67:ASN:HB2	3:E:167:PHE:HZ	1.84	0.43
3:E:164:VAL:HG23	3:E:179:ILE:HD11	2.00	0.43
3:E:660:ARG:HH12	3:E:1112:TRP:HA	1.83	0.43
1:A:373:VAL:HA	1:A:392:ILE:HD13	1.99	0.43
1:A:1564:ILE:HD11	1:A:1599:LEU:HD23	2.00	0.43
1:B:373:VAL:HA	1:B:392:ILE:HD13	1.99	0.43
2:C:113:ASP:HB3	2:C:116:SER:OG	2.18	0.43
3:F:993:LEU:O	3:F:997:ARG:HG2	2.18	0.43
1:A:658:LYS:HA	1:A:661:VAL:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2432:MET:HE1	1:A:2493:LEU:HA	2.00	0.43
1:B:210:VAL:O	1:B:214:ARG:HG3	2.18	0.43
1:B:827:MET:HE2	1:B:827:MET:HB2	1.80	0.43
1:B:865:LEU:HD21	1:B:1583:TYR:CE1	2.54	0.43
1:B:1326:TRP:HZ3	1:B:1369:LYS:HB3	1.84	0.43
1:B:2432:MET:HE1	1:B:2493:LEU:HA	2.00	0.43
3:F:1089:ASP:N	3:F:1089:ASP:OD1	2.51	0.43
1:A:484:THR:O	1:A:487:SER:OG	2.28	0.43
1:A:2034:ALA:HB1	1:A:2047:MET:SD	2.59	0.43
1:B:772:MET:HE3	1:B:772:MET:HB3	1.91	0.43
1:B:855:VAL:HG21	1:B:1625:GLY:HA3	2.01	0.43
1:B:1910:TRP:CD1	1:B:1923:LEU:HD13	2.53	0.43
1:B:2052:GLU:HG3	1:B:2053:PRO:HD3	2.01	0.43
2:D:17:THR:HB	2:D:27:TRP:HE1	1.84	0.43
3:E:1230:VAL:HG22	3:E:1236:VAL:HG22	2.01	0.43
3:F:1040:VAL:HG21	3:F:1325:ILE:HD13	2.01	0.43
1:A:485:CYS:HA	1:A:488:MET:HE2	2.01	0.43
1:A:1297:SER:OG	1:A:1300:LEU:HB2	2.19	0.43
1:A:2278:LEU:HD13	1:A:2282:GLN:HB3	2.01	0.43
1:B:1071:LEU:N	1:B:1072:PRO:HD2	2.34	0.43
3:F:1064:ASN:H	3:F:1064:ASN:HD22	1.67	0.43
1:A:1060:VAL:HG11	1:A:1103:PHE:CD1	2.54	0.43
1:A:1372:LEU:HD12	1:A:1372:LEU:H	1.83	0.43
1:A:2002:SER:O	1:A:2003:ASN:HB3	2.18	0.43
1:B:485:CYS:HA	1:B:488:MET:HE2	2.01	0.43
1:B:660:LEU:HD23	1:B:660:LEU:HA	1.85	0.43
1:B:1372:LEU:H	1:B:1372:LEU:HD12	1.83	0.43
1:B:1628:ARG:HH22	1:B:1655:LYS:HE3	1.83	0.43
1:B:2011:MET:SD	1:B:2012:VAL:HG23	2.59	0.43
2:C:17:THR:HB	2:C:27:TRP:HE1	1.84	0.43
2:D:242:GLN:HB2	2:D:261:LYS:HZ1	1.83	0.43
3:F:479:LEU:HD13	3:F:479:LEU:HA	1.86	0.43
3:F:479:LEU:HD22	4:H:403:MET:HE1	2.00	0.43
3:F:601:HIS:CD2	3:F:640:ILE:HD11	2.54	0.43
3:F:1230:VAL:HG22	3:F:1236:VAL:HG22	2.00	0.43
1:A:855:VAL:HG21	1:A:1625:GLY:HA3	2.01	0.42
1:A:888:LEU:HD23	1:A:888:LEU:HA	1.87	0.42
1:A:1326:TRP:HZ3	1:A:1369:LYS:HB3	1.84	0.42
1:A:1460:LEU:HD11	1:A:1484:LEU:HD22	2.01	0.42
1:A:1650:MET:HE3	1:A:1650:MET:HB3	1.90	0.42
1:A:2197:ARG:NH2	1:A:2424:ASP:OD2	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1297:SER:OG	1:B:1300:LEU:HB2	2.19	0.42
3:F:369:ARG:HA	3:F:369:ARG:HH11	1.83	0.42
3:F:467:ALA:O	3:F:470:VAL:HG12	2.19	0.42
1:A:1071:LEU:N	1:A:1072:PRO:HD2	2.34	0.42
3:E:980:LEU:HG	3:E:981:GLU:HG3	2.00	0.42
3:E:1048:PHE:HE2	3:E:1060:PHE:CE2	2.37	0.42
1:A:521:LEU:HD22	1:A:539:LEU:HD12	2.01	0.42
1:A:1338:ILE:HD11	1:A:1372:LEU:HB3	2.00	0.42
1:A:2011:MET:SD	1:A:2012:VAL:HG23	2.59	0.42
1:B:1578:MET:HE2	1:B:1578:MET:HA	2.01	0.42
1:B:1877:LEU:O	1:B:1881:THR:OG1	2.29	0.42
3:E:314:TRP:HZ2	3:E:439:VAL:HA	1.84	0.42
3:E:337:PHE:CZ	3:E:347:PHE:HB3	2.54	0.42
3:E:467:ALA:O	3:E:470:VAL:HG12	2.19	0.42
3:F:64:LEU:HD11	3:F:106:LEU:HD23	2.01	0.42
3:F:164:VAL:HG23	3:F:179:ILE:HD11	2.00	0.42
3:F:1177:HIS:O	3:F:1177:HIS:CG	2.72	0.42
1:A:774:PRO:HB3	3:F:286:LEU:HA	2.02	0.42
1:A:1065:GLY:N	3:E:361:ASN:OD1	2.52	0.42
1:A:1191:PHE:CD2	1:B:665:THR:HG23	2.54	0.42
1:B:851:THR:HG23	1:B:853:TYR:HB2	2.01	0.42
1:B:1109:ASP:OD1	3:F:282:LYS:NZ	2.53	0.42
1:B:1879:MET:HE3	1:B:1879:MET:HB3	1.90	0.42
3:E:1040:VAL:HG21	3:E:1325:ILE:HD13	2.01	0.42
3:F:337:PHE:CZ	3:F:347:PHE:HB3	2.54	0.42
1:B:2421:PHE:HA	1:B:2424:ASP:HB2	2.00	0.42
2:C:21:ASP:OD2	2:C:23:THR:OG1	2.33	0.42
2:C:291:ASP:O	2:C:293:LEU:HG	2.20	0.42
2:D:129:ALA:HB3	2:D:147:ASP:HB2	2.02	0.42
3:E:358:ARG:HD3	3:E:374:TYR:CG	2.55	0.42
3:F:67:ASN:HB2	3:F:167:PHE:HZ	1.84	0.42
1:A:1628:ARG:HH22	1:A:1655:LYS:HE3	1.83	0.42
1:A:1910:TRP:NE1	1:A:1953:VAL:HG13	2.35	0.42
1:B:615:HIS:ND1	1:B:618:ILE:HG13	2.34	0.42
1:B:1349:ASP:OD1	1:B:1354:THR:OG1	2.38	0.42
1:B:1380:ILE:HD13	1:B:1380:ILE:HA	1.91	0.42
1:B:2204:LEU:HD22	1:B:2417:VAL:HG21	2.02	0.42
1:B:2339:ARG:HH21	1:B:2343:ASN:HB3	1.83	0.42
3:E:479:LEU:HD22	4:G:403:MET:HE1	2.00	0.42
1:A:2526:GLU:O	1:A:2530:LYS:HG2	2.20	0.42
1:B:664:ILE:HD11	1:B:699:ALA:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2208:LEU:HD22	1:B:2410:HIS:CG	2.55	0.42
2:C:129:ALA:HB3	2:C:147:ASP:HB2	2.02	0.42
3:E:258:LEU:HD23	3:E:258:LEU:HA	1.88	0.42
3:E:601:HIS:CD2	3:E:640:ILE:HD11	2.54	0.42
3:F:660:ARG:HH12	3:F:1112:TRP:HA	1.83	0.42
1:A:1578:MET:HA	1:A:1578:MET:HE2	2.01	0.42
1:B:689:GLN:HG2	3:E:422:MET:HE3	2.01	0.42
1:B:1060:VAL:HG11	1:B:1103:PHE:CD1	2.54	0.42
1:B:1460:LEU:HD11	1:B:1484:LEU:HD22	2.01	0.42
1:B:2526:GLU:O	1:B:2530:LYS:HG2	2.20	0.42
2:D:14:ILE:HD13	2:D:70:LEU:HD22	2.02	0.42
3:E:299:PRO:HD3	3:E:397:ILE:HD11	2.02	0.42
3:E:1064:ASN:HD22	3:E:1064:ASN:H	1.67	0.42
3:F:413:LEU:HD13	3:F:413:LEU:HA	1.90	0.42
3:F:658:MET:HE3	3:F:658:MET:HB3	1.86	0.42
3:F:815:LEU:HB3	3:F:834:LEU:HD21	2.02	0.42
1:A:934:MET:HE2	1:A:934:MET:HB2	1.87	0.42
1:B:83:ARG:O	1:B:87:ILE:HG13	2.20	0.42
1:B:1065:GLY:N	3:F:361:ASN:OD1	2.52	0.42
1:B:1721:VAL:HA	1:B:1747:MET:HE1	2.02	0.42
2:C:46:ASN:ND2	2:C:61:GLY:O	2.52	0.42
3:E:64:LEU:HD11	3:E:106:LEU:HD23	2.01	0.42
3:F:314:TRP:HZ2	3:F:439:VAL:HA	1.84	0.42
1:A:689:GLN:HG2	3:F:422:MET:HE3	2.02	0.42
1:A:1490:TRP:CE3	1:A:1519:ALA:HA	2.55	0.42
1:A:2052:GLU:HG3	1:A:2053:PRO:HD3	2.01	0.42
1:A:2366:MET:HG2	1:A:2373:GLU:O	2.20	0.42
1:B:1329:LEU:HD13	1:B:1333:GLN:HB2	2.02	0.42
1:B:1910:TRP:NE1	1:B:1953:VAL:HG13	2.35	0.42
1:B:2034:ALA:HB1	1:B:2047:MET:SD	2.59	0.42
1:B:2366:MET:HG2	1:B:2373:GLU:O	2.20	0.42
1:B:2516:ASP:OD1	1:B:2516:ASP:N	2.53	0.42
3:E:309:LEU:HD12	3:E:309:LEU:HA	1.86	0.42
3:E:395:THR:O	3:E:400:GLY:N	2.40	0.42
3:E:1189:SER:HB3	3:E:1208:ARG:HG2	2.02	0.42
1:A:827:MET:HE3	1:A:835:ALA:HB3	2.00	0.41
1:B:1108:ASP:OD1	1:B:1109:ASP:N	2.53	0.41
1:B:1407:GLY:HA2	1:B:1408:PRO:HD3	1.97	0.41
1:B:1481:MET:HE1	1:B:1512:MET:HE1	2.02	0.41
1:B:1680:ASP:HB2	1:B:1683:ARG:HG2	2.02	0.41
1:B:2047:MET:HE3	1:B:2047:MET:HB3	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2278:LEU:HD13	1:B:2282:GLN:HB3	2.01	0.41
3:E:346:LEU:HD23	3:E:346:LEU:HA	1.90	0.41
3:F:57:THR:HG21	3:F:342:LEU:HD23	2.02	0.41
3:F:105:ASN:O	3:F:109:GLN:HG2	2.20	0.41
3:F:299:PRO:HD3	3:F:397:ILE:HD11	2.02	0.41
1:B:774:PRO:HB3	3:E:286:LEU:HA	2.02	0.41
1:B:1195:VAL:O	1:B:1199:LEU:HG	2.20	0.41
1:B:1322:PHE:CD1	1:B:1360:LEU:HD21	2.55	0.41
3:E:57:THR:HG21	3:E:342:LEU:HD23	2.02	0.41
3:E:815:LEU:HB3	3:E:834:LEU:HD21	2.02	0.41
1:A:195:PHE:CD2	1:A:240:GLU:HG3	2.55	0.41
1:A:851:THR:HG23	1:A:853:TYR:HB2	2.01	0.41
1:A:1195:VAL:O	1:A:1199:LEU:HG	2.20	0.41
1:A:1680:ASP:HB2	1:A:1683:ARG:HG2	2.02	0.41
2:C:14:ILE:HD13	2:C:70:LEU:HD22	2.02	0.41
2:D:291:ASP:O	2:D:293:LEU:HG	2.20	0.41
3:E:319:ILE:HG23	3:E:378:MET:HG3	2.03	0.41
3:E:1320:SER:OG	3:E:1321:ASN:N	2.53	0.41
3:F:1048:PHE:HE2	3:F:1060:PHE:CE2	2.37	0.41
3:F:1218:SER:HB3	3:F:1261:ILE:HD12	2.02	0.41
1:A:1129:LEU:N	1:A:1130:PRO:HD2	2.35	0.41
1:A:1413:LEU:HD13	1:A:1436:ALA:HB2	2.03	0.41
1:A:1481:MET:HE1	1:A:1512:MET:HE1	2.02	0.41
1:A:2208:LEU:HD22	1:A:2410:HIS:CG	2.55	0.41
1:B:195:PHE:CD2	1:B:240:GLU:HG3	2.55	0.41
2:D:46:ASN:ND2	2:D:61:GLY:O	2.52	0.41
3:E:271:ILE:HD11	3:E:313:ASN:HA	2.02	0.41
3:E:1177:HIS:O	3:E:1177:HIS:CG	2.72	0.41
3:F:257:ASP:OD1	3:F:257:ASP:N	2.52	0.41
1:A:820:LEU:HB3	1:A:847:LEU:HD11	2.03	0.41
1:A:1423:LEU:HD12	1:A:1423:LEU:HA	1.87	0.41
1:B:521:LEU:HD22	1:B:539:LEU:HD12	2.01	0.41
3:F:358:ARG:HD3	3:F:374:TYR:CG	2.55	0.41
3:F:533:THR:HG21	3:F:572:LEU:HG	2.02	0.41
3:F:1135:THR:HG23	3:F:1137:LEU:H	1.85	0.41
1:A:1329:LEU:HD13	1:A:1333:GLN:HB2	2.02	0.41
1:A:1349:ASP:OD1	1:A:1354:THR:OG1	2.38	0.41
1:B:1129:LEU:N	1:B:1130:PRO:HD2	2.35	0.41
3:E:1135:THR:HG23	3:E:1137:LEU:H	1.86	0.41
1:A:203:GLN:O	1:A:207:GLU:HG3	2.21	0.41
1:A:2047:MET:HE3	1:A:2047:MET:HB3	1.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:827:MET:HE3	1:B:835:ALA:HB3	2.00	0.41
1:B:1096:LEU:O	1:B:1100:ILE:HG12	2.21	0.41
1:B:1362:GLU:OE2	1:B:2309:SER:HA	2.21	0.41
3:E:257:ASP:OD1	3:E:257:ASP:N	2.52	0.41
3:E:533:THR:HG21	3:E:572:LEU:HG	2.02	0.41
3:E:1147:VAL:HB	3:E:1161:ILE:HD11	2.02	0.41
3:F:271:ILE:HD11	3:F:313:ASN:HA	2.02	0.41
1:A:83:ARG:O	1:A:87:ILE:HG13	2.20	0.41
1:A:795:ILE:HA	1:A:798:VAL:HG12	2.03	0.41
1:A:1322:PHE:CD1	1:A:1360:LEU:HD21	2.55	0.41
1:A:1386:ARG:HA	1:A:1386:ARG:HD3	1.87	0.41
1:A:1484:LEU:HD13	1:A:1484:LEU:HA	1.95	0.41
1:A:2204:LEU:HD22	1:A:2417:VAL:HG21	2.02	0.41
3:E:105:ASN:O	3:E:109:GLN:HG2	2.20	0.41
3:E:519:LEU:HD22	3:E:561:ILE:HD12	2.03	0.41
3:E:1218:SER:HB3	3:E:1261:ILE:HD12	2.02	0.41
3:F:1189:SER:HB3	3:F:1208:ARG:HG2	2.02	0.41
3:F:1236:VAL:HB	3:F:1250:LEU:HB2	2.02	0.41
1:A:199:TRP:HE3	1:A:260:ARG:HG2	1.86	0.41
1:A:664:ILE:HD11	1:A:699:ALA:HB2	2.01	0.41
1:A:693:LEU:HA	1:A:693:LEU:HD12	1.73	0.41
1:A:1096:LEU:O	1:A:1100:ILE:HG12	2.21	0.41
1:A:1671:LYS:O	1:A:1674:VAL:HG12	2.21	0.41
1:B:199:TRP:HE3	1:B:260:ARG:HG2	1.86	0.41
1:B:203:GLN:O	1:B:207:GLU:HG3	2.21	0.41
1:B:675:VAL:O	1:B:678:SER:OG	2.31	0.41
1:B:1027:ILE:HD12	1:B:1027:ILE:HA	1.98	0.41
1:B:1627:GLN:O	1:B:1633:TRP:NE1	2.47	0.41
1:B:1753:LYS:HD3	1:B:1753:LYS:HA	1.96	0.41
3:E:1083:LEU:HD13	3:E:1095:VAL:HG11	2.03	0.41
3:F:156:PRO:HG2	3:F:165:TRP:CG	2.56	0.41
3:F:188:SER:OG	3:F:189:PRO:HD3	2.21	0.41
3:F:304:ASP:O	3:F:307:THR:OG1	2.33	0.41
3:F:1320:SER:OG	3:F:1321:ASN:N	2.53	0.41
1:A:1721:VAL:HA	1:A:1747:MET:HE1	2.02	0.41
1:A:1974:TYR:O	1:A:1978:VAL:HG23	2.20	0.41
1:A:1981:LYS:NZ	1:A:2014:GLU:OE2	2.48	0.41
1:A:2516:ASP:OD1	1:A:2516:ASP:N	2.53	0.41
2:C:68:TYR:HE1	2:C:77:PRO:HG3	1.86	0.41
2:C:136:LEU:HD21	2:C:140:GLN:OE1	2.21	0.41
3:F:309:LEU:HD12	3:F:309:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:420:VAL:O	1:B:424:VAL:HG23	2.21	0.40
1:B:1671:LYS:O	1:B:1674:VAL:HG12	2.21	0.40
1:B:2006:VAL:O	1:B:2010:MET:HG2	2.21	0.40
2:D:136:LEU:HD21	2:D:140:GLN:OE1	2.21	0.40
3:E:188:SER:OG	3:E:189:PRO:HD3	2.21	0.40
3:F:570:HIS:HB3	3:F:573:LEU:HB3	2.03	0.40
3:F:984:ILE:H	3:F:984:ILE:HD12	1.86	0.40
3:F:1147:VAL:HB	3:F:1161:ILE:HD11	2.02	0.40
1:A:420:VAL:O	1:A:424:VAL:HG23	2.21	0.40
1:A:856:GLU:H	1:A:856:GLU:HG3	1.73	0.40
1:A:1012:LEU:HD23	1:A:1012:LEU:HA	1.86	0.40
1:A:1759:LEU:HD11	1:A:1796:MET:SD	2.62	0.40
1:A:2122:GLU:OE1	1:A:2159:SER:OG	2.34	0.40
1:B:1208:ARG:HA	1:B:1211:VAL:HG12	2.04	0.40
1:B:1974:TYR:O	1:B:1978:VAL:HG23	2.20	0.40
3:F:298:ILE:HG13	3:F:393:LEU:HD21	2.04	0.40
3:F:319:ILE:HG23	3:F:378:MET:HG3	2.03	0.40
1:A:1787:TYR:CE1	1:A:1895:SER:HB2	2.57	0.40
1:A:2194:GLN:HE22	1:A:2421:PHE:HE1	1.69	0.40
1:A:2280:LEU:HD12	1:A:2280:LEU:HA	1.95	0.40
1:B:825:MET:HB2	1:B:825:MET:HE3	1.77	0.40
1:B:1197:LYS:HE3	1:B:1197:LYS:HB3	1.96	0.40
1:B:1490:TRP:CE3	1:B:1519:ALA:HA	2.55	0.40
2:C:224:LEU:HG	2:C:274:TRP:HB3	2.03	0.40
3:E:527:MET:HG3	3:E:528:PRO:HD2	2.03	0.40
3:F:605:TYR:HB3	3:F:647:MET:HE3	2.03	0.40
1:A:1108:ASP:OD1	1:A:1109:ASP:N	2.53	0.40
1:A:2006:VAL:O	1:A:2010:MET:HG2	2.21	0.40
1:B:1888:PHE:O	1:B:1892:ILE:HG13	2.21	0.40
1:B:2002:SER:C	1:B:2004:THR:H	2.30	0.40
2:C:139:ASN:ND2	2:C:141:ALA:HB3	2.37	0.40
3:E:570:HIS:HB3	3:E:573:LEU:HB3	2.03	0.40
3:F:169:LYS:HD2	3:F:169:LYS:HA	1.94	0.40
3:F:1083:LEU:HD13	3:F:1095:VAL:HG11	2.03	0.40
3:F:1128:VAL:HG11	3:F:1172:LEU:N	2.37	0.40
3:F:1139:MET:HE3	3:F:1149:ILE:HG23	2.03	0.40
1:A:449:LYS:HA	1:A:452:LEU:HG	2.03	0.40
1:A:774:PRO:CB	3:F:286:LEU:HA	2.51	0.40
1:A:2002:SER:C	1:A:2004:THR:H	2.30	0.40
1:A:2254:ARG:NH1	1:A:2261:LEU:O	2.54	0.40
1:B:390:MET:HE2	1:B:390:MET:HB3	2.00	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:449:LYS:HA	1:B:452:LEU:HG	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	A	2150/2549 (84%)	2101 (98%)	48 (2%)	1 (0%)	100	100
1	B	2150/2549 (84%)	2101 (98%)	48 (2%)	1 (0%)	100	100
2	C	315/326 (97%)	302 (96%)	13 (4%)	0	100	100
2	D	315/326 (97%)	302 (96%)	13 (4%)	0	100	100
3	E	1040/1363 (76%)	1007 (97%)	33 (3%)	0	100	100
3	F	1040/1363 (76%)	1007 (97%)	33 (3%)	0	100	100
4	G	7/13 (54%)	7 (100%)	0	0	100	100
4	H	7/13 (54%)	7 (100%)	0	0	100	100
All	All	7024/8502 (83%)	6834 (97%)	188 (3%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	379	LYS
1	B	379	LYS

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	1876/2220 (84%)	1844 (98%)	32 (2%)	53	67
1	B	1876/2220 (84%)	1844 (98%)	32 (2%)	53	67
2	C	269/276 (98%)	265 (98%)	4 (2%)	57	68
2	D	269/276 (98%)	265 (98%)	4 (2%)	57	68
3	E	928/1191 (78%)	913 (98%)	15 (2%)	55	67
3	F	928/1191 (78%)	913 (98%)	15 (2%)	55	67
4	G	9/13 (69%)	9 (100%)	0	100	100
4	H	9/13 (69%)	9 (100%)	0	100	100
All	All	6164/7400 (83%)	6062 (98%)	102 (2%)	52	67

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	548	MET
1	A	665	THR
1	A	671	ILE
1	A	693	LEU
1	A	794	VAL
1	A	820	LEU
1	A	825	MET
1	A	826	ASP
1	A	827	MET
1	A	968	HIS
1	A	1001	ILE
1	A	1097	LEU
1	A	1119	VAL
1	A	1160	VAL
1	A	1220	TYR
1	A	1262	THR
1	A	1284	LEU
1	A	1311	ASN
1	A	1372	LEU
1	A	1477	MET
1	A	1484	LEU
1	A	1624	GLN
1	A	1673	LEU
1	A	1743	LEU

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Mol	Chain	Res	Type
1	A	1808	ASN
1	A	1946	ILE
1	A	2057	MET
1	A	2138	LEU
1	A	2172	LEU
1	A	2173	THR
1	A	2303	LEU
1	A	2540	GLN
1	B	548	MET
1	B	665	THR
1	B	671	ILE
1	B	693	LEU
1	B	794	VAL
1	B	820	LEU
1	B	825	MET
1	B	826	ASP
1	B	827	MET
1	B	968	HIS
1	B	1001	ILE
1	B	1097	LEU
1	B	1119	VAL
1	B	1160	VAL
1	B	1220	TYR
1	B	1262	THR
1	B	1284	LEU
1	B	1311	ASN
1	B	1372	LEU
1	B	1477	MET
1	B	1484	LEU
1	B	1624	GLN
1	B	1673	LEU
1	B	1743	LEU
1	B	1808	ASN
1	B	1946	ILE
1	B	2057	MET
1	B	2138	LEU
1	B	2172	LEU
1	B	2173	THR
1	B	2303	LEU
1	B	2540	GLN
2	C	128	ASN
2	C	233	THR

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Mol	Chain	Res	Type
2	C	274	TRP
2	C	292	ASN
2	D	128	ASN
2	D	233	THR
2	D	274	TRP
2	D	292	ASN
3	E	78	THR
3	E	186	MET
3	E	211	LEU
3	E	257	ASP
3	E	336	LEU
3	E	479	LEU
3	E	527	MET
3	E	551	GLU
3	E	646	MET
3	E	1011	THR
3	E	1014	ASP
3	E	1064	ASN
3	E	1139	MET
3	E	1323	TYR
3	E	1330	VAL
3	F	78	THR
3	F	186	MET
3	F	211	LEU
3	F	257	ASP
3	F	336	LEU
3	F	479	LEU
3	F	527	MET
3	F	551	GLU
3	F	646	MET
3	F	1011	THR
3	F	1014	ASP
3	F	1064	ASN
3	F	1139	MET
3	F	1323	TYR
3	F	1330	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	A	418	ASN

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Mol	Chain	Res	Type
1	A	536	GLN
1	A	689	GLN
1	A	703	GLN
1	A	1058	GLN
1	A	1308	GLN
1	A	1446	GLN
1	A	1992	ASN
1	A	2093	ASN
1	A	2194	GLN
1	A	2428	ASN
1	B	70	HIS
1	B	419	HIS
1	B	536	GLN
1	B	689	GLN
1	B	703	GLN
1	B	1308	GLN
1	B	1446	GLN
1	B	1920	ASN
1	B	1992	ASN
1	B	2082	GLN
1	B	2093	ASN
1	B	2194	GLN
1	B	2428	ASN
1	B	2537	ASN
2	C	39	GLN
2	C	148	GLN
2	D	39	GLN
2	D	148	GLN
3	E	281	GLN
3	E	349	ASN
3	E	380	GLN
3	E	438	GLN
3	E	586	GLN
3	E	601	HIS
3	E	643	ASN
3	E	677	ASN
3	E	1064	ASN
3	E	1233	ASN
3	E	1289	ASN
3	E	1310	HIS
3	F	151	ASN
3	F	281	GLN

Continued on next page...

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Mol	Chain	Res	Type
3	F	380	GLN
3	F	586	GLN
3	F	601	HIS
3	F	643	ASN
3	F	669	HIS
3	F	677	ASN
3	F	1233	ASN
3	F	1289	ASN
3	F	1310	HIS

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](#)

2 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	$\# Z > 2$
5	IHP	B	2601	-	36,36,36	0.78	2 (5%)	60,60,60	0.57	1 (1%)
5	IHP	A	2601	-	36,36,36	0.79	2 (5%)	60,60,60	0.57	1 (1%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	IHP	B	2601	-	-	1/30/54/54	0/1/1/1
5	IHP	A	2601	-	-	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)
5	A	2601	IHP	P3-O13	3.08	1.64	1.59
5	B	2601	IHP	P3-O13	3.02	1.64	1.59
5	A	2601	IHP	P1-O11	2.07	1.63	1.59
5	B	2601	IHP	P1-O11	2.03	1.63	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	2601	IHP	P2-O12-C2	2.88	131.11	123.43
5	A	2601	IHP	P2-O12-C2	2.87	131.11	123.43

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	2601	IHP	C2-O12-P2-O22
5	B	2601	IHP	C2-O12-P2-O22

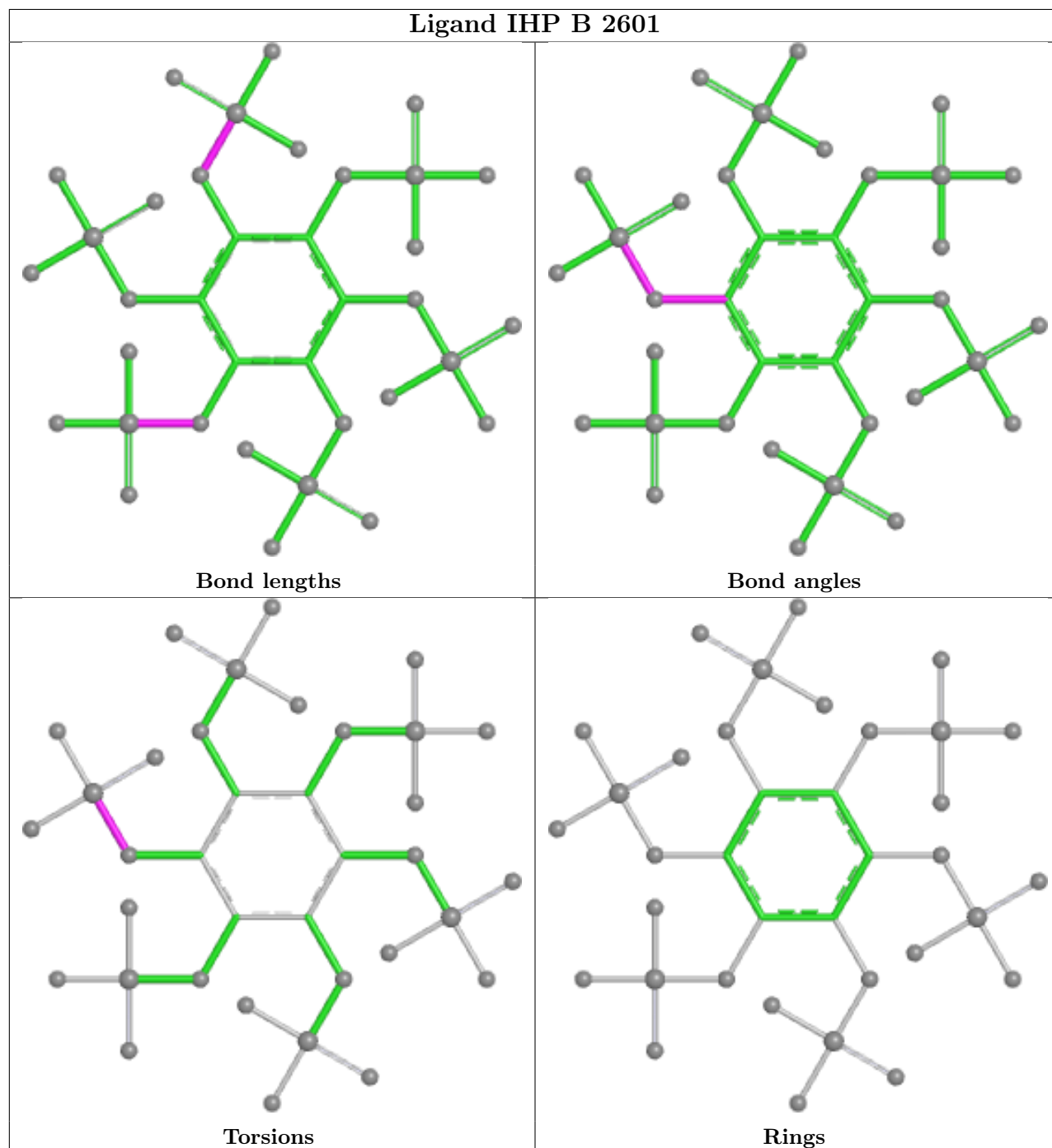
There are no ring outliers.

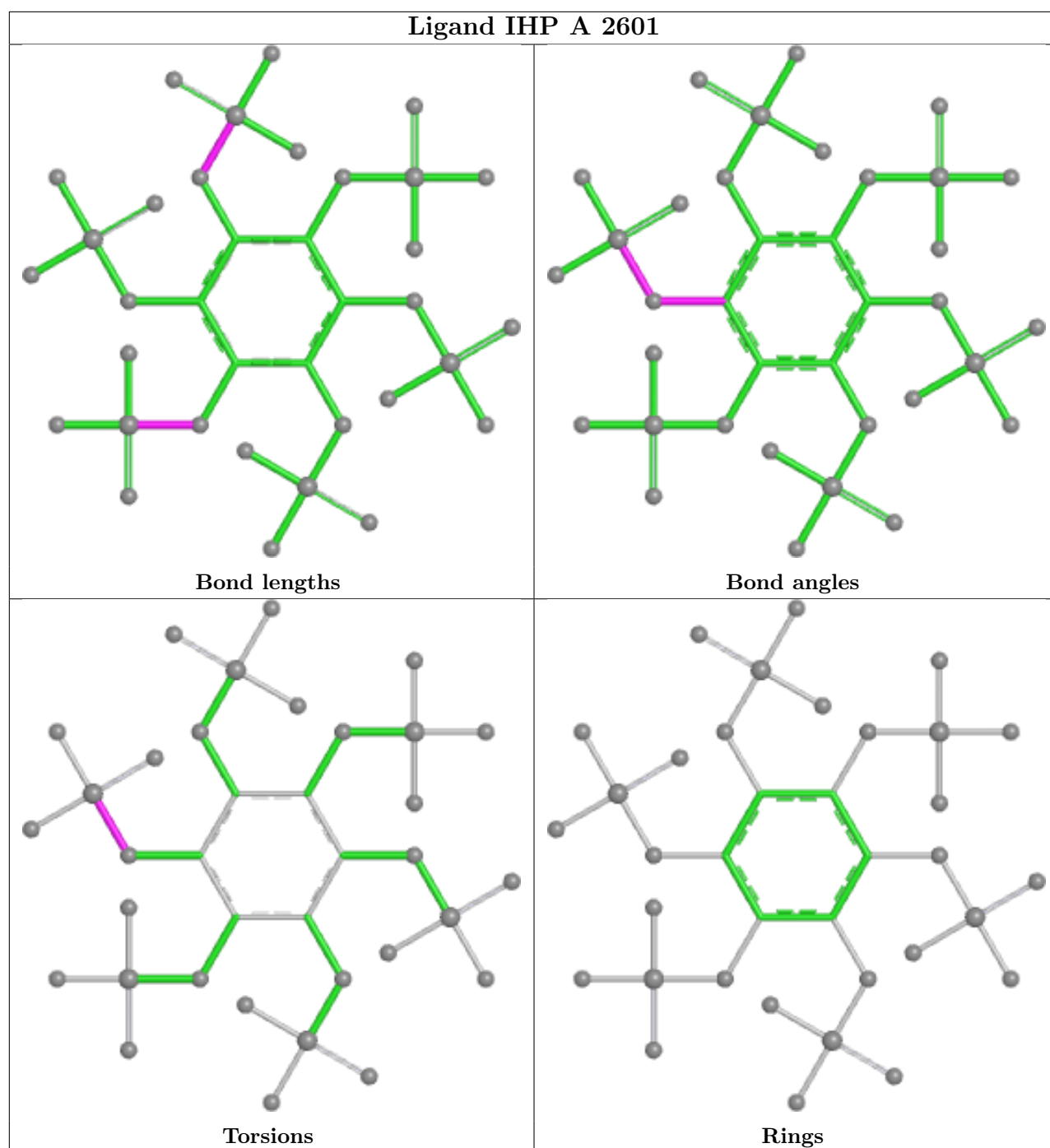
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2601	IHP	2	0
5	A	2601	IHP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

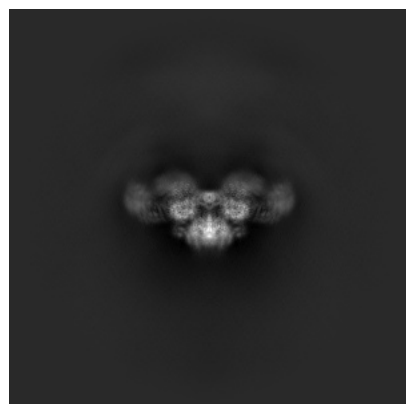
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-50183. These allow visual inspection of the internal detail of the map and identification of artifacts.

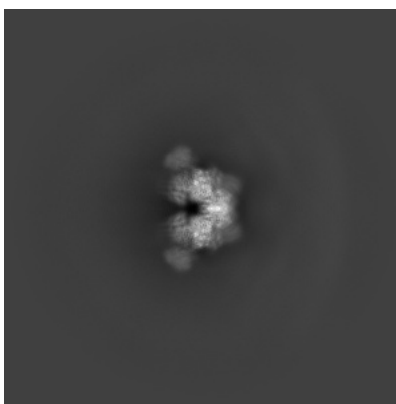
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

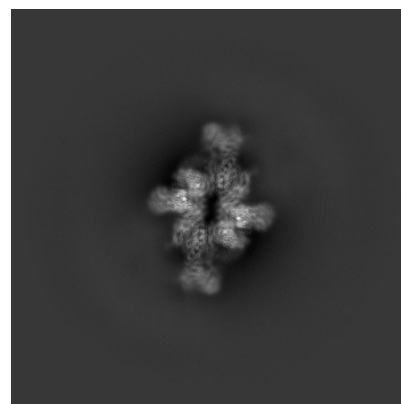
6.1.1 Primary map



X

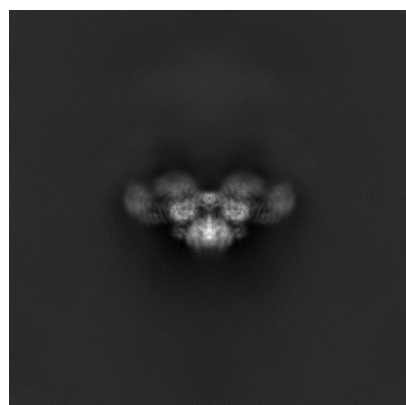


Y

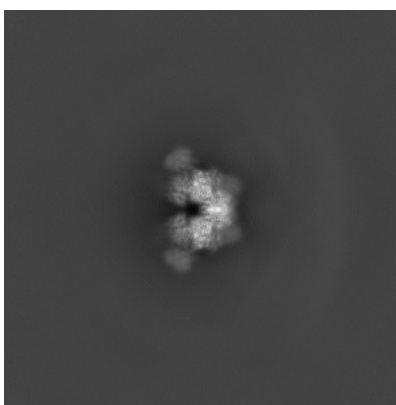


Z

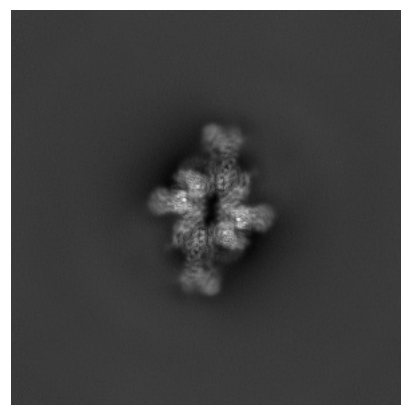
6.1.2 Raw map



X



Y

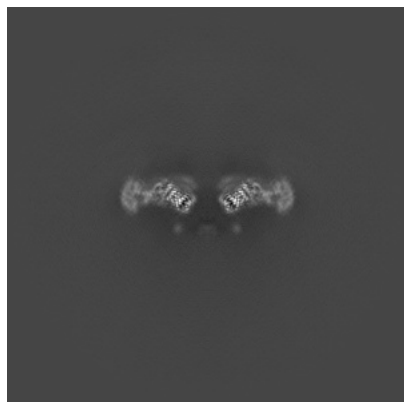


Z

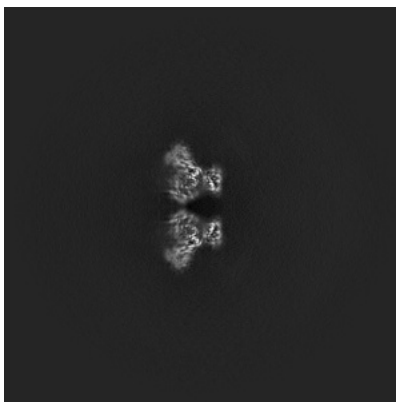
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

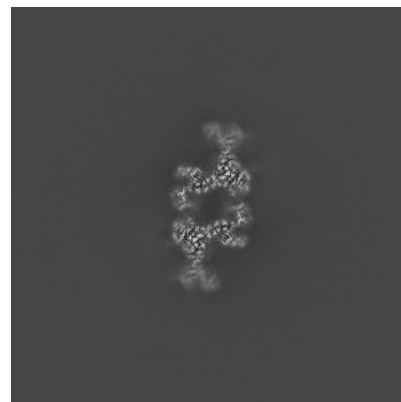
6.2.1 Primary map



X Index: 256

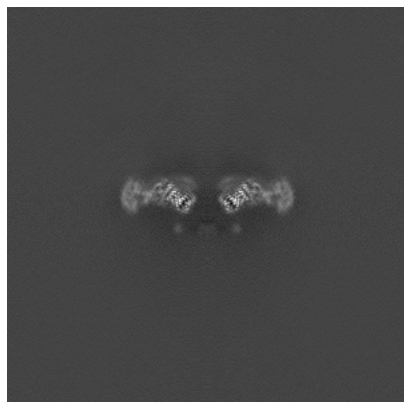


Y Index: 256

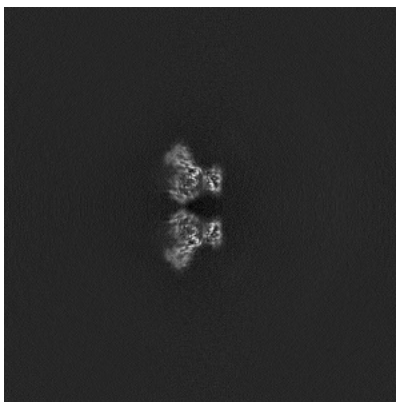


Z Index: 256

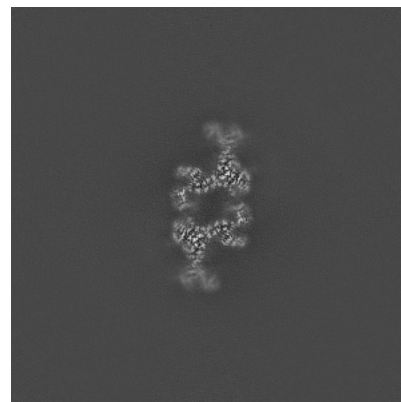
6.2.2 Raw map



X Index: 256



Y Index: 256

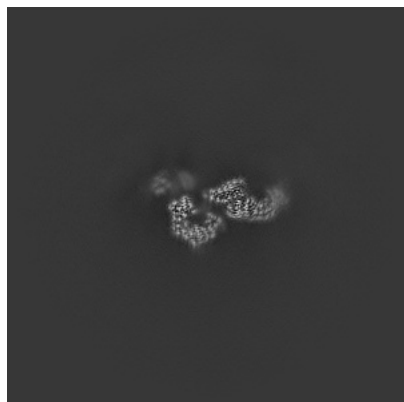


Z Index: 256

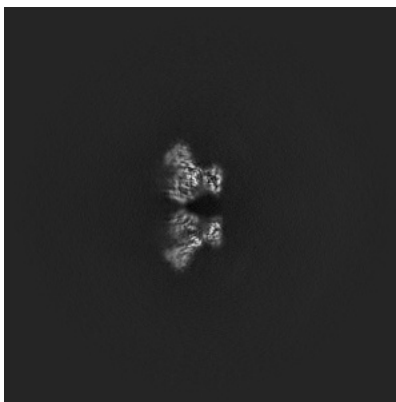
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

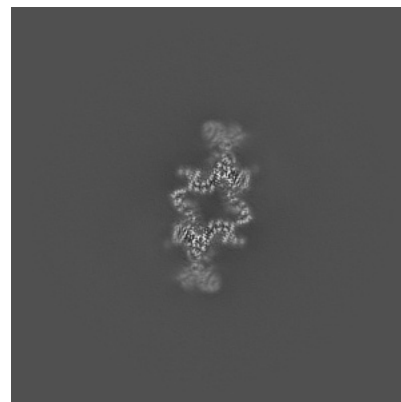
6.3.1 Primary map



X Index: 274

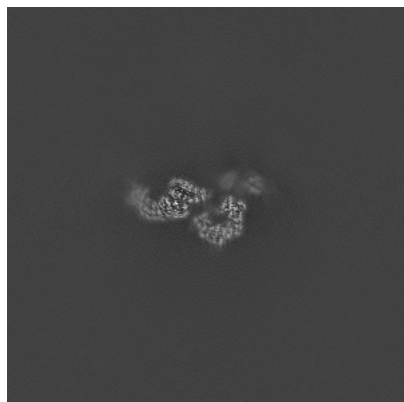


Y Index: 255

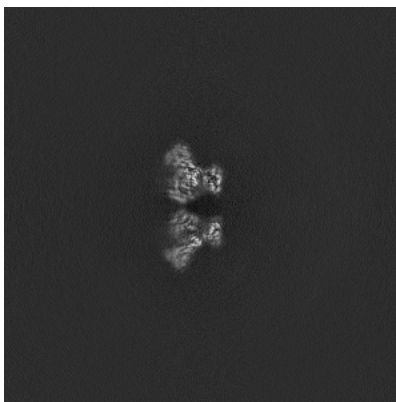


Z Index: 260

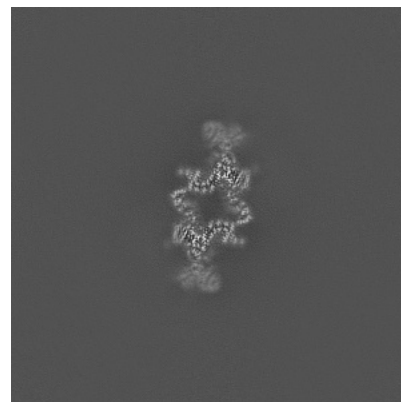
6.3.2 Raw map



X Index: 239



Y Index: 255

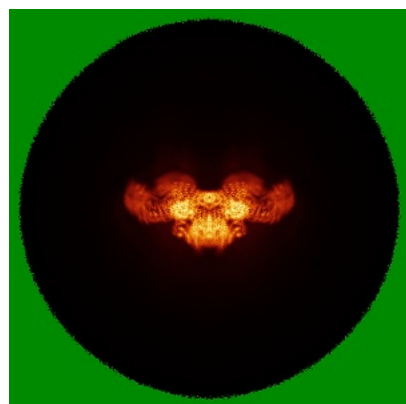


Z Index: 260

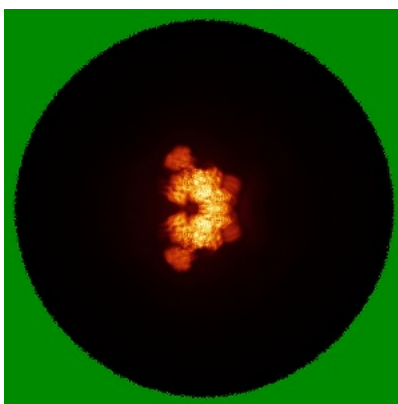
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



X

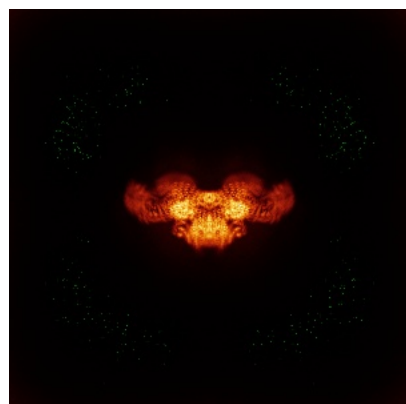


Y

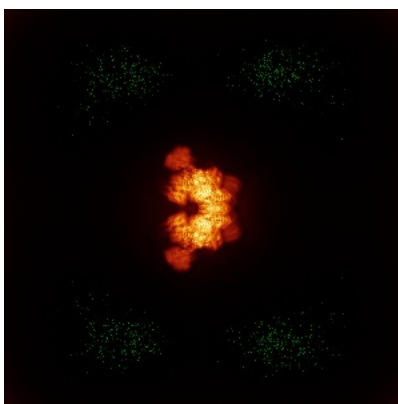


Z

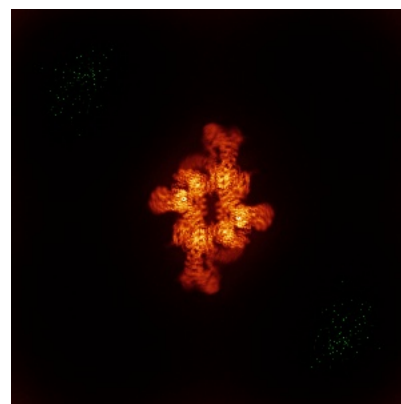
6.4.2 Raw map



X



Y

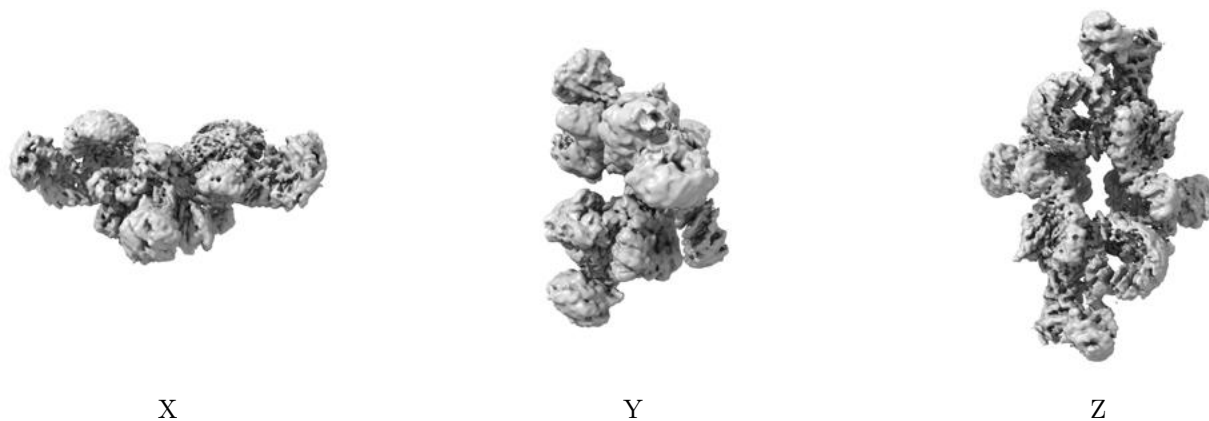


Z

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.

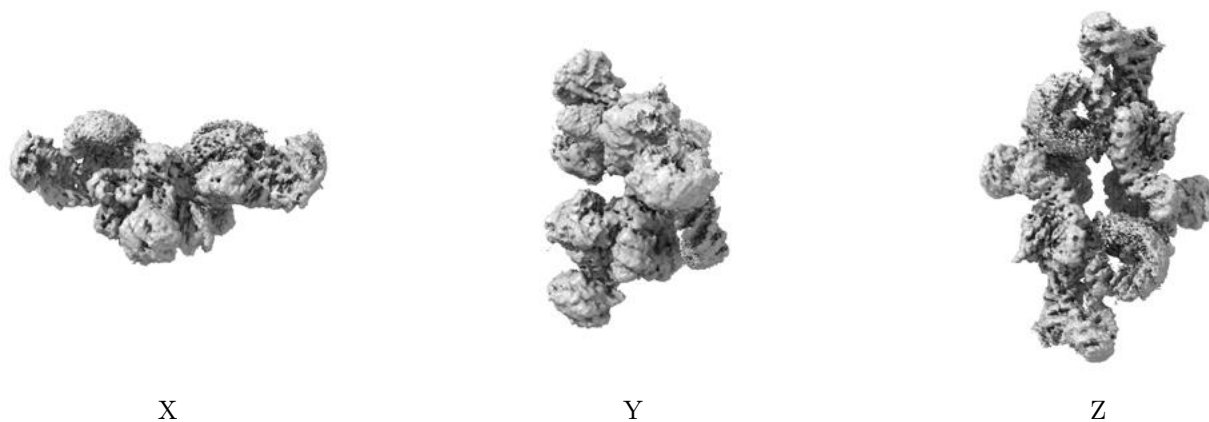
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.13. 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.

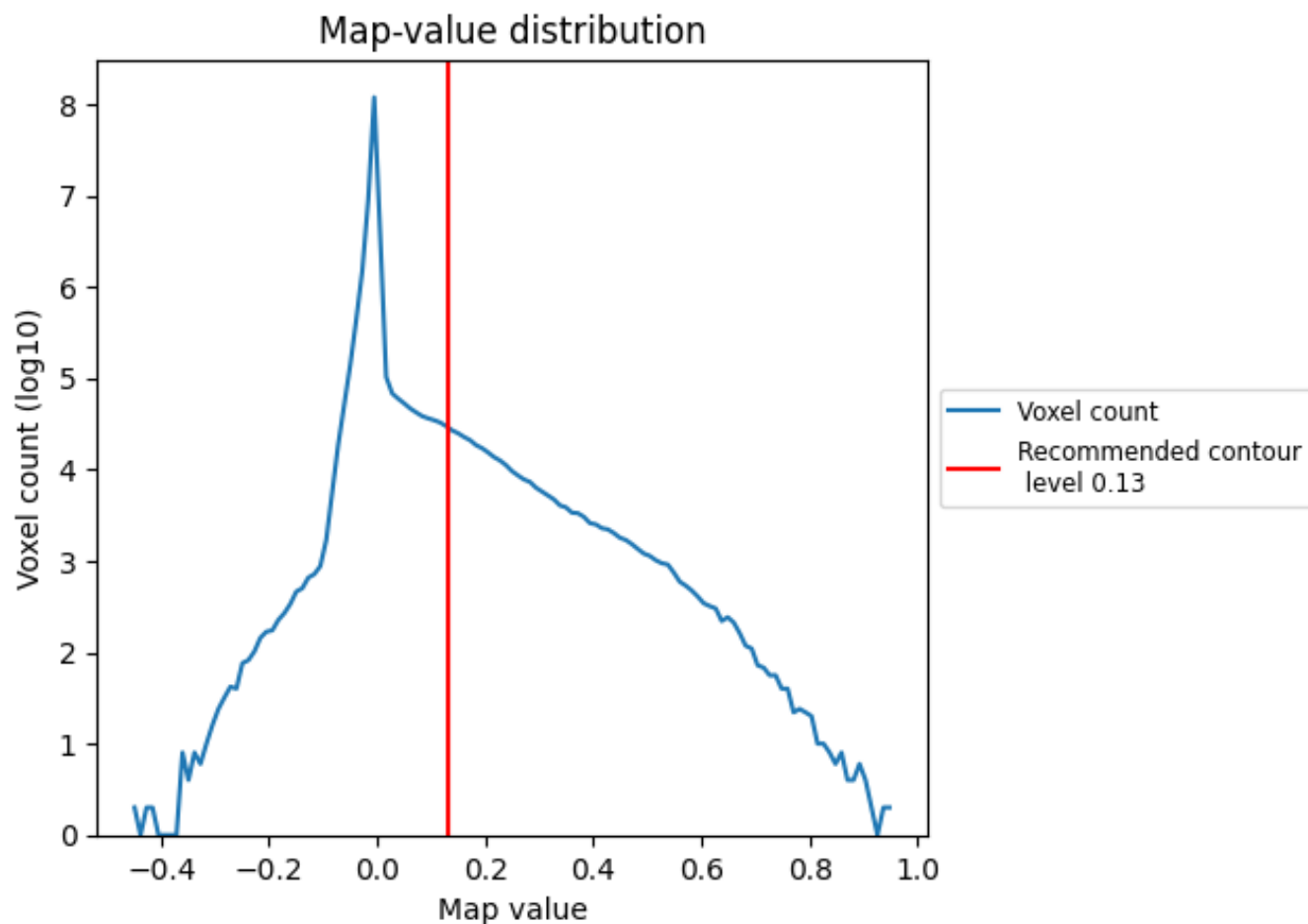
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

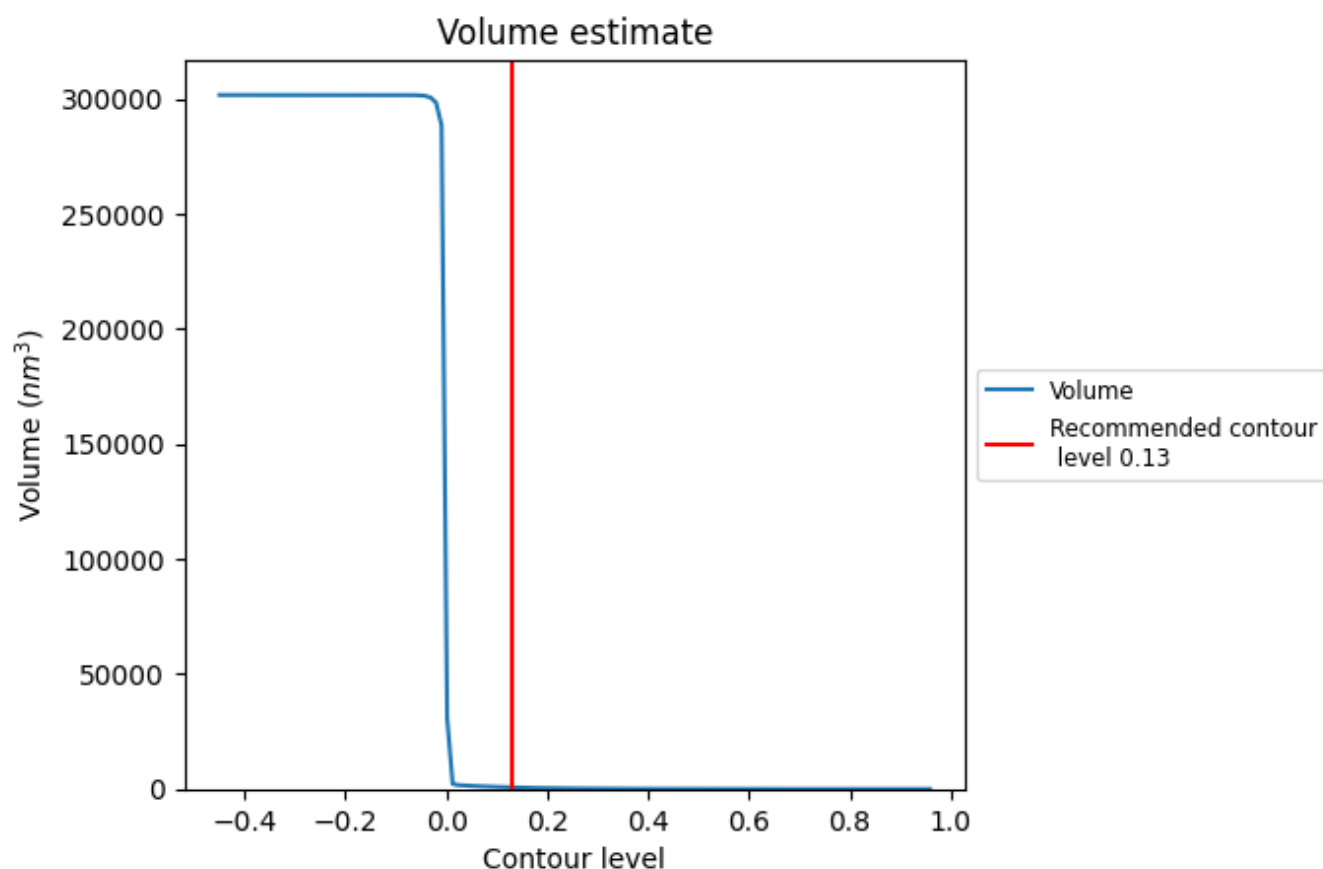
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

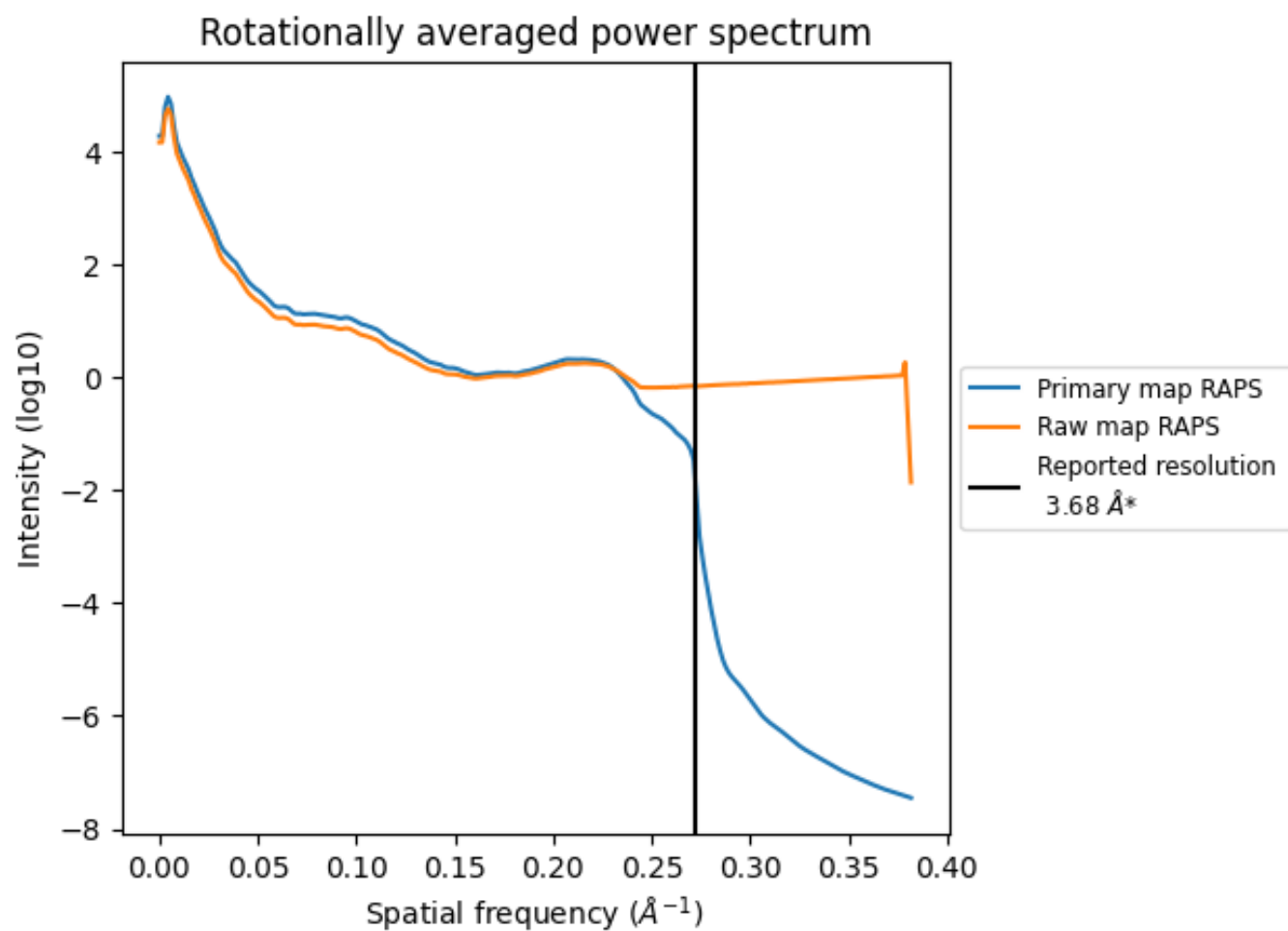
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 694 nm³; this corresponds to an approximate mass of 627 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 ⓘ

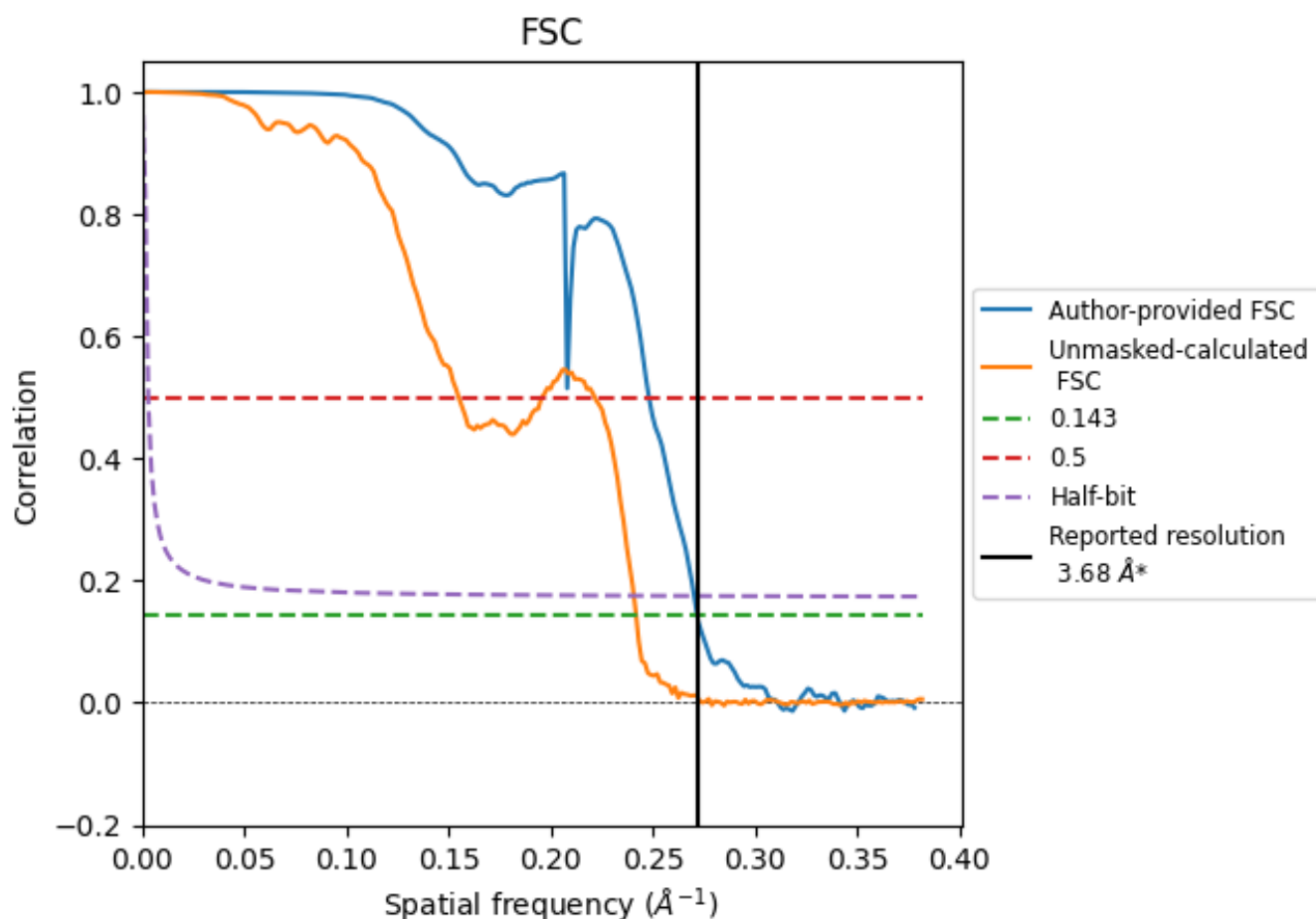


*Reported resolution corresponds to spatial frequency of 0.272 Å⁻¹

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 0.272 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

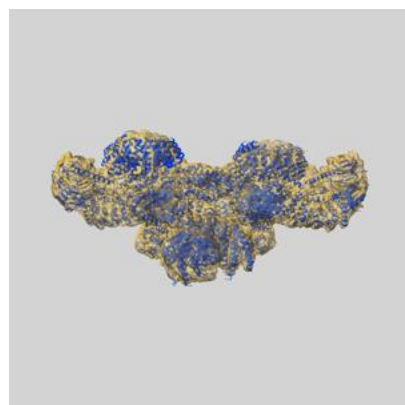
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.68	-	-
Author-provided FSC curve	3.68	4.03	3.70
Unmasked-calculated*	4.14	6.46	4.16

*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 4.14 differs from the reported value 3.68 by more than 10 %

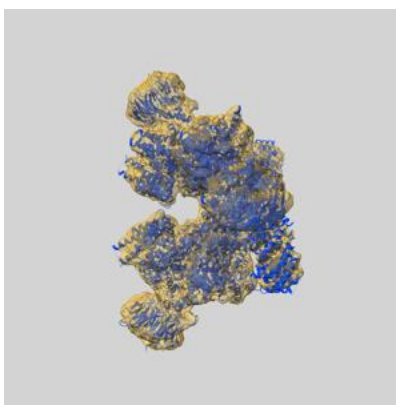
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-50183 and PDB model 9F44. Per-residue inclusion information can be found in section 3 on page 7.

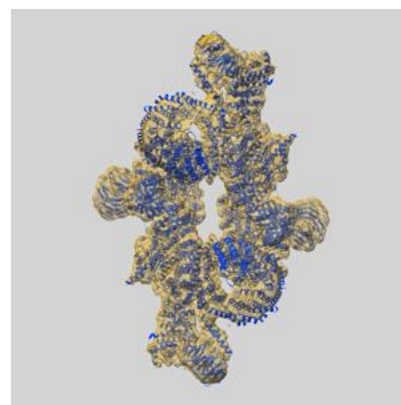
9.1 Map-model overlay [i](#)



X



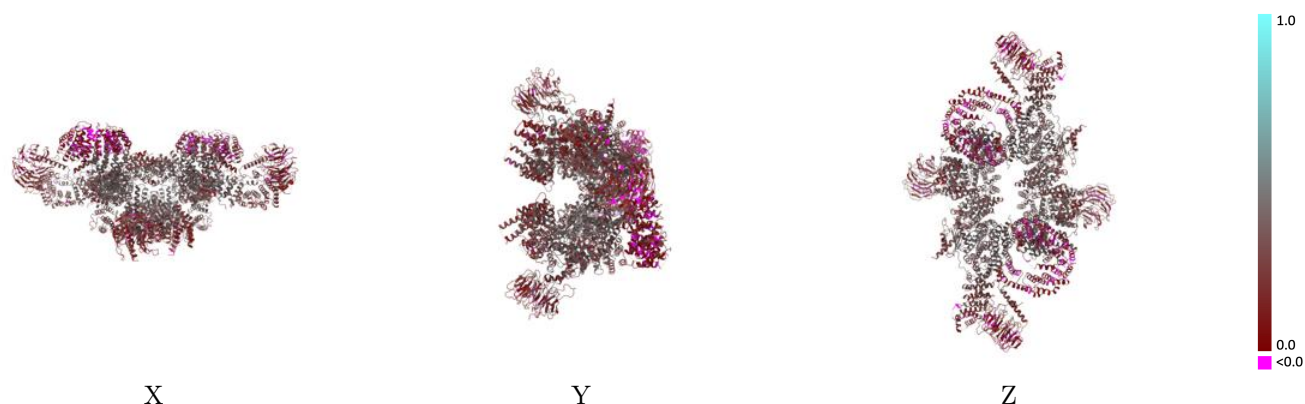
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.13 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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

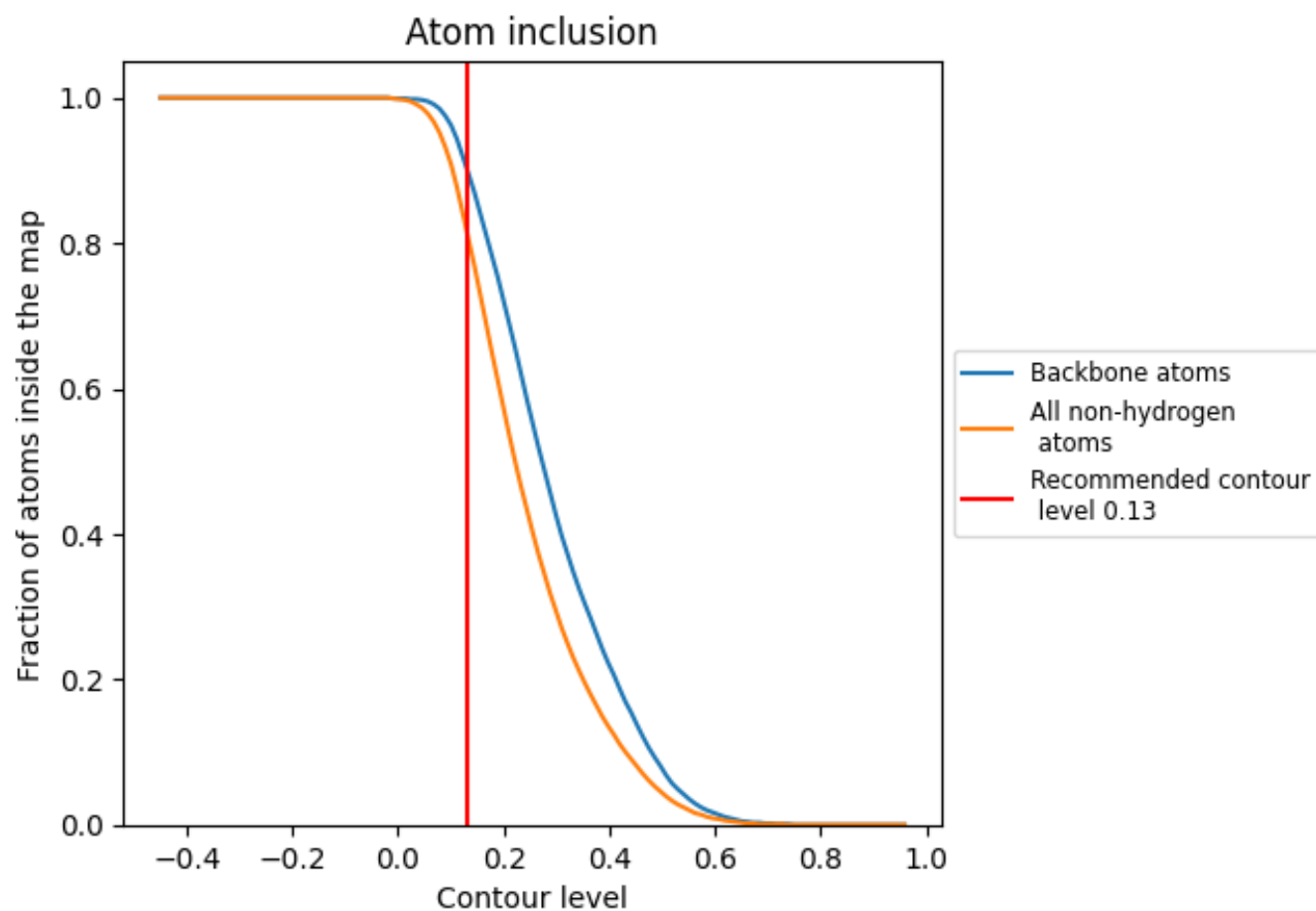


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 82% 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 (0.13) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8180	0.2800
A	0.8030	0.2910
B	0.8020	0.2920
C	0.8750	0.2080
D	0.8730	0.2070
E	0.8350	0.2760
F	0.8350	0.2750
G	0.8210	0.3510
H	0.8210	0.3460

