



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 04:07 AM UTC

PDB ID : 9BI4 / pdb_00009bi4
EMDB ID : EMD-44558
Title : cryo EM structure of dsDNA bound Mre11-Rad50 complex
Authors : Yu, Y.; Patel, D.J.
Deposited on : 2024-04-22
Resolution : 3.20 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

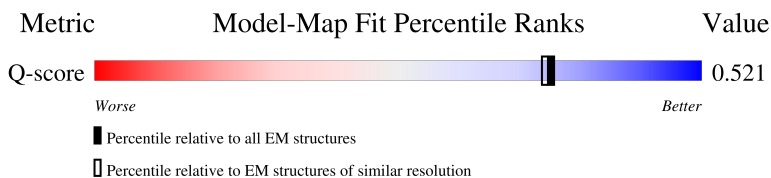
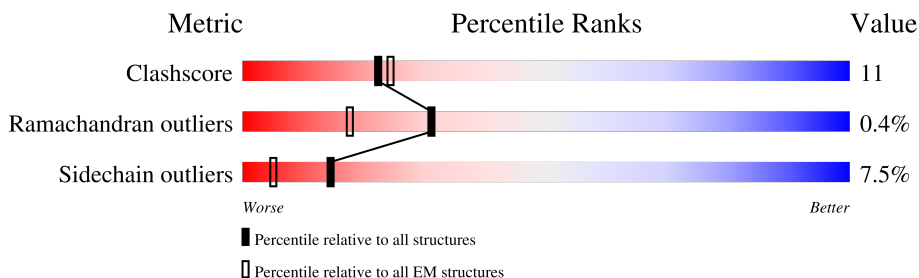
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 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	15020 (2.70 - 3.70)

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	1312	
1	D	1312	
2	A	706	
2	B	706	

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Mol	Chain	Length	Quality of chain
3	E	83	 14% 12% 73%
4	F	83	 17% 10% 73%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15869 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 DNA repair protein RAD50.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	453	Total	C	N	O	S	0	0
			3620	2278	631	694	17		
1	D	459	Total	C	N	O	S	0	0
			3676	2315	642	702	17		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1235	GLN	GLU	engineered mutation	UNP P12753
D	1235	GLN	GLU	engineered mutation	UNP P12753

- Molecule 2 is a protein called Double-strand break repair protein MRE11.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	476	Total	C	N	O	S	0	0
			3748	2372	636	720	20		
2	A	493	Total	C	N	O	S	0	0
			3855	2436	656	742	21		

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	693	TYR	-	expression tag	UNP P32829
B	694	ASP	-	expression tag	UNP P32829
B	695	TYR	-	expression tag	UNP P32829
B	696	LYS	-	expression tag	UNP P32829
B	697	ASP	-	expression tag	UNP P32829
B	698	ASP	-	expression tag	UNP P32829
B	699	ASP	-	expression tag	UNP P32829
B	700	ASP	-	expression tag	UNP P32829
B	701	LYS	-	expression tag	UNP P32829
B	702	HIS	-	expression tag	UNP P32829
B	703	HIS	-	expression tag	UNP P32829

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Chain	Residue	Modelled	Actual	Comment	Reference
B	704	HIS	-	expression tag	UNP P32829
B	705	HIS	-	expression tag	UNP P32829
B	706	HIS	-	expression tag	UNP P32829
A	693	TYR	-	expression tag	UNP P32829
A	694	ASP	-	expression tag	UNP P32829
A	695	TYR	-	expression tag	UNP P32829
A	696	LYS	-	expression tag	UNP P32829
A	697	ASP	-	expression tag	UNP P32829
A	698	ASP	-	expression tag	UNP P32829
A	699	ASP	-	expression tag	UNP P32829
A	700	ASP	-	expression tag	UNP P32829
A	701	LYS	-	expression tag	UNP P32829
A	702	HIS	-	expression tag	UNP P32829
A	703	HIS	-	expression tag	UNP P32829
A	704	HIS	-	expression tag	UNP P32829
A	705	HIS	-	expression tag	UNP P32829
A	706	HIS	-	expression tag	UNP P32829

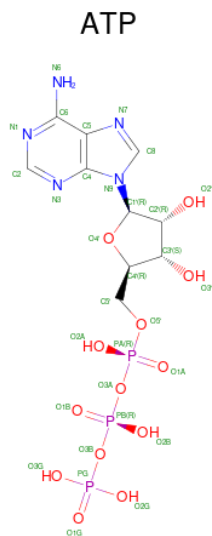
- Molecule 3 is a DNA chain called one strand of dsDNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	22	Total	C	N	O	P	0	0
			462	220	110	110	22		

- Molecule 4 is a DNA chain called second strand of dsDNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	22	Total	C	N	O	P	0	0
			440	220	44	154	22		

- Molecule 5 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					AltConf
5	C	1	Total 31	C 10	N 5	O 13	P 3	0
5	D	1	Total 31	C 10	N 5	O 13	P 3	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	AltConf
6	C	1	Total Mg 1 1	0
6	D	1	Total Mg 1 1	0

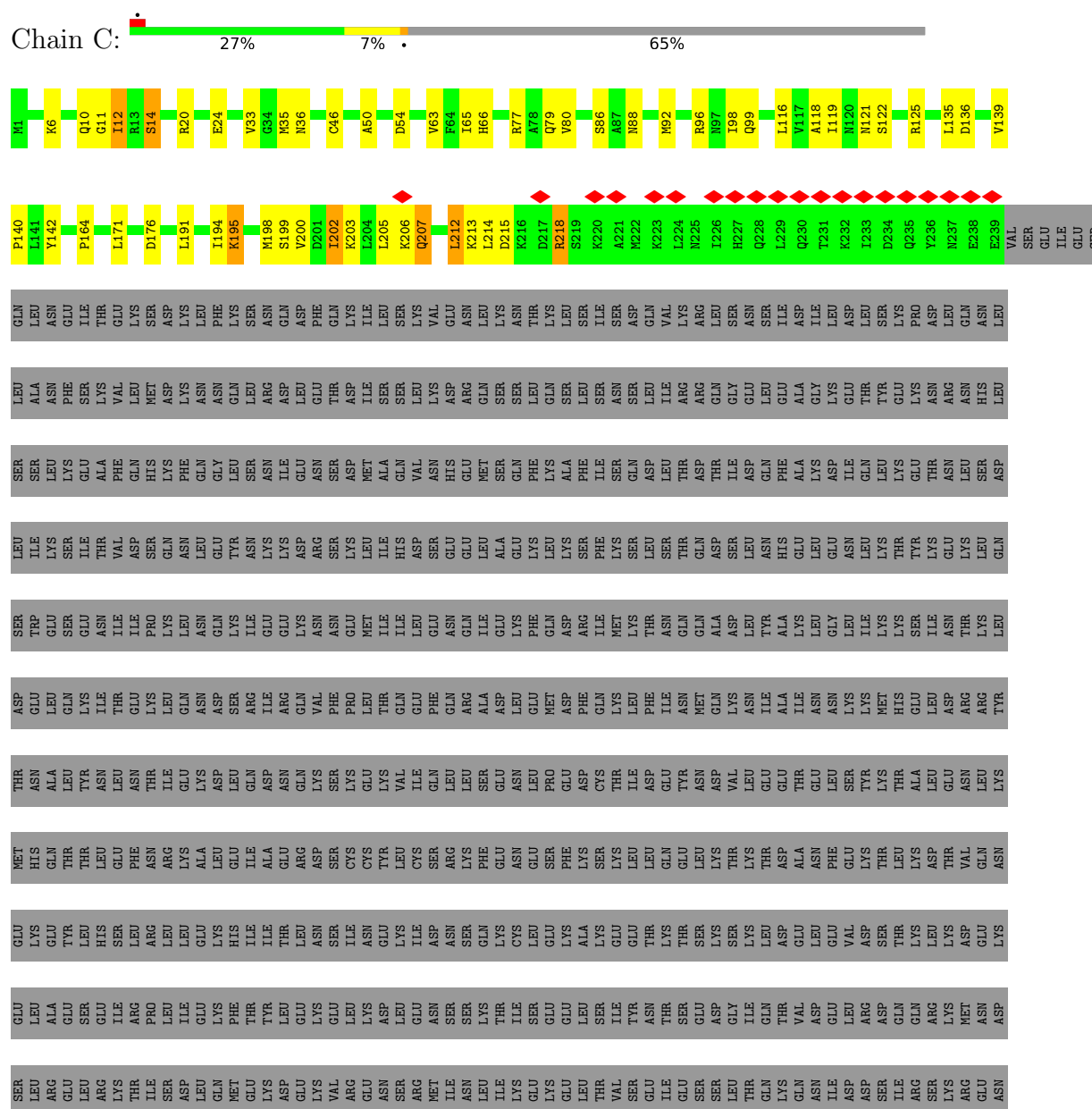
- Molecule 7 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

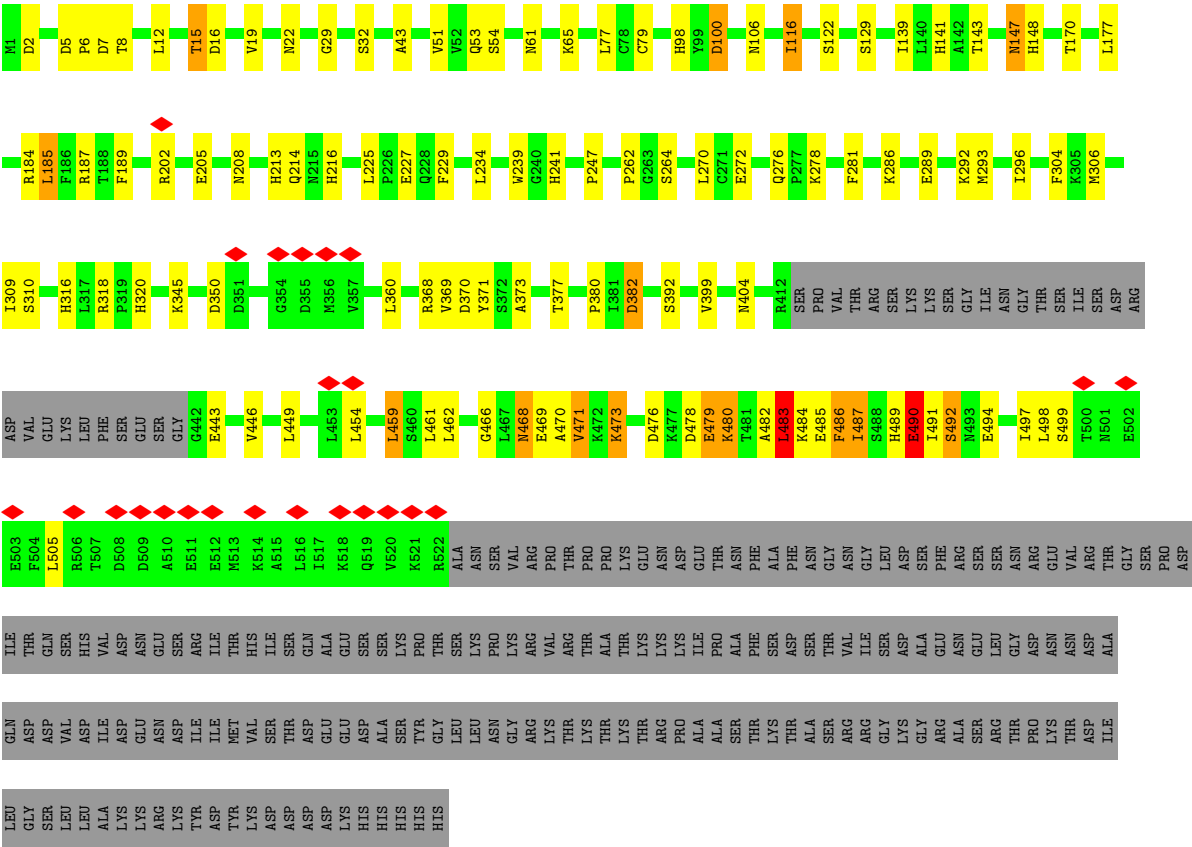
Mol	Chain	Residues	Atoms		AltConf
7	B	2	Total 2	Mn 2	0
7	A	2	Total 2	Mn 2	0

3 Residue-property plots

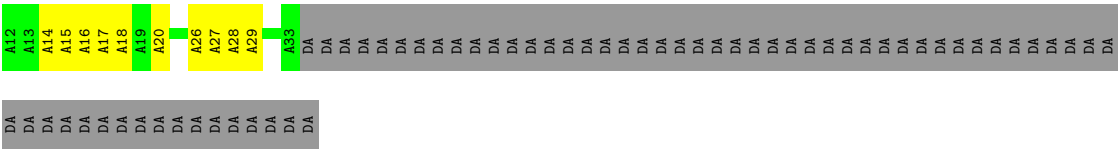
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: DNA repair protein RAD50

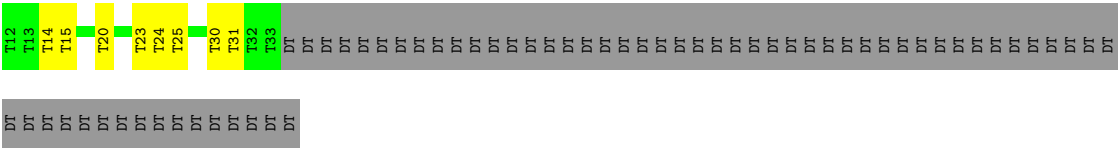




● Molecule 3: one strand of dsDNA



● Molecule 4: second strand of dsDNA



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	242558	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 ($e^-/\text{\AA}^2$)	51.22	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.809	Depositor
Minimum map value	-1.837	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.086	Depositor
Recommended contour level	0.28	Depositor
Map size ($\text{\AA}$)	303.24, 303.24, 303.24	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.083, 1.083, 1.083	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: MG, ATP, MN

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	0.48	1/3673 (0.0%)	0.76	3/4943 (0.1%)
1	D	0.44	0/3730	0.73	4/5017 (0.1%)
2	A	0.54	6/3938 (0.2%)	0.70	10/5343 (0.2%)
2	B	0.46	0/3832	0.70	6/5198 (0.1%)
3	E	0.24	0/527	0.32	0/810
4	F	0.25	0/483	0.53	0/744
All	All	0.47	7/16183 (0.0%)	0.70	23/22055 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	483	LEU	CA-C	-7.47	1.43	1.52
2	A	486	PHE	C-O	-6.30	1.16	1.24
2	A	479	GLU	CA-C	-6.07	1.45	1.53
1	C	207	GLN	C-O	5.68	1.30	1.24
2	A	489	HIS	C-O	-5.56	1.17	1.24
2	A	486	PHE	CA-C	-5.48	1.45	1.52
2	A	485	GLU	C-O	-5.43	1.17	1.24

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	489	HIS	N-CA-C	-11.14	97.81	111.40
2	A	476	ASP	N-CA-C	9.33	121.53	111.36
1	D	202	ILE	N-CA-CB	6.62	117.86	110.51
1	C	1109	ASN	N-CA-C	-6.60	104.01	111.07
1	D	202	ILE	CB-CA-C	-6.40	103.69	111.88
2	B	475	VAL	N-CA-C	6.27	116.31	110.42
2	A	499	SER	N-CA-C	-5.99	105.82	113.01
2	A	480	LYS	N-CA-C	5.95	123.48	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	1129	VAL	N-CA-C	5.94	116.60	110.36
2	A	490	GLU	N-CA-C	5.93	120.05	112.34
2	A	483	LEU	N-CA-C	-5.76	104.37	111.40
1	D	217	ASP	N-CA-C	-5.71	105.03	112.23
2	B	468	ASN	N-CA-C	-5.41	105.54	111.82
2	A	468	ASN	N-CA-C	-5.35	105.61	111.82
2	A	486	PHE	CB-CA-C	-5.32	101.64	110.68
2	A	482	ALA	N-CA-C	5.29	117.80	111.71
2	A	492	SER	N-CA-C	-5.27	99.58	110.80
2	B	32	SER	N-CA-C	5.20	117.02	111.36
1	C	202	ILE	CA-C-N	5.10	127.62	120.28
1	C	202	ILE	C-N-CA	5.10	127.62	120.28
2	B	140	LEU	N-CA-C	-5.05	107.25	113.41
2	B	495	VAL	CA-C-N	5.03	125.52	119.94
2	B	495	VAL	C-N-CA	5.03	125.52	119.94

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	C	3620	0	3657	99	0
1	D	3676	0	3726	93	0
2	A	3855	0	3711	86	0
2	B	3748	0	3620	113	0
3	E	462	0	243	9	0
4	F	440	0	265	6	0
5	C	31	0	12	1	0
5	D	31	0	12	8	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
All	All	15869	0	15246	354	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 11.

All (354) 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:C:1119:THR:CG2	2:B:491:ILE:HG23	1.74	1.18
2:A:473:LYS:HB3	2:A:479:GLU:OE2	1.46	1.13
1:C:1119:THR:HG21	2:B:491:ILE:HG23	1.13	1.13
1:C:191:LEU:HG	2:B:483:LEU:HD21	1.20	1.12
2:A:471:VAL:HG22	2:A:483:LEU:HD21	1.37	1.01
2:A:454:LEU:CB	2:A:468:ASN:HD21	1.73	1.00
1:C:1119:THR:HG21	2:B:491:ILE:CG2	1.93	0.97
1:D:1159:ARG:HG3	1:D:1159:ARG:HH11	1.30	0.95
2:B:232:ASP:OD1	2:B:255:LYS:HG2	1.67	0.93
1:C:1126:ASP:OD1	2:B:459:LEU:HB3	1.69	0.92
2:B:14:THR:HG21	2:B:39:VAL:HG11	1.56	0.88
2:B:65:LYS:HG2	2:A:139:ILE:HD11	1.52	0.87
1:C:1123:VAL:HG21	2:B:491:ILE:HD11	1.56	0.87
2:B:14:THR:HG21	2:B:39:VAL:CG1	2.05	0.86
1:C:191:LEU:HG	2:B:483:LEU:CD2	2.06	0.83
2:A:471:VAL:CG2	2:A:483:LEU:HD21	2.08	0.83
2:A:61:ASN:HD21	2:A:129:SER:HB3	1.44	0.82
2:A:498:LEU:HD23	2:A:498:LEU:O	1.79	0.82
1:C:218:ARG:HG3	1:C:218:ARG:HH11	1.46	0.81
2:A:484:LYS:HD2	2:A:484:LYS:O	1.81	0.81
2:B:366:ARG:HH12	2:B:404:ASN:HB2	1.46	0.81
1:C:65:ILE:HD11	1:C:98:ILE:CD1	2.13	0.79
1:D:1159:ARG:HG3	1:D:1159:ARG:NH1	1.96	0.79
1:D:1120:ARG:CA	1:D:1120:ARG:HH11	1.96	0.77
2:A:12:LEU:HD22	2:A:43:ALA:HB2	1.64	0.77
2:A:471:VAL:HG22	2:A:483:LEU:CD2	2.14	0.77
1:C:12:ILE:CD1	1:C:80:VAL:CG2	2.63	0.76
1:C:218:ARG:HG3	1:C:218:ARG:NH1	2.00	0.76
2:A:454:LEU:CB	2:A:468:ASN:ND2	2.47	0.75
1:D:205:LEU:CB	1:D:1117:LEU:CD2	2.65	0.74
1:C:50:ALA:O	1:C:96:ARG:NH1	2.24	0.71
2:A:147:ASN:N	2:A:147:ASN:OD1	2.22	0.71
1:C:1205:SER:HB2	5:D:1401:ATP:O3G	1.91	0.70
1:D:1120:ARG:HH11	1:D:1120:ARG:HA	1.56	0.70
1:C:1123:VAL:HG21	2:B:491:ILE:CD1	2.21	0.70
1:D:204:LEU:HD12	1:D:204:LEU:O	1.91	0.69
1:D:206:LYS:HD3	1:D:1117:LEU:HD11	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:ASN:HD22	1:D:1244:ASN:HD22	1.41	0.69
1:D:206:LYS:CD	1:D:1117:LEU:HD11	2.23	0.69
2:A:216:HIS:NE2	2:A:227:GLU:OE1	2.25	0.69
1:C:12:ILE:CD1	1:C:80:VAL:HG23	2.23	0.68
1:D:1118:GLN:OE1	1:D:1118:GLN:HA	1.93	0.68
1:D:204:LEU:HD12	1:D:204:LEU:C	2.18	0.68
2:B:139:ILE:CD1	2:A:65:LYS:HG2	2.24	0.68
1:C:1244:ASN:HD22	1:D:36:ASN:HD22	1.42	0.67
1:C:1118:GLN:HA	1:C:1118:GLN:OE1	1.93	0.67
1:D:199:SER:HA	1:D:202:ILE:HG13	1.76	0.67
1:D:198:MET:SD	1:D:198:MET:N	2.68	0.67
2:B:18:HIS:HE1	2:B:125:HIS:HD2	1.43	0.67
2:B:106:ASN:OD1	2:B:147:ASN:ND2	2.28	0.67
2:A:8:THR:HG22	2:A:286:LYS:HB2	1.77	0.67
1:C:1196:VAL:HG11	1:D:70:ILE:HG22	1.78	0.66
2:A:484:LYS:HD2	2:A:484:LYS:C	2.20	0.66
2:B:182:ASP:OD2	2:B:215:ASN:ND2	2.29	0.66
2:A:202:ARG:HH11	2:A:205:GLU:HB2	1.61	0.65
2:B:330:ILE:HD11	2:B:394:ARG:HH11	1.61	0.65
2:B:121:ILE:O	2:B:137:MET:HE3	1.97	0.64
1:C:205:LEU:CB	1:C:1117:LEU:HD21	2.28	0.64
1:D:79:GLN:HE21	1:D:97:ASN:HD21	1.43	0.64
1:C:12:ILE:HD13	1:C:80:VAL:CG2	2.26	0.64
2:A:310:SER:HA	2:A:370:ASP:HB3	1.80	0.64
1:C:65:ILE:HD11	1:C:98:ILE:HD12	1.78	0.64
2:A:100:ASP:N	2:A:100:ASP:OD1	2.28	0.63
2:A:471:VAL:HG23	2:A:483:LEU:HD11	1.80	0.63
1:C:205:LEU:CB	1:C:1117:LEU:CD2	2.77	0.63
1:D:1159:ARG:HG3	1:D:1159:ARG:O	1.98	0.63
1:C:77:ARG:NH2	1:C:99:GLN:OE1	2.31	0.63
1:D:39:GLY:HA2	5:D:1401:ATP:O2A	1.99	0.63
1:D:202:ILE:HG22	1:D:206:LYS:HE2	1.80	0.62
2:B:399:VAL:HG22	2:B:401:ASN:H	1.64	0.62
1:D:1151:ARG:NH1	2:A:380:PRO:O	2.33	0.62
2:B:315:PRO:O	2:B:316:HIS:ND1	2.32	0.62
2:B:257:PHE:HD1	2:B:257:PHE:O	1.83	0.62
1:C:194:ILE:CD1	2:B:484:LYS:HA	2.30	0.62
2:B:122:SER:HG	2:B:135:CYS:HG	1.47	0.61
1:C:194:ILE:HD13	2:B:484:LYS:HA	1.83	0.60
2:B:250:VAL:HG12	2:B:250:VAL:O	2.00	0.60
1:D:205:LEU:CB	1:D:1117:LEU:HD23	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:193:SER:O	1:D:197:ASP:OD1	2.20	0.59
2:B:305:LYS:NZ	2:B:343:GLU:OE2	2.29	0.59
1:D:205:LEU:CB	1:D:1117:LEU:HD21	2.32	0.59
2:B:152:VAL:HG23	2:B:152:VAL:O	2.03	0.59
2:B:300:THR:O	2:B:300:THR:HG22	2.00	0.59
1:C:1205:SER:CB	5:D:1401:ATP:O3G	2.49	0.59
1:D:1172:ARG:NH1	1:D:1174:ASP:OD2	2.36	0.59
1:C:202:ILE:HG22	1:C:206:LYS:HE2	1.83	0.59
2:B:139:ILE:HD11	2:A:65:LYS:CB	2.34	0.58
2:A:491:ILE:O	2:A:491:ILE:HG22	2.03	0.58
2:B:365:ILE:O	2:B:401:ASN:ND2	2.30	0.58
2:A:498:LEU:HD23	2:A:498:LEU:C	2.28	0.58
1:C:1149:ILE:HG13	1:C:1223:THR:HG21	1.85	0.58
1:C:1119:THR:HG23	2:B:491:ILE:HG23	1.76	0.57
2:B:61:ASN:ND2	2:B:127:ASP:OD1	2.37	0.57
1:D:197:ASP:OD1	1:D:197:ASP:N	2.37	0.57
2:B:33:TRP:CD2	2:B:70:GLN:HG2	2.39	0.57
1:C:164:PRO:HA	1:C:171:LEU:HD21	1.86	0.57
1:C:212:LEU:HD13	1:C:1110:TYR:HB2	1.86	0.57
1:D:10:GLN:O	1:D:10:GLN:HG2	2.04	0.57
2:B:345:LYS:HB2	2:B:360:LEU:HD11	1.87	0.57
1:C:12:ILE:HD11	1:C:80:VAL:HG23	1.86	0.56
2:B:66:LYS:O	2:B:66:LYS:HG3	2.04	0.56
1:C:1108:LYS:HA	1:C:1111:HIS:HB2	1.87	0.56
2:B:260:LEU:HD22	2:B:293:MET:HE1	1.87	0.56
2:B:257:PHE:HD1	2:B:257:PHE:C	2.13	0.56
1:C:212:LEU:CD1	1:C:1110:TYR:HA	2.35	0.56
1:D:1123:VAL:HG21	2:A:491:ILE:HD11	1.87	0.56
1:D:1221:SER:O	1:D:1259:ARG:NH1	2.39	0.56
2:B:466:GLY:O	2:B:470:ALA:HB2	2.06	0.56
1:D:1231:ILE:HG23	1:D:1267:LEU:HD12	1.88	0.56
1:D:1140:LYS:NZ	2:A:449:LEU:CB	2.69	0.55
2:A:486:PHE:CD1	2:A:486:PHE:C	2.85	0.55
2:B:258:ASP:OD1	2:B:258:ASP:N	2.39	0.55
1:C:65:ILE:CD1	1:C:98:ILE:CD1	2.83	0.55
1:D:28:PRO:HB3	1:D:1260:ARG:HD3	1.87	0.55
1:C:194:ILE:O	1:C:198:MET:HB2	2.06	0.55
1:C:1140:LYS:HZ1	2:B:449:LEU:CB	2.20	0.55
1:C:92:MET:HG2	1:C:118:ALA:HA	1.88	0.55
1:D:1183:LYS:NZ	4:F:20:DT:H5"	2.22	0.55
2:B:69:TYR:CD1	2:B:69:TYR:C	2.85	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:454:LEU:CB	2:B:468:ASN:HD21	2.20	0.55
1:C:1140:LYS:NZ	2:B:449:LEU:CB	2.70	0.54
2:B:474:PHE:CD1	2:B:474:PHE:C	2.85	0.54
1:C:63:VAL:O	5:C:1401:ATP:N6	2.37	0.54
1:C:1240:LEU:HA	1:D:36:ASN:HD21	1.72	0.54
3:E:26:DA:H2''	3:E:27:DA:C8	2.42	0.54
2:A:466:GLY:O	2:A:470:ALA:HB2	2.06	0.54
1:C:92:MET:HE1	1:C:142:TYR:HB3	1.89	0.54
2:A:5:ASP:OD1	2:A:6:PRO:HD2	2.07	0.54
1:D:117:VAL:HG21	1:D:125:ARG:HH21	1.73	0.54
2:B:227:GLU:O	2:B:227:GLU:HG2	2.08	0.54
1:C:1111:HIS:CD2	2:B:505:LEU:HD13	2.43	0.54
1:C:20:ARG:NH1	1:C:79:GLN:OE1	2.40	0.54
1:C:1110:TYR:CD1	1:C:1110:TYR:C	2.86	0.54
2:B:100:ASP:OD1	2:B:100:ASP:N	2.38	0.54
2:A:270:LEU:HD22	2:A:304:PHE:CE1	2.43	0.54
2:B:231:PRO:HB2	2:B:233:PHE:CE1	2.43	0.53
1:C:1185:TYR:O	1:C:1187:TYR:N	2.42	0.53
1:C:119:ILE:HD12	1:C:125:ARG:HE	1.71	0.53
4:F:24:DT:H2''	4:F:25:DT:H5''	1.88	0.53
2:A:2:ASP:OD1	2:A:2:ASP:N	2.34	0.53
1:D:1123:VAL:HG13	1:D:1127:ILE:HD12	1.91	0.53
1:D:60:LYS:NZ	3:E:20:DA:N3	2.51	0.53
2:B:139:ILE:HD11	2:A:65:LYS:HB3	1.89	0.53
2:A:141:HIS:HD2	2:A:148:HIS:H	1.55	0.53
2:B:257:PHE:C	2:B:257:PHE:CD1	2.85	0.52
1:D:1154:ASP:O	1:D:1158:LYS:HG3	2.09	0.52
2:A:19:VAL:HA	2:A:32:SER:HB2	1.90	0.52
1:C:11:GLY:O	1:C:66:HIS:HB2	2.09	0.52
1:D:1120:ARG:HH11	1:D:1120:ARG:N	2.08	0.52
2:B:481:THR:O	2:B:481:THR:OG1	2.27	0.52
1:D:1165:ASP:OD1	1:D:1165:ASP:N	2.40	0.52
1:D:1140:LYS:HZ3	2:A:449:LEU:CB	2.21	0.51
4:F:30:DT:H1'	4:F:31:DT:H5'	1.92	0.51
1:C:1146:MET:HE2	1:C:1171:ILE:HB	1.92	0.51
2:B:205:GLU:OE2	2:B:205:GLU:N	2.44	0.51
3:E:16:DA:H2''	3:E:17:DA:H5''	1.92	0.51
1:C:176:ASP:OD1	1:C:1142:HIS:NE2	2.39	0.51
2:B:388:PRO:HG3	2:B:408:PHE:HD2	1.75	0.51
2:B:271:CYS:SG	2:B:274:GLU:HG3	2.51	0.51
1:D:1205:SER:OG	1:D:1208:GLN:NE2	2.31	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:139:ILE:HD11	2:A:65:LYS:HG2	1.92	0.51
1:C:198:MET:HE3	1:C:1123:VAL:HG11	1.91	0.51
2:B:334:GLU:OE2	2:B:398:ARG:NH2	2.43	0.51
1:D:1126:ASP:HA	2:A:459:LEU:HD23	1.93	0.51
1:D:1132:LYS:HB3	1:D:1132:LYS:HZ2	1.76	0.51
2:A:306:MET:HE1	2:A:368:ARG:HG3	1.91	0.51
1:C:195:LYS:HB2	1:C:1127:ILE:HG21	1.93	0.50
2:B:65:LYS:HG2	2:A:139:ILE:CD1	2.36	0.50
1:D:206:LYS:HD2	1:D:1117:LEU:CD1	2.42	0.50
1:C:96:ARG:NH2	1:C:136:ASP:OD1	2.44	0.50
5:D:1401:ATP:H8	5:D:1401:ATP:H3'	1.76	0.50
2:A:202:ARG:HD3	2:A:205:GLU:HG2	1.94	0.50
1:D:219:SER:OG	1:D:1103:TYR:HB3	2.12	0.50
1:C:14:SER:OG	1:C:1300:SER:OG	2.29	0.50
1:D:206:LYS:CD	1:D:1117:LEU:CD1	2.90	0.50
2:B:359:GLU:CD	2:B:359:GLU:H	2.20	0.49
1:D:1111:HIS:CG	2:A:505:LEU:HD22	2.47	0.49
2:B:231:PRO:HB2	2:B:233:PHE:CD1	2.47	0.49
2:B:471:VAL:HG23	2:B:483:LEU:HD13	1.93	0.49
2:B:184:ARG:HH21	2:B:187:ARG:HE	1.60	0.49
1:C:200:VAL:O	1:C:203:LYS:HB3	2.11	0.49
1:D:1120:ARG:N	1:D:1120:ARG:NH1	2.60	0.49
1:D:67:ASP:OD1	1:D:67:ASP:N	2.44	0.49
1:C:205:LEU:CB	1:C:1117:LEU:HD23	2.42	0.49
1:D:157:HIS:CE1	1:D:159:GLU:HG2	2.46	0.49
2:B:166:GLN:HG3	2:B:171:LYS:HG2	1.95	0.49
2:B:498:LEU:HD22	2:B:504:PHE:CE1	2.47	0.49
1:C:218:ARG:HH11	1:C:218:ARG:CG	2.15	0.49
2:A:213:HIS:CD2	2:A:241:HIS:HB2	2.48	0.49
1:D:1272:HIS:O	1:D:1272:HIS:ND1	2.43	0.49
1:C:54:ASP:OD1	1:C:54:ASP:N	2.43	0.48
3:E:16:DA:H2''	3:E:17:DA:C8	2.48	0.48
3:E:28:DA:H2''	3:E:29:DA:C8	2.47	0.48
2:A:272:GLU:OE2	2:A:276:GLN:NE2	2.46	0.48
2:B:56:ASP:OD1	2:B:125:HIS:HB2	2.13	0.48
2:A:462:LEU:HD22	2:A:486:PHE:HE1	1.78	0.48
1:D:12:ILE:HG21	1:D:46:CYS:SG	2.54	0.48
1:D:1183:LYS:HZ3	4:F:20:DT:H5''	1.78	0.48
1:C:1287:ASP:OD1	1:C:1287:ASP:N	2.40	0.48
1:D:1120:ARG:CA	1:D:1120:ARG:NH1	2.74	0.48
2:B:139:ILE:O	2:B:139:ILE:HG22	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:203:GLU:OE1	2:B:204:GLY:N	2.46	0.48
1:C:6:LYS:NZ	1:C:24:GLU:OE1	2.47	0.48
1:C:65:ILE:CD1	1:C:98:ILE:HD11	2.44	0.48
1:C:1275:LYS:HG2	1:C:1279:HIS:HE1	1.79	0.47
2:B:14:THR:HG21	2:B:39:VAL:HG13	1.93	0.47
2:B:391:PHE:O	2:B:394:ARG:HG2	2.14	0.47
2:B:122:SER:OG	2:B:135:CYS:SG	2.60	0.47
1:C:86:SER:HB3	1:C:92:MET:HE2	1.95	0.47
2:B:401:ASN:HB3	2:B:405:VAL:HG22	1.97	0.47
2:A:7:ASP:HB2	2:A:286:LYS:HE2	1.97	0.47
2:A:61:ASN:ND2	2:A:129:SER:HB3	2.23	0.47
1:D:1123:VAL:HG13	1:D:1127:ILE:CD1	2.45	0.47
2:B:228:GLN:HG2	2:B:255:LYS:NZ	2.29	0.47
1:D:48:LYS:HE3	1:D:151:GLU:HA	1.97	0.47
2:B:18:HIS:CE1	2:B:125:HIS:HD2	2.27	0.47
1:C:36:ASN:HD21	1:D:1240:LEU:HA	1.79	0.46
1:D:54:ASP:OD1	1:D:54:ASP:N	2.47	0.46
1:C:191:LEU:CG	2:B:483:LEU:HD21	2.14	0.46
1:D:31:LEU:HD11	1:D:1282:ALA:HB1	1.98	0.46
2:B:184:ARG:NH2	2:B:187:ARG:HH21	2.13	0.46
2:A:316:HIS:O	2:A:316:HIS:ND1	2.48	0.46
1:C:10:GLN:OE1	1:C:20:ARG:NH1	2.48	0.46
1:C:1117:LEU:HD22	1:C:1117:LEU:HA	1.67	0.46
1:C:1205:SER:OG	5:D:1401:ATP:O3G	2.33	0.46
1:D:1127:ILE:HG12	2:A:487:ILE:CD1	2.46	0.46
1:D:1132:LYS:HB3	1:D:1132:LYS:NZ	2.31	0.46
2:B:28:THR:O	2:B:28:THR:OG1	2.24	0.46
1:C:1164:THR:O	1:C:1164:THR:OG1	2.33	0.46
1:D:1193:LYS:O	1:D:1196:VAL:HG12	2.15	0.46
1:D:2:SER:OG	1:D:3:ALA:N	2.48	0.46
2:B:139:ILE:HD11	2:A:65:LYS:CG	2.45	0.46
1:D:1153:ILE:HD13	1:D:1216:ILE:HG12	1.98	0.46
2:B:476:ASP:N	2:B:476:ASP:OD1	2.47	0.46
1:D:93:ILE:HB	1:D:117:VAL:HG22	1.98	0.45
1:D:1126:ASP:HB3	2:A:462:LEU:HD12	1.97	0.45
1:D:1126:ASP:HB3	2:A:459:LEU:HD22	1.99	0.45
2:B:101:GLU:OE1	2:A:77:LEU:HD11	2.16	0.45
2:A:213:HIS:CD2	2:A:213:HIS:O	2.70	0.45
1:C:1262:GLN:HG2	1:C:1263:LYS:H	1.81	0.45
1:C:212:LEU:HD22	1:C:212:LEU:HA	1.70	0.45
2:B:270:LEU:HD22	2:B:401:ASN:HD21	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:106:ASN:OD1	2:A:106:ASN:N	2.50	0.45
2:B:184:ARG:HH21	2:B:187:ARG:HH21	1.64	0.45
1:D:1149:ILE:O	1:D:1153:ILE:HG13	2.17	0.45
2:B:18:HIS:HE1	2:B:125:HIS:CD2	2.28	0.45
1:C:1111:HIS:ND1	2:B:505:LEU:HD22	2.31	0.45
2:A:289:GLU:N	2:A:289:GLU:OE2	2.50	0.45
2:A:292:LYS:HB3	2:A:292:LYS:HE2	1.85	0.45
2:A:79:CYS:HA	2:A:116:ILE:HG23	1.98	0.45
2:A:262:PRO:HA	2:A:281:PHE:CE1	2.52	0.45
1:C:1218:LEU:HD23	1:C:1231:ILE:HD13	1.99	0.44
5:D:1401:ATP:C3'	5:D:1401:ATP:C8	3.01	0.44
2:A:241:HIS:O	2:A:241:HIS:CG	2.70	0.44
2:A:469:GLU:O	2:A:469:GLU:HG3	2.16	0.44
1:D:1111:HIS:CD2	2:A:505:LEU:HB3	2.52	0.44
2:B:388:PRO:HG3	2:B:408:PHE:CD2	2.52	0.44
2:A:241:HIS:O	2:A:241:HIS:CD2	2.70	0.44
2:A:318:ARG:HB3	2:A:320:HIS:CE1	2.53	0.44
1:C:198:MET:HB2	1:C:198:MET:HE2	1.85	0.44
1:C:198:MET:SD	1:C:1123:VAL:HG11	2.58	0.44
1:C:1111:HIS:CG	2:B:505:LEU:HD13	2.53	0.44
1:D:104:LYS:HD3	1:D:104:LYS:HA	1.77	0.44
2:B:469:GLU:HG3	2:B:469:GLU:O	2.16	0.44
1:D:1233:LEU:HB3	1:D:1236:PRO:HG3	2.00	0.44
2:A:213:HIS:NE2	2:A:241:HIS:HB2	2.32	0.44
2:A:22:ASN:O	2:A:29:GLY:HA2	2.17	0.44
2:A:189:PHE:CD1	2:A:229:PHE:HB3	2.52	0.44
1:C:86:SER:OG	1:C:88:ASN:OD1	2.26	0.44
1:C:194:ILE:O	1:C:198:MET:HE2	2.18	0.44
1:D:29:LEU:HD13	1:D:1267:LEU:HD23	2.00	0.44
3:E:14:DA:H2''	3:E:15:DA:C8	2.53	0.44
1:D:59:SER:HA	1:D:63:VAL:HG22	1.99	0.44
2:A:264:SER:O	2:A:278:LYS:HD3	2.18	0.44
1:C:195:LYS:HB3	1:C:195:LYS:HE2	1.51	0.44
1:D:79:GLN:HE21	1:D:97:ASN:ND2	2.14	0.44
2:B:141:HIS:ND1	2:B:147:ASN:HA	2.33	0.44
2:B:151:LYS:HE2	2:B:151:LYS:HB3	1.88	0.44
2:B:179:ALA:HB2	2:B:213:HIS:HB3	1.99	0.44
1:D:152:TYR:OH	1:D:177:GLU:OE2	2.31	0.43
3:E:14:DA:H2''	3:E:15:DA:H8	1.83	0.43
1:D:183:LYS:HE3	1:D:183:LYS:HB2	1.87	0.43
2:B:14:THR:HG22	2:B:280:VAL:CG1	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:177:LEU:HD21	2:A:185:LEU:HD11	2.00	0.43
2:B:71:VAL:O	2:B:71:VAL:HG12	2.17	0.43
1:D:5:TYR:HB3	1:D:83:ALA:O	2.17	0.43
1:D:176:ASP:OD1	1:D:1142:HIS:NE2	2.50	0.43
1:D:1306:ASP:OD1	1:D:1307:ILE:N	2.51	0.43
2:A:214:GLN:O	2:A:239:TRP:NE1	2.46	0.43
4:F:23:DT:H2'	4:F:24:DT:C5	2.54	0.43
1:D:1246:GLU:O	1:D:1250:LYS:HG2	2.19	0.43
2:B:15:THR:HB	2:B:54:SER:HB2	2.01	0.43
2:B:126:ASP:HB2	2:B:136:PRO:HG2	2.00	0.43
2:B:454:LEU:CB	2:B:468:ASN:ND2	2.80	0.43
1:D:1127:ILE:HG12	2:A:487:ILE:HD11	2.00	0.43
5:D:1401:ATP:H3'	5:D:1401:ATP:C8	2.54	0.43
3:E:27:DA:H1'	3:E:28:DA:H5'	2.00	0.43
2:A:247:PRO:HA	2:A:281:PHE:CZ	2.54	0.43
2:A:483:LEU:HD12	2:A:483:LEU:HA	1.81	0.43
1:C:1165:ASP:OD1	1:C:1166:ILE:HG13	2.19	0.43
1:C:33:VAL:O	1:C:1291:LYS:HA	2.19	0.42
1:D:127:THR:O	1:D:127:THR:OG1	2.36	0.42
2:B:201:MET:SD	2:B:201:MET:N	2.89	0.42
2:A:490:GLU:OE2	2:A:490:GLU:N	2.51	0.42
1:C:1193:LYS:O	1:C:1196:VAL:HG12	2.19	0.42
1:C:12:ILE:HG21	1:C:46:CYS:SG	2.60	0.42
1:D:1252:LEU:O	1:D:1256:ILE:HG13	2.20	0.42
1:D:1158:LYS:NZ	2:A:371:TYR:CZ	2.88	0.42
1:D:1191:MET:HE3	1:D:1193:LYS:HD2	2.02	0.42
1:D:183:LYS:NZ	2:A:443:GLU:CB	2.83	0.42
2:B:180:VAL:HG23	2:B:185:LEU:HB2	2.01	0.42
2:B:81:GLY:O	2:B:113:ASN:HB3	2.19	0.42
2:B:231:PRO:CB	2:B:233:PHE:CE1	3.02	0.42
1:C:194:ILE:HD11	2:B:484:LYS:HA	2.01	0.42
2:B:14:THR:HG22	2:B:280:VAL:HG12	2.02	0.42
1:C:1244:ASN:HD22	1:D:36:ASN:ND2	2.14	0.42
2:B:330:ILE:HD11	2:B:394:ARG:NH1	2.33	0.42
1:C:1164:THR:HA	2:B:387:ASN:ND2	2.36	0.41
2:B:318:ARG:HA	2:B:318:ARG:HD3	1.88	0.41
2:B:498:LEU:HD22	2:B:504:PHE:CZ	2.55	0.41
2:A:345:LYS:HB2	2:A:360:LEU:HD11	2.02	0.41
1:C:1283:ALA:O	2:B:181:ARG:HD2	2.20	0.41
2:A:309:ILE:O	2:A:369:VAL:HA	2.20	0.41
1:C:214:LEU:HA	1:C:214:LEU:HD23	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:7:ASP:O	2:A:170:THR:OG1	2.37	0.41
2:A:98:HIS:O	2:A:98:HIS:ND1	2.51	0.41
1:D:1112:LYS:HB2	1:D:1112:LYS:HE3	1.48	0.41
1:C:212:LEU:HD12	1:C:1110:TYR:HA	2.03	0.41
2:A:15:THR:HB	2:A:54:SER:OG	2.21	0.41
1:C:213:LYS:HE2	1:C:213:LYS:HB3	1.89	0.41
2:A:184:ARG:O	2:A:184:ARG:HD3	2.21	0.41
1:C:139:VAL:HB	1:C:140:PRO:HD3	2.02	0.41
2:B:11:ILE:HA	2:B:50:MET:O	2.20	0.41
2:A:15:THR:OG1	2:A:16:ASP:N	2.53	0.41
4:F:14:DT:H2'	4:F:15:DT:H71	2.02	0.41
1:C:116:LEU:HB2	1:C:135:LEU:HD11	2.02	0.41
1:D:1193:LYS:HD3	1:D:1198:LEU:HD12	2.03	0.41
1:C:36:ASN:HD22	1:D:1244:ASN:ND2	2.13	0.40
1:C:121:ASN:OD1	1:C:122:SER:N	2.54	0.40
1:C:1123:VAL:CG2	2:B:491:ILE:HD11	2.37	0.40
1:C:1192:TYR:HA	1:C:1196:VAL:O	2.21	0.40
1:D:39:GLY:CA	5:D:1401:ATP:O2A	2.69	0.40
1:D:203:LYS:HB3	1:D:203:LYS:HE3	1.79	0.40
2:B:228:GLN:NE2	2:B:228:GLN:O	2.54	0.40
2:B:376:ASN:OD1	2:B:377:THR:N	2.55	0.40
2:A:208:ASN:HB2	2:A:234:LEU:HA	2.03	0.40
2:A:373:ALA:HB2	2:A:382:ASP:OD1	2.22	0.40
1:C:1126:ASP:HB3	2:B:462:LEU:HD12	2.04	0.40
1:D:1111:HIS:NE2	2:A:505:LEU:HB3	2.36	0.40
2:B:143:THR:HG23	2:A:143:THR:HB	2.03	0.40
2:B:228:GLN:O	2:B:228:GLN:CG	2.70	0.40
3:E:17:DA:H2''	3:E:18:DA:H5'	2.03	0.40
1:C:1108:LYS:HA	1:C:1111:HIS:CG	2.56	0.40
2:B:12:LEU:HD22	2:B:43:ALA:HB2	2.04	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	447/1312 (34%)	419 (94%)	27 (6%)	1 (0%)	43	73
1	D	453/1312 (34%)	424 (94%)	28 (6%)	1 (0%)	43	73
2	A	489/706 (69%)	460 (94%)	26 (5%)	3 (1%)	21	56
2	B	472/706 (67%)	449 (95%)	20 (4%)	3 (1%)	21	56
All	All	1861/4036 (46%)	1752 (94%)	101 (5%)	8 (0%)	31	62

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	445	GLU
2	B	457	MET
2	A	478	ASP
2	B	446	VAL
2	A	480	LYS
2	A	446	VAL
1	C	1104	LYS
1	D	1104	LYS

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	405/1234 (33%)	381 (94%)	24 (6%)	18	50
1	D	412/1234 (33%)	372 (90%)	40 (10%)	8	32
2	A	422/635 (66%)	394 (93%)	28 (7%)	15	47
2	B	415/635 (65%)	383 (92%)	32 (8%)	12	41
All	All	1654/3738 (44%)	1530 (92%)	124 (8%)	14	42

All (124) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	12	ILE
1	C	14	SER
1	C	35	MET
1	C	195	LYS
1	C	199	SER
1	C	207	GLN
1	C	212	LEU
1	C	215	ASP
1	C	218	ARG
1	C	1106	ILE
1	C	1113	GLU
1	C	1116	GLU
1	C	1117	LEU
1	C	1118	GLN
1	C	1119	THR
1	C	1123	VAL
1	C	1124	THR
1	C	1148	ASP
1	C	1152	ILE
1	C	1156	LEU
1	C	1165	ASP
1	C	1227	ASN
1	C	1238	THR
1	C	1277	LEU
1	D	33	VAL
1	D	76	ILE
1	D	85	THR
1	D	117	VAL
1	D	171	LEU
1	D	176	ASP
1	D	177	GLU
1	D	196	LYS
1	D	197	ASP
1	D	198	MET
1	D	199	SER
1	D	202	ILE
1	D	203	LYS
1	D	204	LEU
1	D	206	LYS
1	D	207	GLN
1	D	208	SER
1	D	210	GLU
1	D	211	HIS

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Mol	Chain	Res	Type
1	D	213	LYS
1	D	214	LEU
1	D	216	LYS
1	D	219	SER
1	D	1106	ILE
1	D	1108	LYS
1	D	1111	HIS
1	D	1112	LYS
1	D	1117	LEU
1	D	1118	GLN
1	D	1119	THR
1	D	1120	ARG
1	D	1121	SER
1	D	1123	VAL
1	D	1124	THR
1	D	1158	LYS
1	D	1159	ARG
1	D	1189	VAL
1	D	1267	LEU
1	D	1273	ASP
1	D	1310	VAL
2	B	15	THR
2	B	16	ASP
2	B	28	THR
2	B	35	THR
2	B	51	VAL
2	B	60	VAL
2	B	66	LYS
2	B	70	GLN
2	B	74	THR
2	B	85	CYS
2	B	100	ASP
2	B	105	VAL
2	B	137	MET
2	B	146	ILE
2	B	147	ASN
2	B	184	ARG
2	B	191	ASP
2	B	198	VAL
2	B	254	ILE
2	B	257	PHE
2	B	258	ASP

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Mol	Chain	Res	Type
2	B	296	ILE
2	B	306	MET
2	B	307	LYS
2	B	314	VAL
2	B	357	VAL
2	B	382	ASP
2	B	459	LEU
2	B	461	LEU
2	B	471	VAL
2	B	473	LYS
2	B	476	ASP
2	A	15	THR
2	A	51	VAL
2	A	53	GLN
2	A	100	ASP
2	A	116	ILE
2	A	122	SER
2	A	147	ASN
2	A	185	LEU
2	A	187	ARG
2	A	225	LEU
2	A	293	MET
2	A	296	ILE
2	A	350	ASP
2	A	377	THR
2	A	382	ASP
2	A	392	SER
2	A	399	VAL
2	A	404	ASN
2	A	459	LEU
2	A	461	LEU
2	A	471	VAL
2	A	473	LYS
2	A	483	LEU
2	A	487	ILE
2	A	490	GLU
2	A	492	SER
2	A	494	GLU
2	A	497	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	36	ASN
1	C	66	HIS
1	C	91	ASN
1	C	190	ASN
1	C	230	GLN
1	C	1089	GLN
1	C	1208	GLN
1	C	1235	GLN
1	C	1253	HIS
1	C	1264	ASN
1	C	1279	HIS
1	D	18	ASN
1	D	36	ASN
1	D	79	GLN
1	D	115	GLN
1	D	207	GLN
1	D	230	GLN
1	D	1089	GLN
1	D	1208	GLN
1	D	1239	ASN
1	D	1253	HIS
2	B	45	ASN
2	B	46	ASN
2	B	125	HIS
2	B	166	GLN
2	B	228	GLN
2	B	261	GLN
2	A	17	ASN
2	A	37	HIS
2	A	46	ASN
2	A	61	ASN
2	A	141	HIS
2	A	148	HIS
2	A	261	GLN
2	A	320	HIS
2	A	376	ASN
2	A	384	GLN
2	A	393	ASN
2	A	404	ASN
2	A	468	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	ATP	D	1401	6	32,33,33	1.43	4 (12%)	48,52,52	1.85	10 (20%)
5	ATP	C	1401	6	32,33,33	1.40	4 (12%)	48,52,52	1.80	9 (18%)

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	ATP	D	1401	6	-	5/22/38/38	0/3/3/3
5	ATP	C	1401	6	-	7/22/38/38	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	1401	ATP	C5-C4	3.42	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	1401	ATP	C5-C4	3.30	1.45	1.39
5	C	1401	ATP	C5-N7	-3.19	1.33	1.39
5	D	1401	ATP	C5-N7	-3.17	1.33	1.39
5	D	1401	ATP	C8-N9	-2.95	1.32	1.37
5	C	1401	ATP	C8-N9	-2.71	1.32	1.37
5	D	1401	ATP	C4-N9	-2.55	1.32	1.37
5	C	1401	ATP	C4-N9	-2.52	1.32	1.37

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	1401	ATP	C5-C4-N3	-5.88	118.62	126.72
5	C	1401	ATP	C5-C4-N3	-5.83	118.69	126.72
5	D	1401	ATP	N3-C4-N9	4.80	135.32	127.17
5	C	1401	ATP	N3-C4-N9	4.59	134.97	127.17
5	C	1401	ATP	C2-N3-C4	3.65	120.75	111.83
5	D	1401	ATP	C2-N3-C4	3.64	120.71	111.83
5	C	1401	ATP	N3-C2-N1	-3.53	123.24	128.58
5	C	1401	ATP	C4-C5-N7	-3.48	106.60	110.58
5	D	1401	ATP	C4-C5-N7	-3.45	106.64	110.58
5	D	1401	ATP	C4-N9-C8	3.28	109.19	105.74
5	D	1401	ATP	N3-C2-N1	-3.26	123.65	128.58
5	D	1401	ATP	O3G-PG-O2G	3.05	119.25	107.80
5	D	1401	ATP	C2'-C1'-N9	-2.89	106.13	113.30
5	C	1401	ATP	C4-N9-C8	2.88	108.76	105.74
5	D	1401	ATP	C5-N7-C8	2.65	107.61	103.45
5	C	1401	ATP	C5-N7-C8	2.60	107.54	103.45
5	D	1401	ATP	N9-C8-N7	-2.42	110.50	113.94
5	C	1401	ATP	O3G-PG-O2G	2.38	116.73	107.80
5	C	1401	ATP	N9-C8-N7	-2.24	110.76	113.94

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	1401	ATP	C5'-O5'-PA-O1A
5	C	1401	ATP	C5'-O5'-PA-O2A
5	C	1401	ATP	C5'-O5'-PA-O3A
5	C	1401	ATP	O4'-C4'-C5'-O5'
5	C	1401	ATP	C3'-C4'-C5'-O5'
5	D	1401	ATP	C5'-O5'-PA-O2A
5	D	1401	ATP	C5'-O5'-PA-O3A

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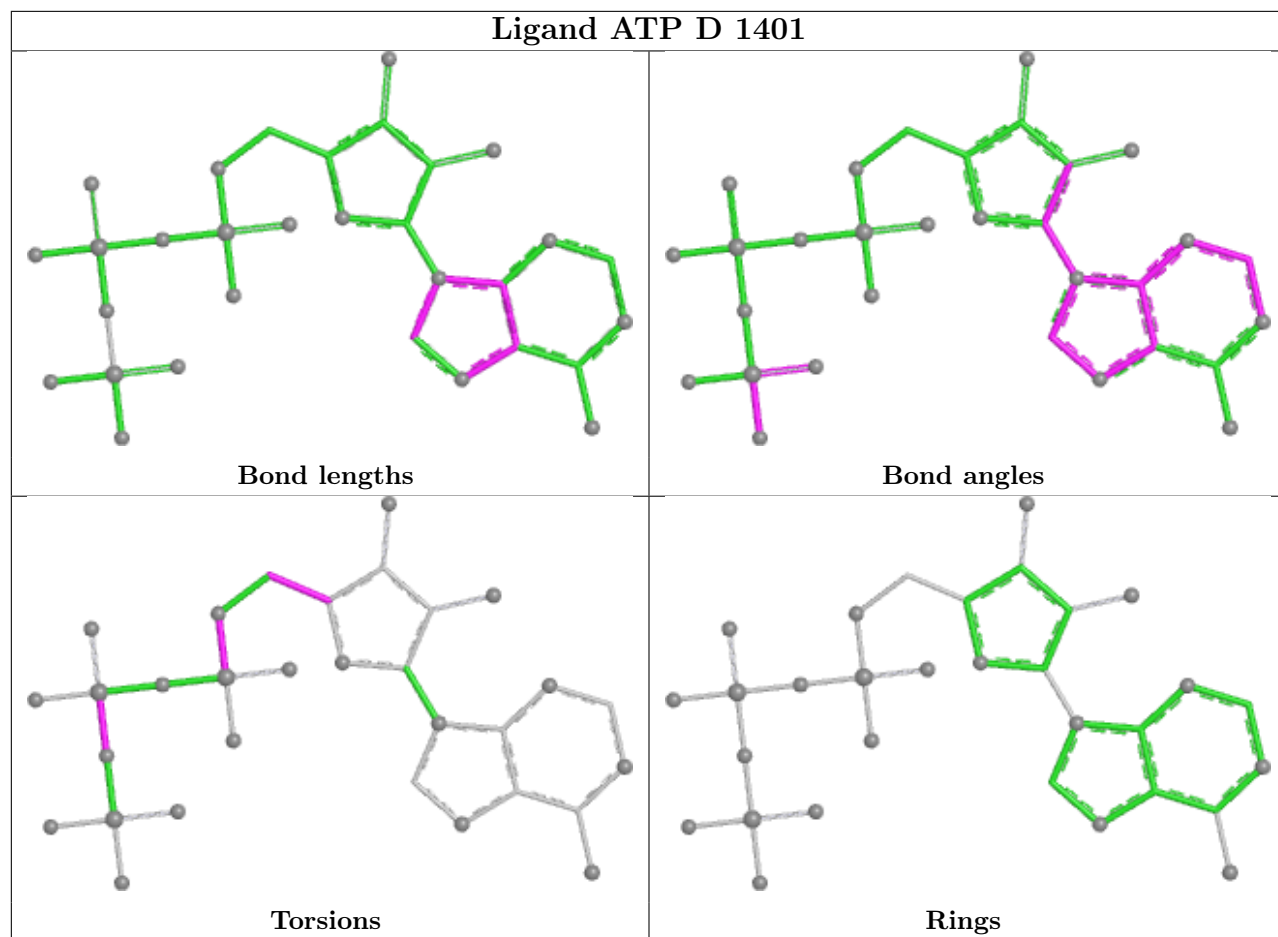
Mol	Chain	Res	Type	Atoms
5	D	1401	ATP	O4'-C4'-C5'-O5'
5	D	1401	ATP	C5'-O5'-PA-O1A
5	C	1401	ATP	PA-O3A-PB-O2B
5	D	1401	ATP	PG-O3B-PB-O2B
5	C	1401	ATP	PA-O3A-PB-O1B

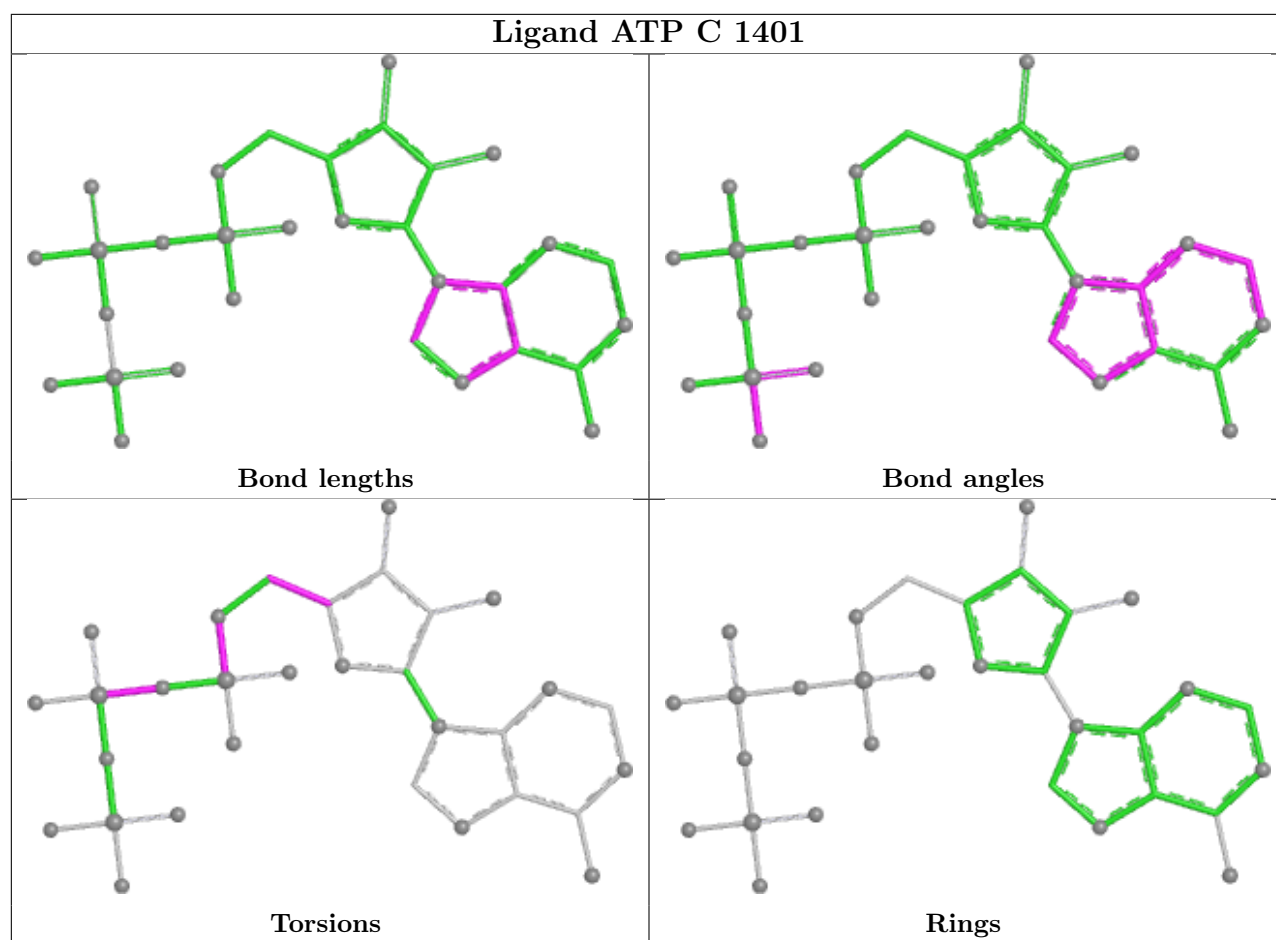
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	1401	ATP	8	0
5	C	1401	ATP	1	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.

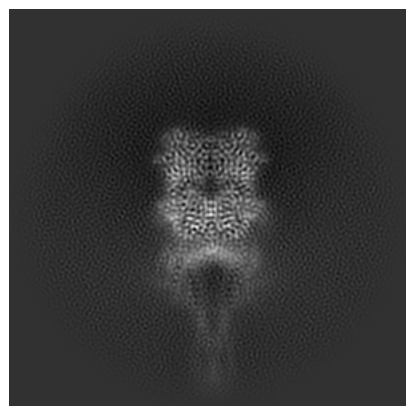
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44558. These allow visual inspection of the internal detail of the map and identification of artifacts.

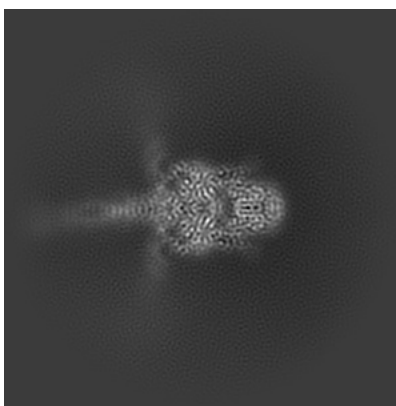
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

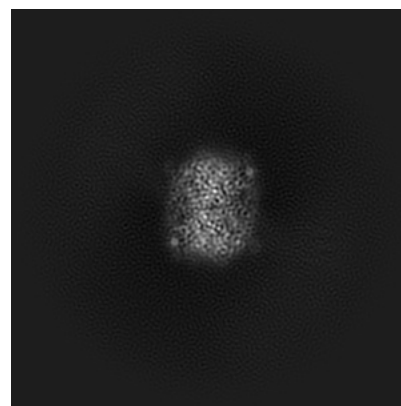
6.1.1 Primary map



X

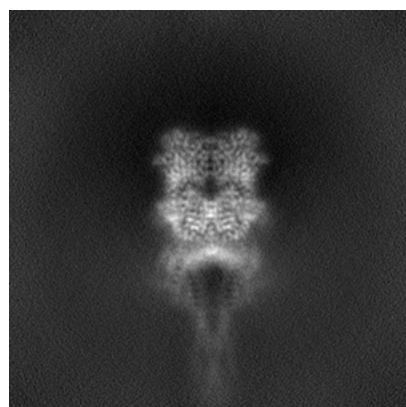


Y

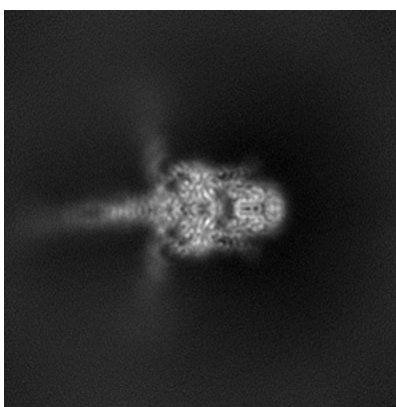


Z

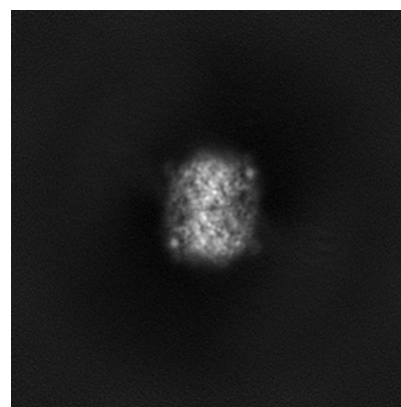
6.1.2 Raw map



X



Y

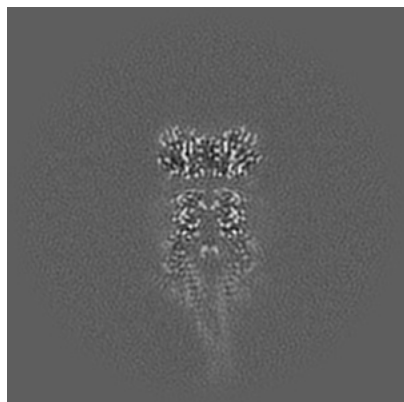


Z

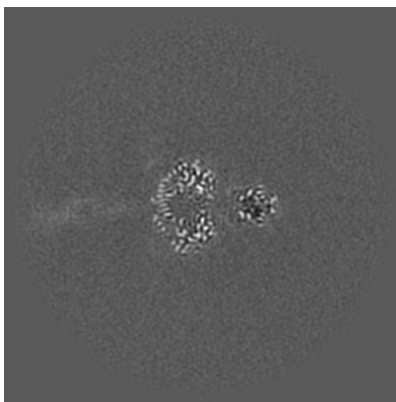
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

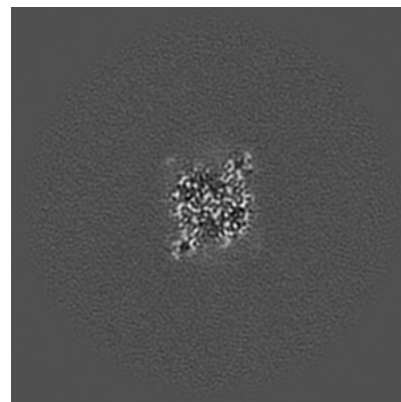
6.2.1 Primary map



X Index: 140

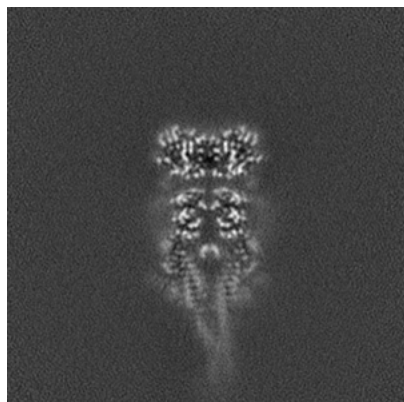


Y Index: 140

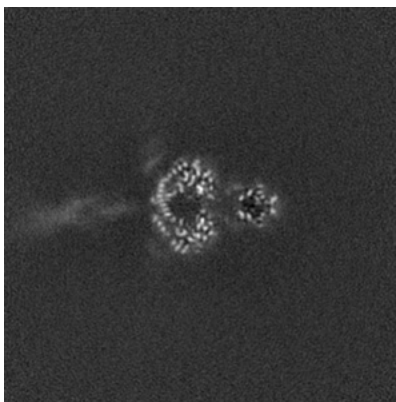


Z Index: 140

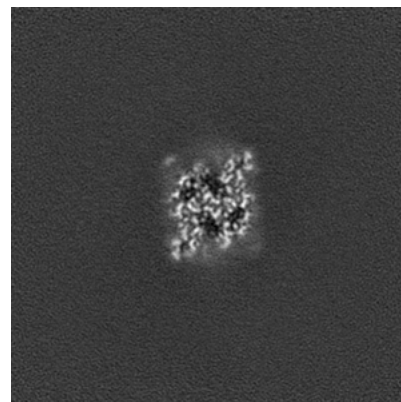
6.2.2 Raw map



X Index: 140



Y Index: 140

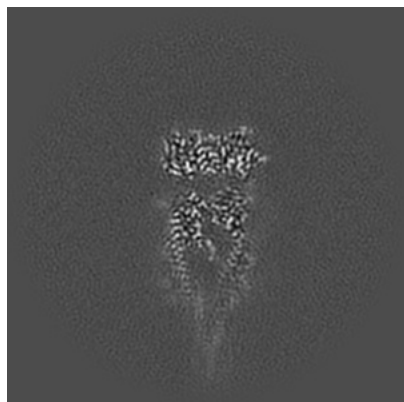


Z Index: 140

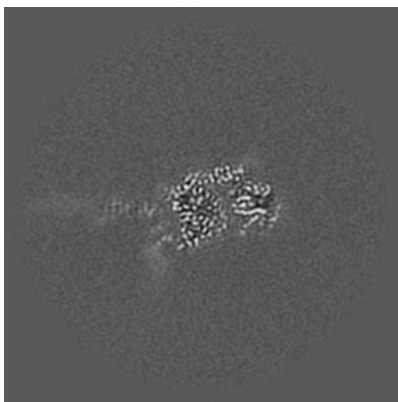
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

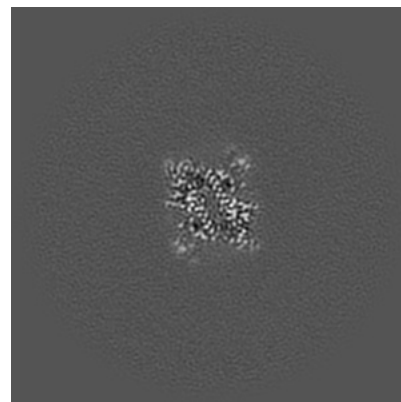
6.3.1 Primary map



X Index: 135

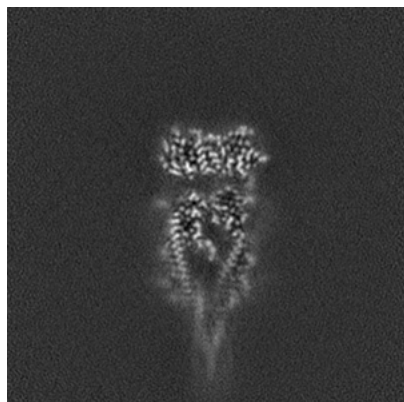


Y Index: 155

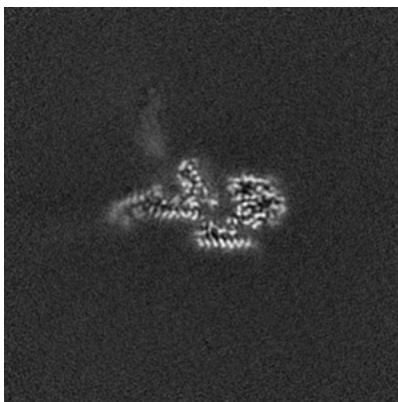


Z Index: 135

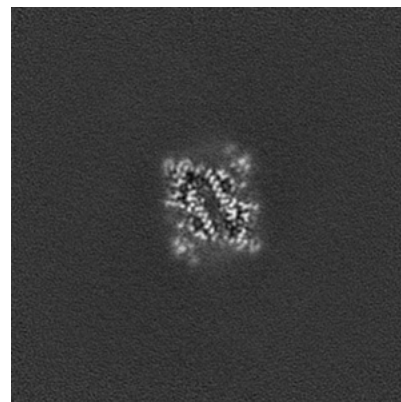
6.3.2 Raw map



X Index: 135



Y Index: 117

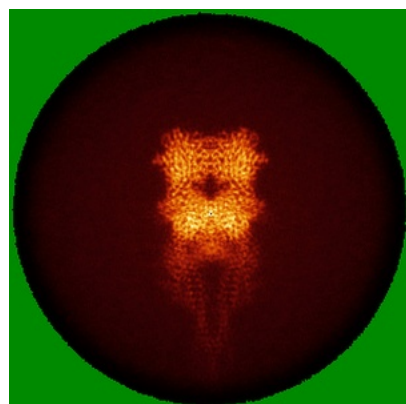


Z Index: 135

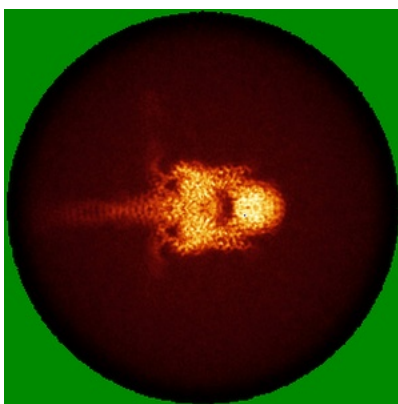
The images above show the largest variance slices of the map in three orthogonal directions.

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

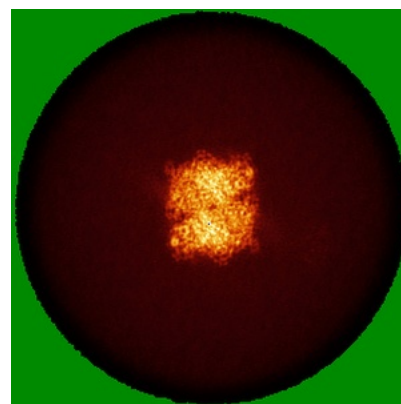
6.4.1 Primary map



X

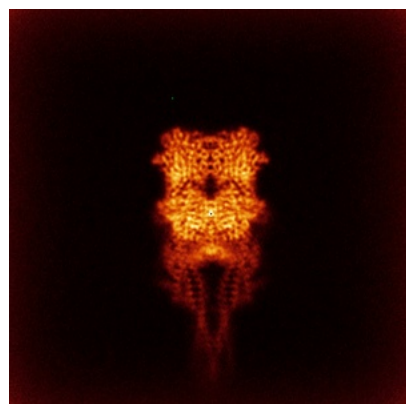


Y

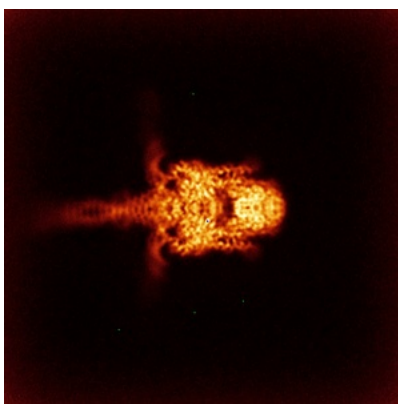


Z

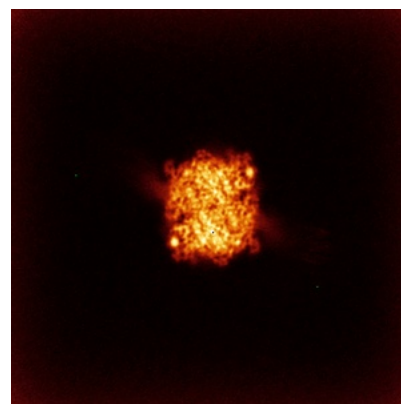
6.4.2 Raw map



X



Y

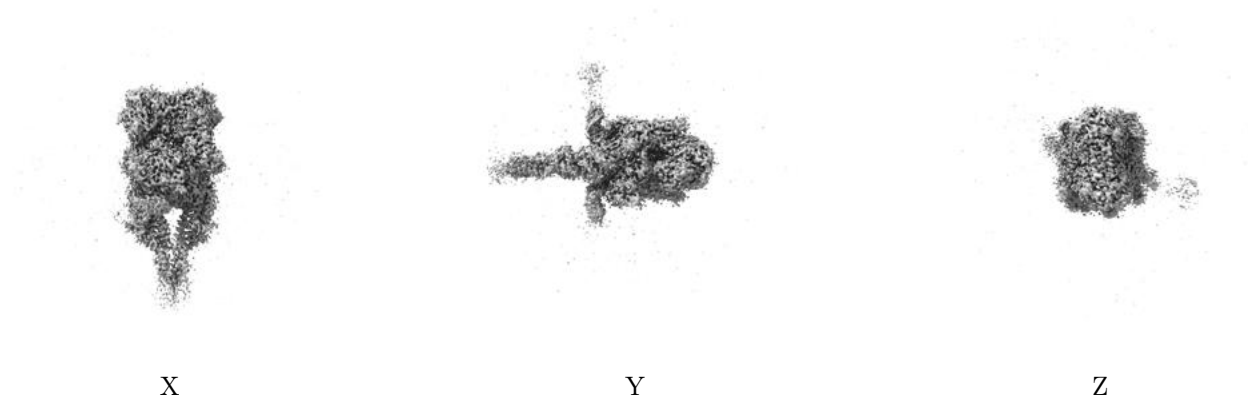


Z

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.

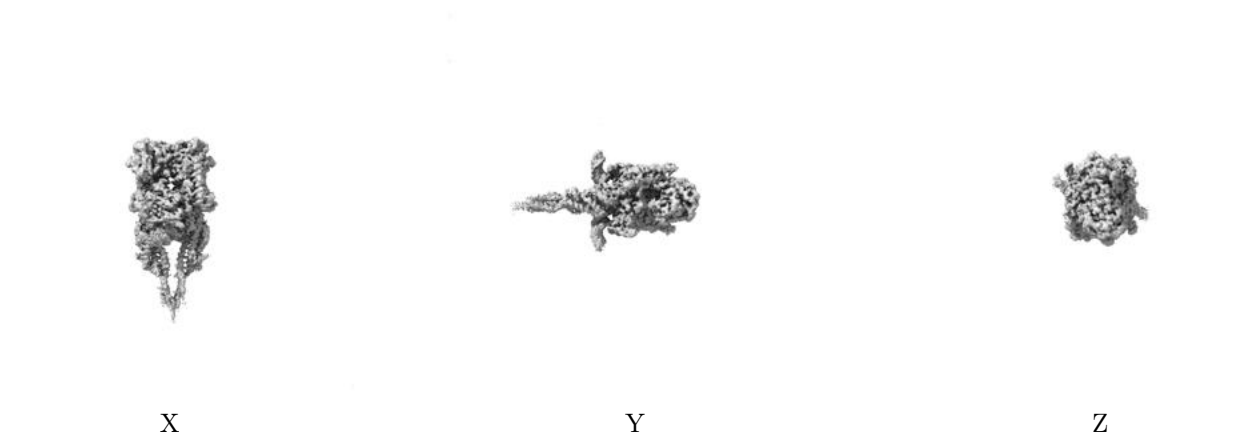
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.28. 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.

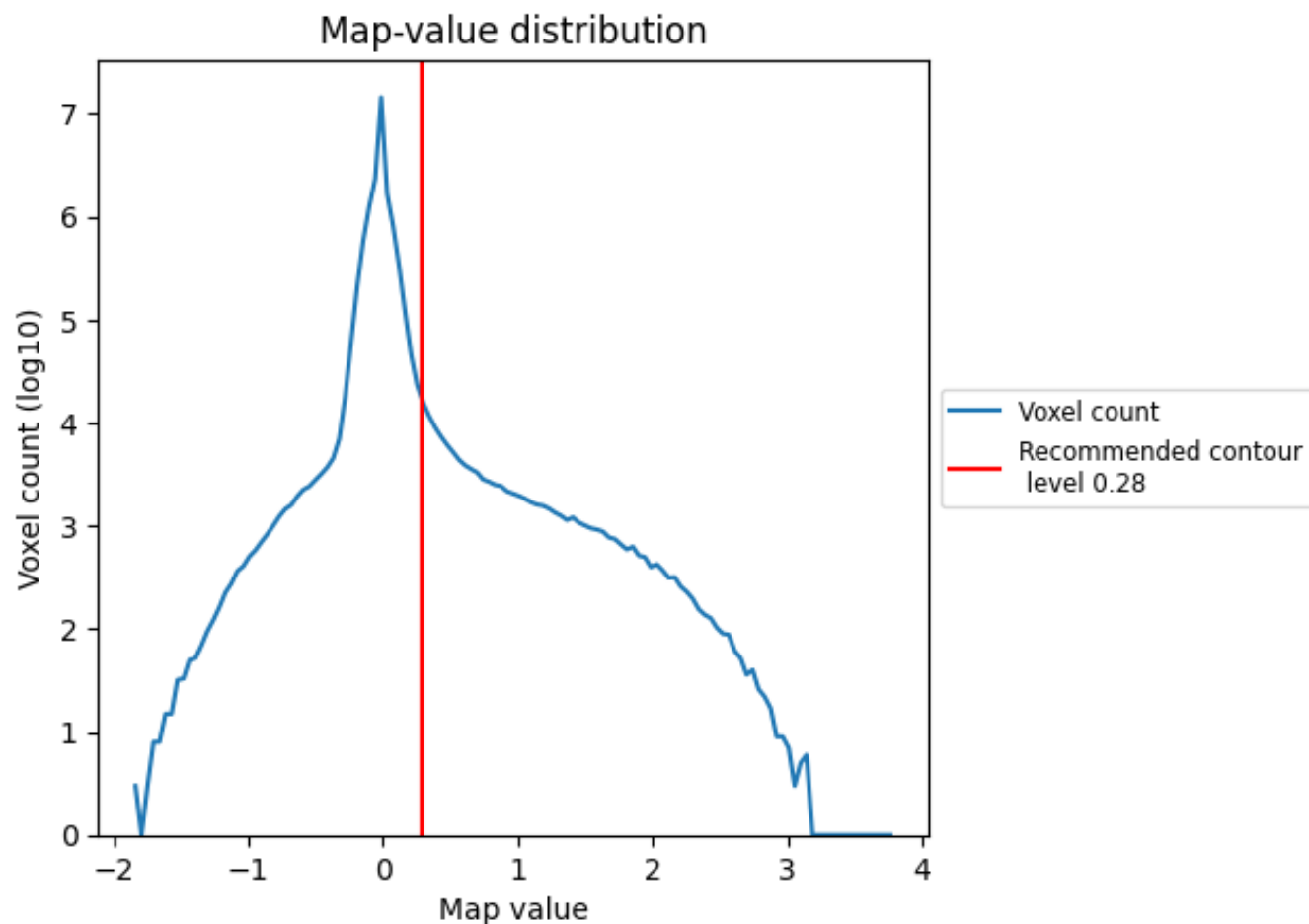
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

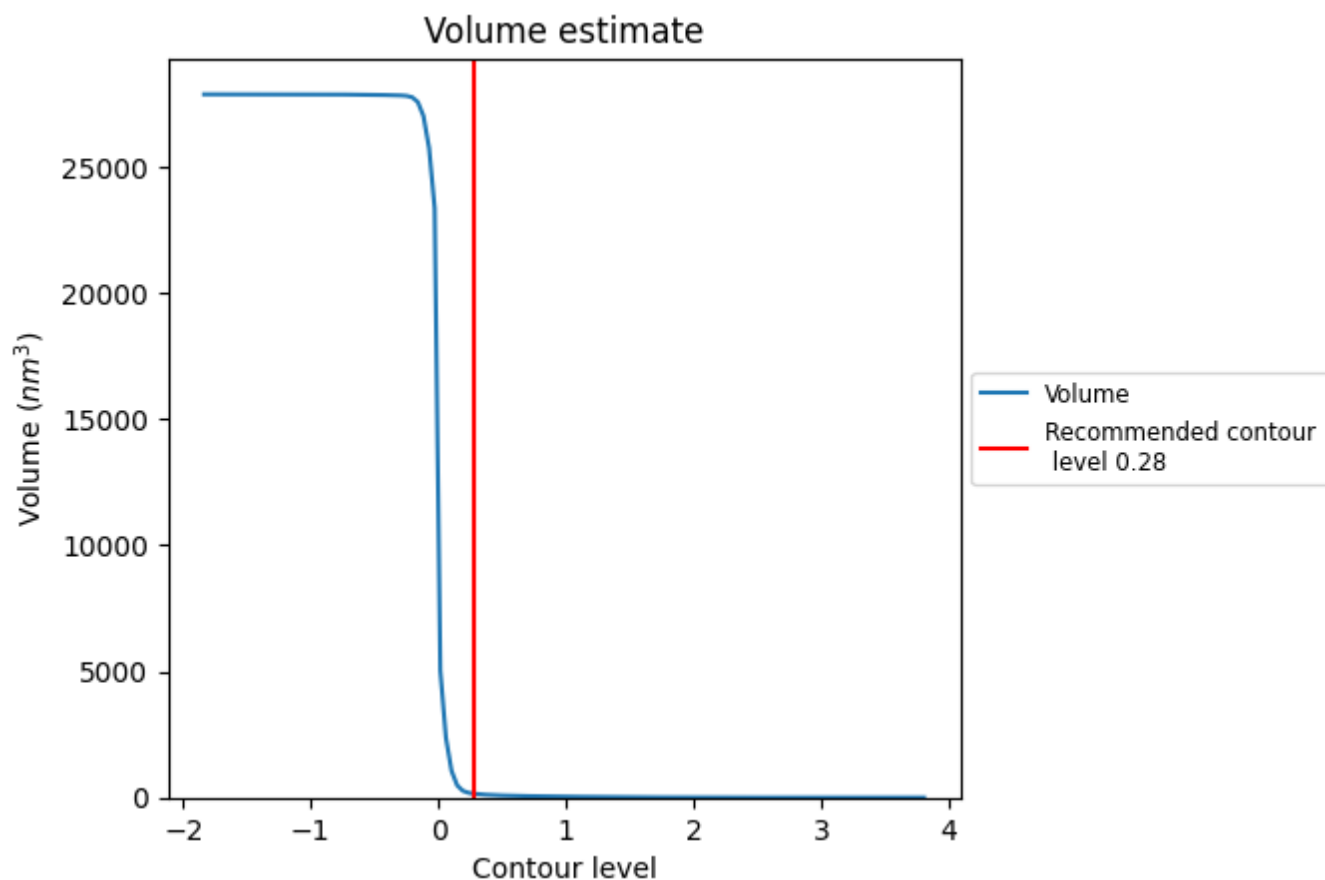
This section contains the results of statistical analysis of the map.

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.

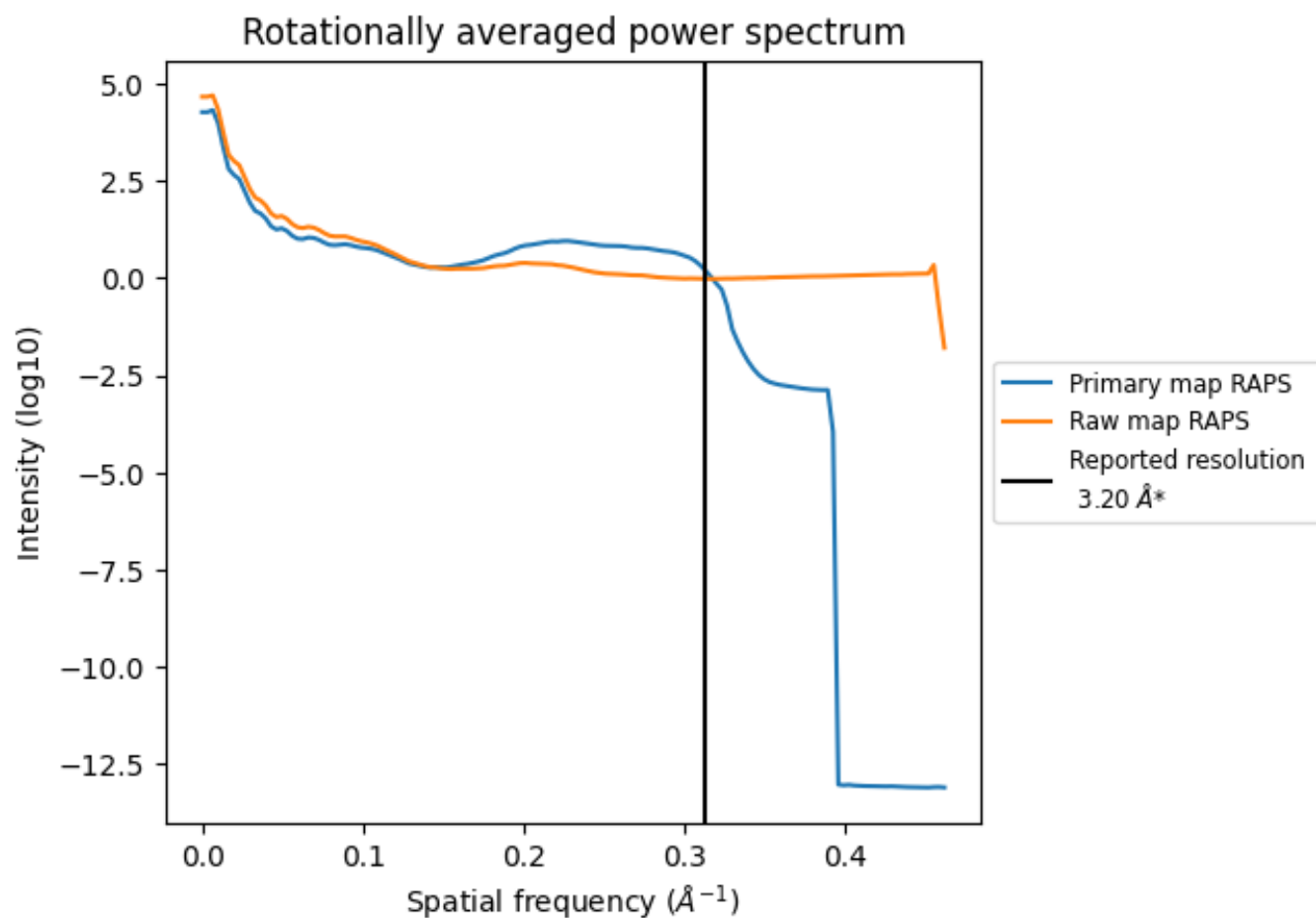
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 152 nm³; this corresponds to an approximate mass of 138 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 ⓘ

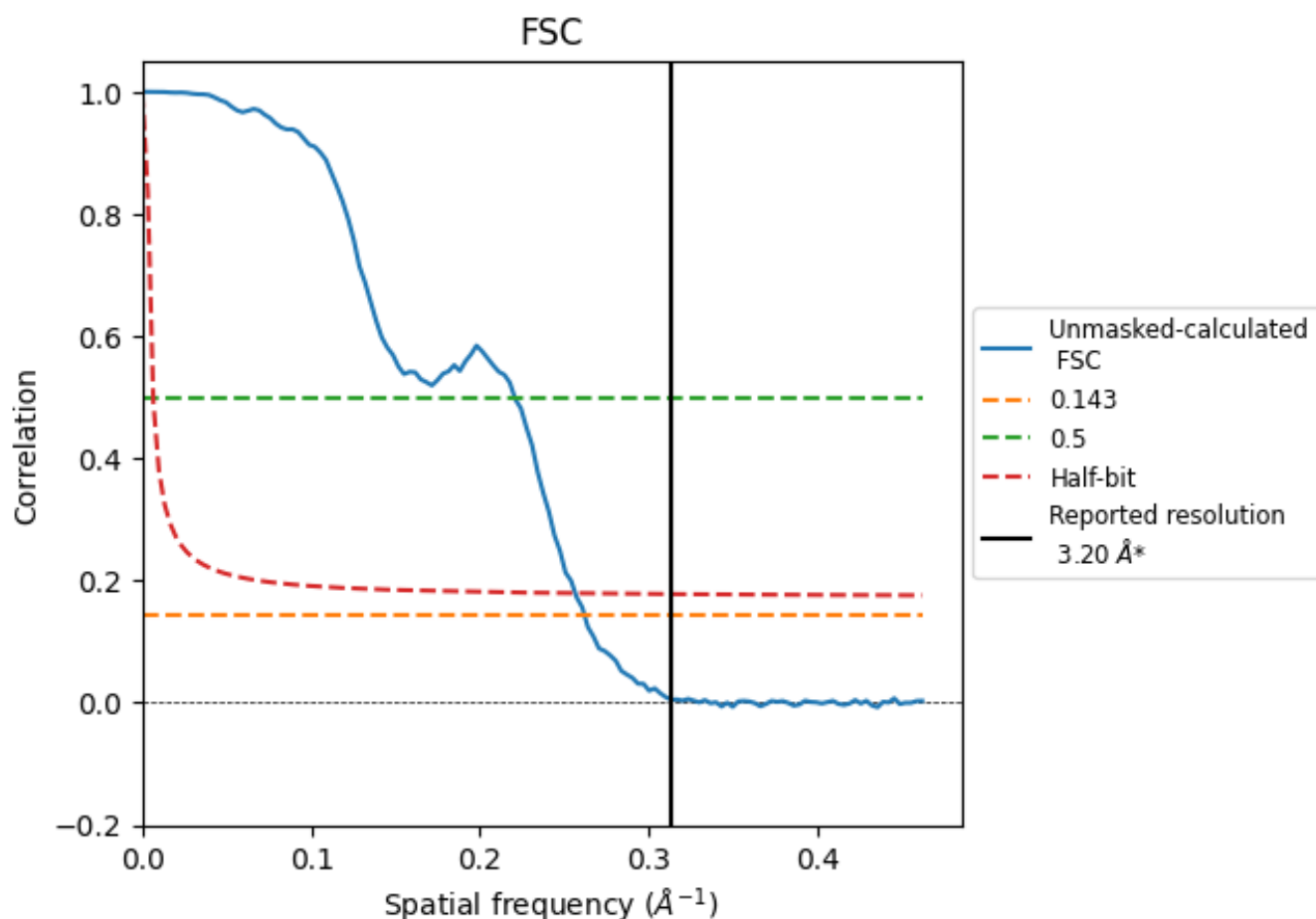


*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

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 $0.312 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

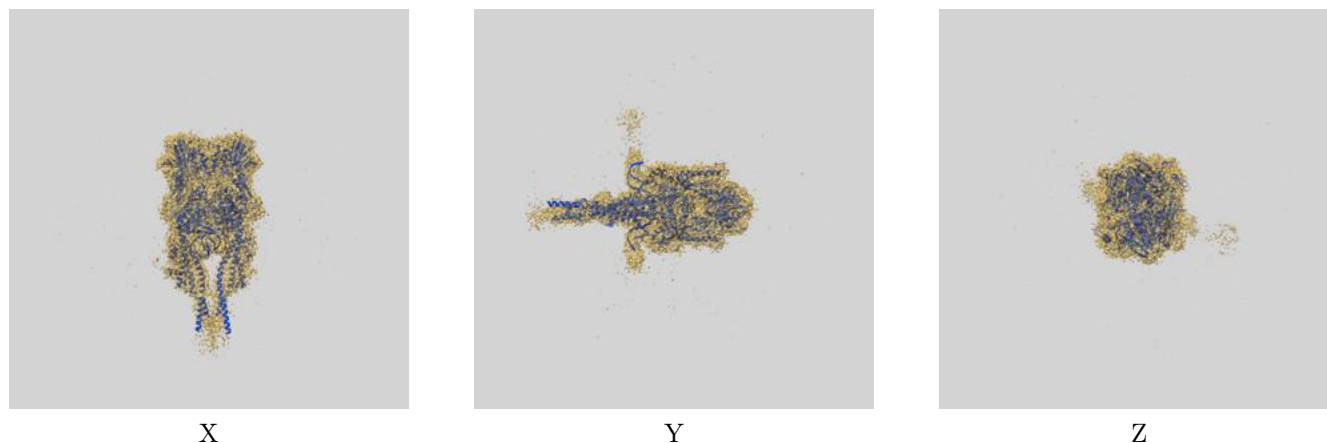
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.82	4.54	3.90

*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 3.82 differs from the reported value 3.2 by more than 10 %

9 Map-model fit [i](#)

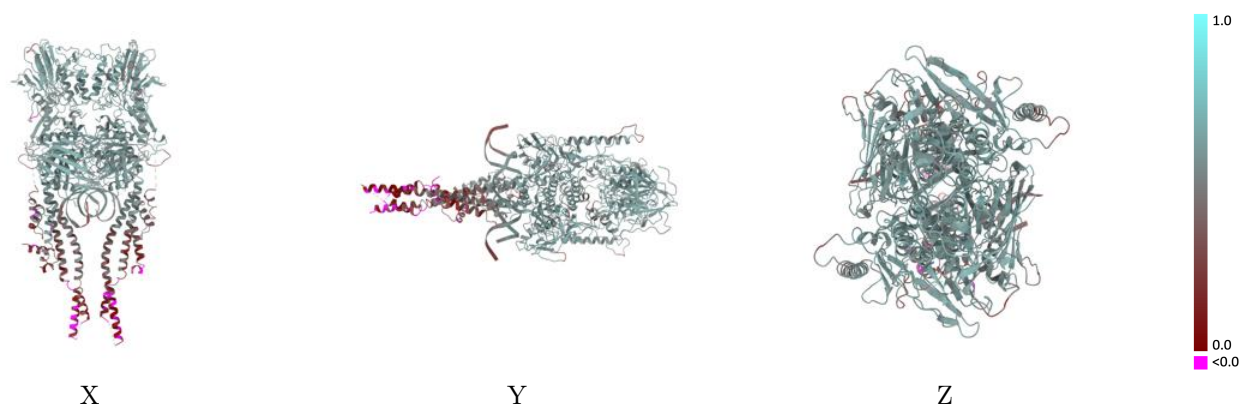
This section contains information regarding the fit between EMDB map EMD-44558 and PDB model 9BI4. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



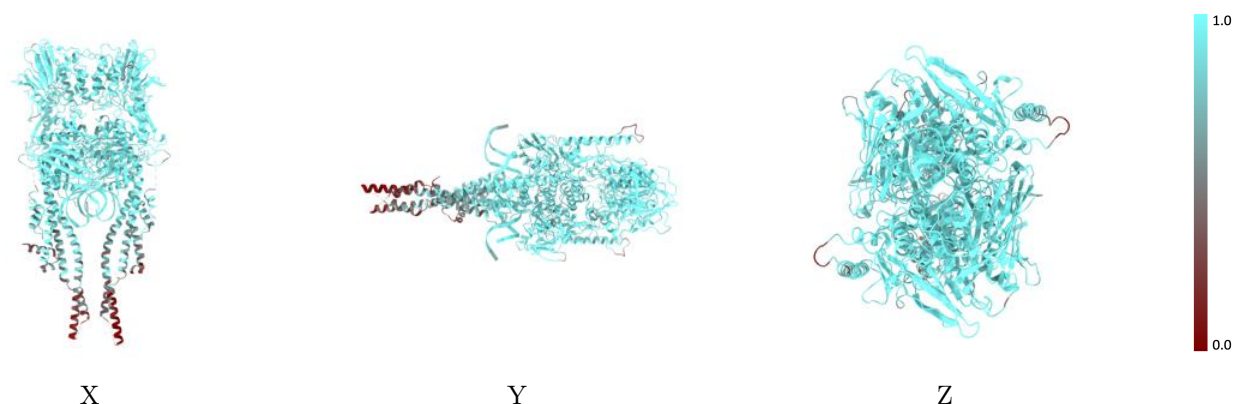
The images above show the 3D surface view of the map at the recommended contour level 0.28 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



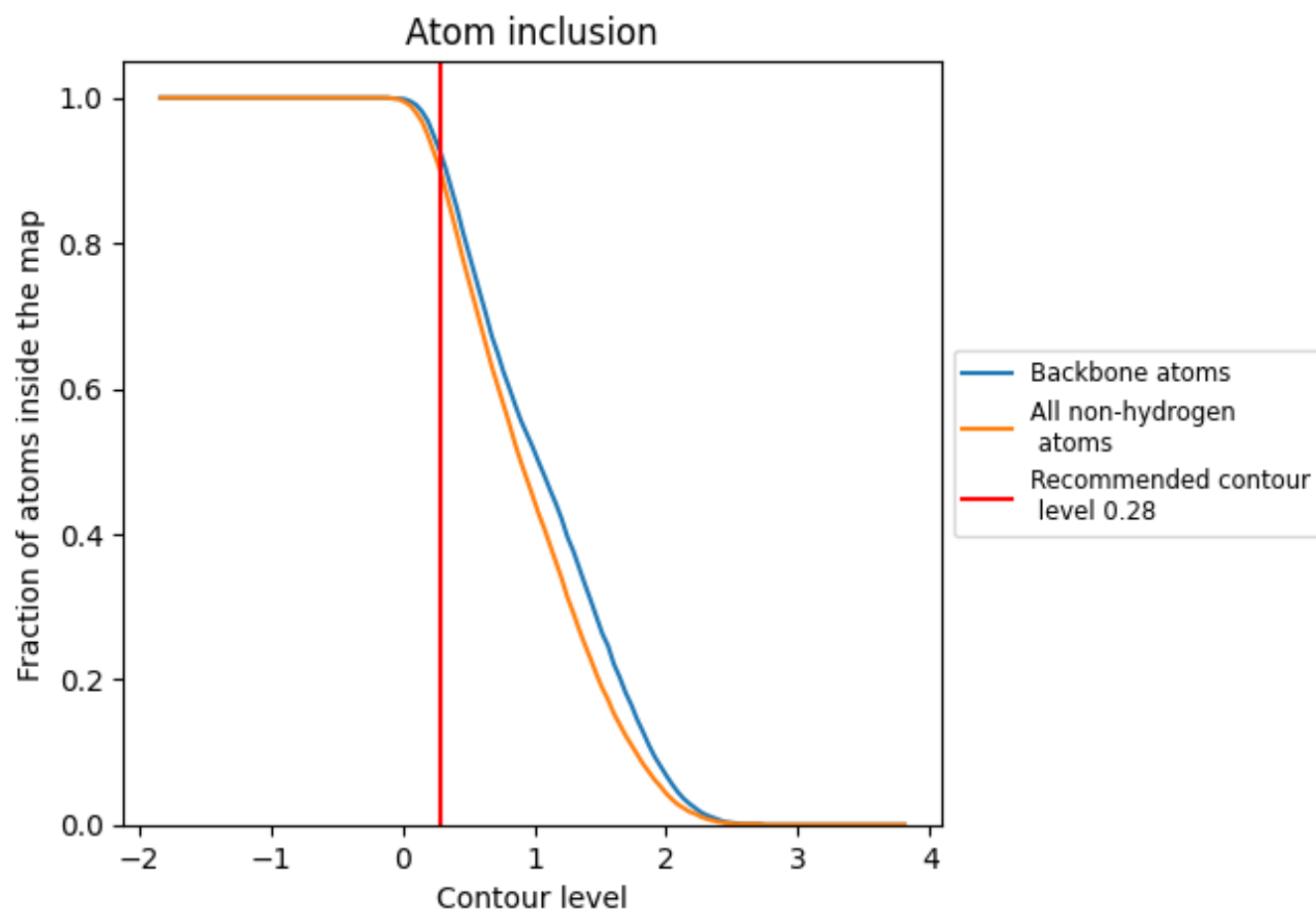
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.28).

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 90% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.28) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9030	0.5210
A	0.9170	0.5280
B	0.9200	0.5360
C	0.8740	0.5080
D	0.8910	0.5150
E	0.9480	0.5170
F	0.9230	0.4930

1.0

0.0

<0.0