



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 01:23 PM UTC

PDB ID : 9I3I / pdb_00009i3i
EMDB ID : EMD-4980
Title : Cryo-EM structure of the MCM-ORC (MO) complex featuring an ORC2 regulatory domain involved in cell cycle regulation of MCM-DH loading for DNA replication.
Authors : Miller, T.C.R.; Lim, C.T.; Diffley, J.F.X.; Costa, A.
Deposited on : 2025-01-23
Resolution : 4.40 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:

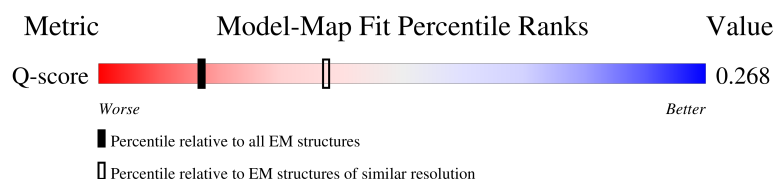
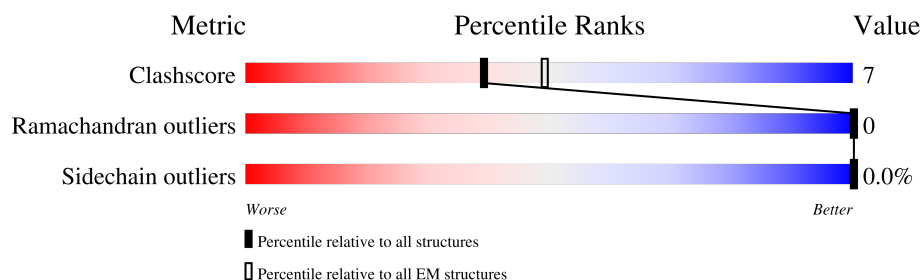
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 4.40 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	3132 (3.91 - 4.90)

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	949	39% 45% 7% 48%
2	B	620	9% 37% 10% 53%
3	C	616	23% 79% 16% 6%
4	D	529	36% 68% 15% 17%

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Mol	Chain	Length	Quality of chain
5	E	479	
6	F	435	
7	2	868	
8	3	1006	
9	4	933	
10	5	775	
11	6	1017	
12	7	845	
13	X	88	
14	Y	88	

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 105645 atoms, of which 52062 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 Origin recognition complex subunit 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	494	Total	C	H	N	O	S	0	0
			8028	2527	4070	676	736	19		

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-34	MET	-	initiating methionine	UNP P54784
A	-33	LYS	-	expression tag	UNP P54784
A	-32	ARG	-	expression tag	UNP P54784
A	-31	ARG	-	expression tag	UNP P54784
A	-30	TRP	-	expression tag	UNP P54784
A	-29	LYS	-	expression tag	UNP P54784
A	-28	LYS	-	expression tag	UNP P54784
A	-27	ASN	-	expression tag	UNP P54784
A	-26	PHE	-	expression tag	UNP P54784
A	-25	ILE	-	expression tag	UNP P54784
A	-24	ALA	-	expression tag	UNP P54784
A	-23	VAL	-	expression tag	UNP P54784
A	-22	SER	-	expression tag	UNP P54784
A	-21	ALA	-	expression tag	UNP P54784
A	-20	ALA	-	expression tag	UNP P54784
A	-19	ASN	-	expression tag	UNP P54784
A	-18	ARG	-	expression tag	UNP P54784
A	-17	PHE	-	expression tag	UNP P54784
A	-16	LYS	-	expression tag	UNP P54784
A	-15	LYS	-	expression tag	UNP P54784
A	-14	ILE	-	expression tag	UNP P54784
A	-13	SER	-	expression tag	UNP P54784
A	-12	SER	-	expression tag	UNP P54784
A	-11	SER	-	expression tag	UNP P54784
A	-10	GLY	-	expression tag	UNP P54784
A	-9	ALA	-	expression tag	UNP P54784
A	-8	LEU	-	expression tag	UNP P54784
A	-7	GLU	-	expression tag	UNP P54784

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	ASN	-	expression tag	UNP P54784
A	-5	LEU	-	expression tag	UNP P54784
A	-4	TYR	-	expression tag	UNP P54784
A	-3	PHE	-	expression tag	UNP P54784
A	-2	GLN	-	expression tag	UNP P54784
A	-1	GLY	-	expression tag	UNP P54784
A	0	GLU	-	expression tag	UNP P54784

- Molecule 2 is a protein called Origin recognition complex subunit 2.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	289	Total	C	H	N	O	S	0	0
			4405	1546	2018	397	435	9		

- Molecule 3 is a protein called Origin recognition complex subunit 3.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	582	Total	C	H	N	O	S	0	0
			9579	3102	4766	794	901	16		

- Molecule 4 is a protein called Origin recognition complex subunit 4.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	D	437	Total	C	H	N	O	S	0	0
			7193	2285	3627	605	663	13		

- Molecule 5 is a protein called Origin recognition complex subunit 5.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	E	459	Total	C	H	N	O	S	0	0
			7529	2431	3780	609	695	14		

- Molecule 6 is a protein called Origin recognition complex subunit 6.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	F	242	Total	C	H	N	O	S	0	0
			4046	1274	2051	343	358	20		

- Molecule 7 is a protein called DNA replication licensing factor MCM2.

Mol	Chain	Residues	Atoms						AltConf	Trace
7	2	601	Total	C	H	N	O	S	0	0
			9573	2998	4820	845	891	19		

- Molecule 8 is a protein called DNA replication licensing factor MCM3.

Mol	Chain	Residues	Atoms						AltConf	Trace
8	3	599	Total	C	H	N	O	S	0	0
			9448	2956	4752	840	887	13		

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	-34	MET	-	initiating methionine	UNP P24279
3	-33	LYS	-	expression tag	UNP P24279
3	-32	ARG	-	expression tag	UNP P24279
3	-31	ARG	-	expression tag	UNP P24279
3	-30	TRP	-	expression tag	UNP P24279
3	-29	LYS	-	expression tag	UNP P24279
3	-28	LYS	-	expression tag	UNP P24279
3	-27	ASN	-	expression tag	UNP P24279
3	-26	PHE	-	expression tag	UNP P24279
3	-25	ILE	-	expression tag	UNP P24279
3	-24	ALA	-	expression tag	UNP P24279
3	-23	VAL	-	expression tag	UNP P24279
3	-22	SER	-	expression tag	UNP P24279
3	-21	ALA	-	expression tag	UNP P24279
3	-20	ALA	-	expression tag	UNP P24279
3	-19	ASN	-	expression tag	UNP P24279
3	-18	ARG	-	expression tag	UNP P24279
3	-17	PHE	-	expression tag	UNP P24279
3	-16	LYS	-	expression tag	UNP P24279
3	-15	LYS	-	expression tag	UNP P24279
3	-14	ILE	-	expression tag	UNP P24279
3	-13	SER	-	expression tag	UNP P24279
3	-12	SER	-	expression tag	UNP P24279
3	-11	SER	-	expression tag	UNP P24279
3	-10	GLY	-	expression tag	UNP P24279
3	-9	ALA	-	expression tag	UNP P24279
3	-8	LEU	-	expression tag	UNP P24279
3	-7	GLU	-	expression tag	UNP P24279
3	-6	ASN	-	expression tag	UNP P24279
3	-5	LEU	-	expression tag	UNP P24279
3	-4	TYR	-	expression tag	UNP P24279

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Chain	Residue	Modelled	Actual	Comment	Reference
3	-3	PHE	-	expression tag	UNP P24279
3	-2	GLN	-	expression tag	UNP P24279
3	-1	GLY	-	expression tag	UNP P24279
3	0	GLU	-	expression tag	UNP P24279

- Molecule 9 is a protein called DNA replication licensing factor MCM4.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	4	641	Total	C	H	N	O	S	0	0
			10282	3210	5172	883	989	28		

- Molecule 10 is a protein called Minichromosome maintenance protein 5.

Mol	Chain	Residues	Atoms						AltConf	Trace
10	5	602	Total	C	H	N	O	S	0	0
			9532	2976	4802	813	917	24		

- Molecule 11 is a protein called DNA replication licensing factor MCM6.

Mol	Chain	Residues	Atoms						AltConf	Trace
11	6	616	Total	C	H	N	O	S	0	0
			9782	3072	4908	853	924	25		

- Molecule 12 is a protein called DNA replication licensing factor MCM7.

Mol	Chain	Residues	Atoms						AltConf	Trace
12	7	654	Total	C	H	N	O	S	0	0
			10353	3244	5215	886	979	29		

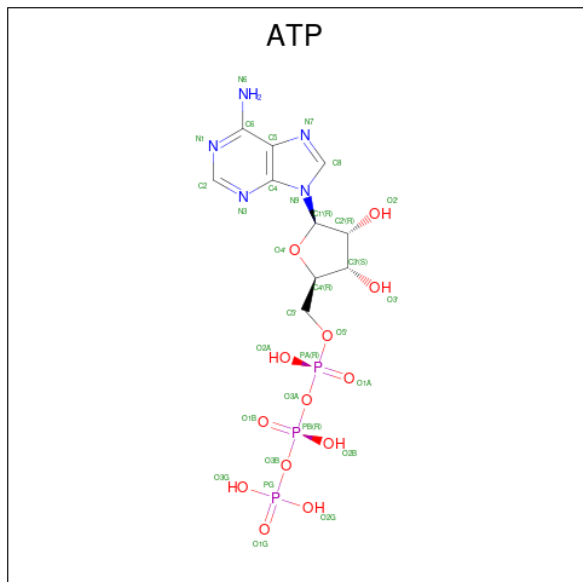
- Molecule 13 is a DNA chain called DNA (88-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
13	X	88	Total	C	H	N	O	P	0	0
			2806	867	1010	294	548	87		

- Molecule 14 is a DNA chain called DNA (88-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
14	Y	88	Total	C	H	N	O	P	0	0
			2796	863	987	352	506	88		

- Molecule 15 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).

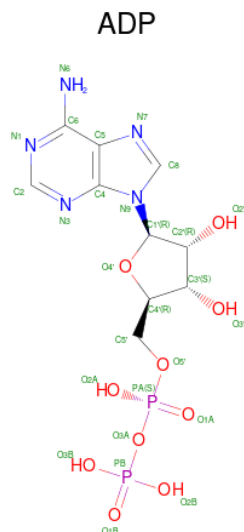


Mol	Chain	Residues	Atoms						AltConf
15	A	1	Total	C	H	N	O	P	0
			43	10	12	5	13	3	
15	D	1	Total	C	H	N	O	P	0
			43	10	12	5	13	3	
15	E	1	Total	C	H	N	O	P	0
			43	10	12	5	13	3	

- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
16	A	1	Total	Mg	0
			1	1	
16	D	1	Total	Mg	0
			1	1	
16	E	1	Total	Mg	0
			1	1	

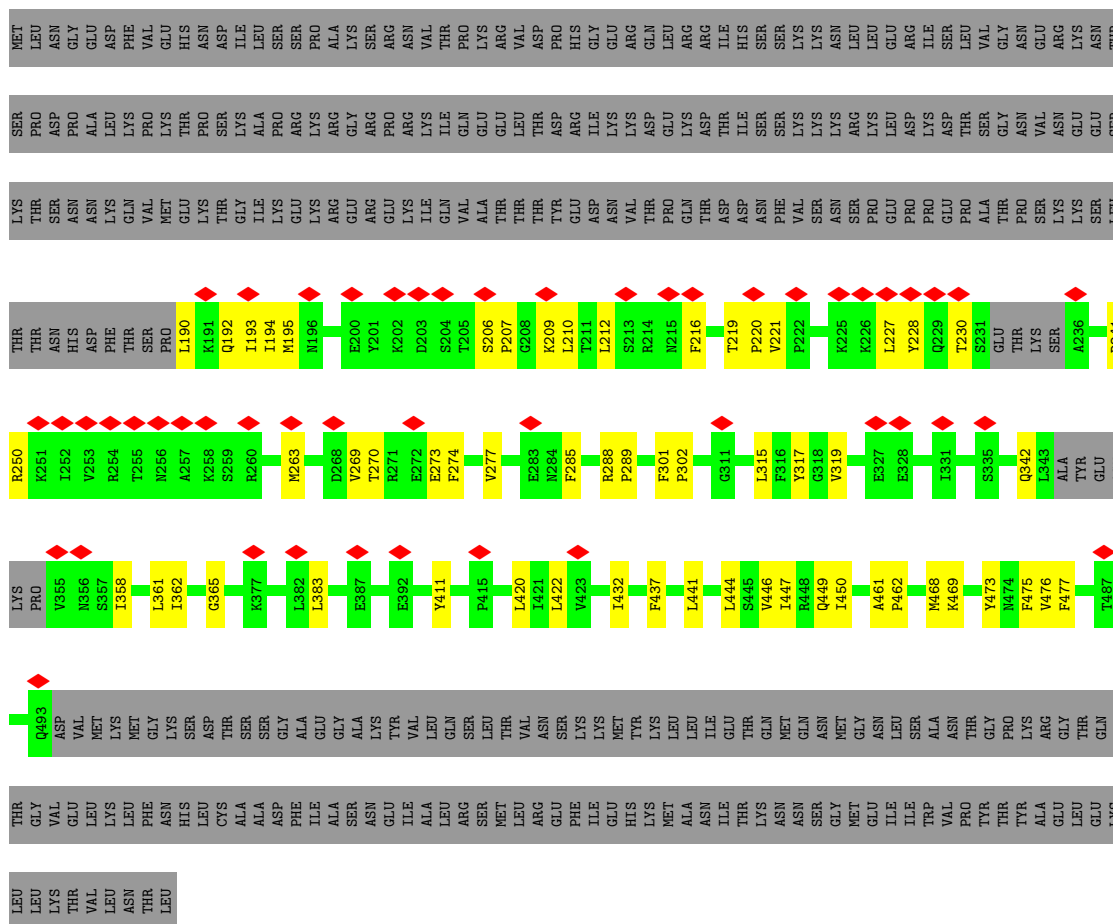
- Molecule 17 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf	
17	2	1	Total 39	C 10	H 12	N 5	O 10	P 2	0
17	3	1	Total 39	C 10	H 12	N 5	O 10	P 2	0
17	5	1	Total 39	C 10	H 12	N 5	O 10	P 2	0
17	7	1	Total 39	C 10	H 12	N 5	O 10	P 2	0

- Molecule 18 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

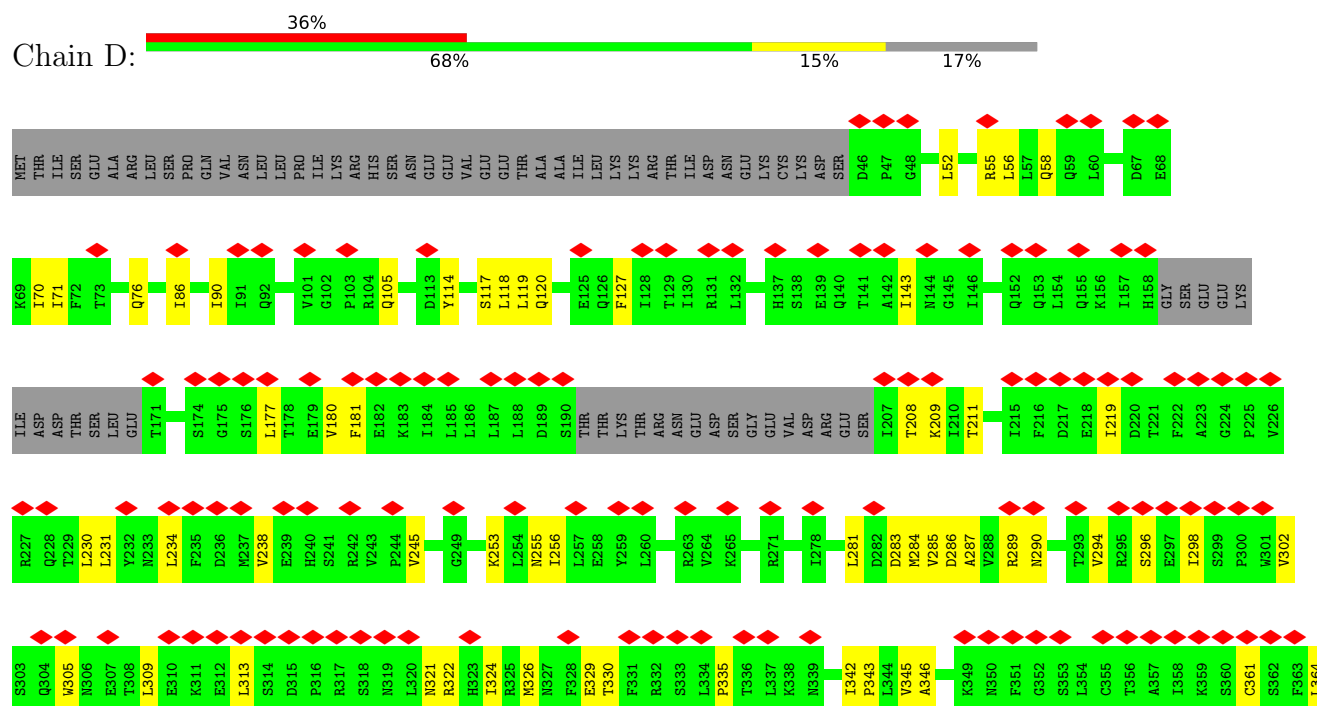
Mol	Chain	Residues	Atoms	AltConf
18	2	1	Total Zn 1 1	0
18	4	1	Total Zn 1 1	0
18	5	1	Total Zn 1 1	0
18	6	1	Total Zn 1 1	0
18	7	1	Total Zn 1 1	0

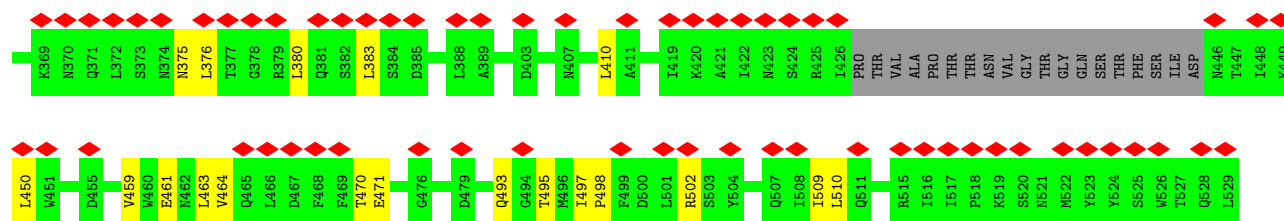


- Molecule 3: Origin recognition complex subunit 3

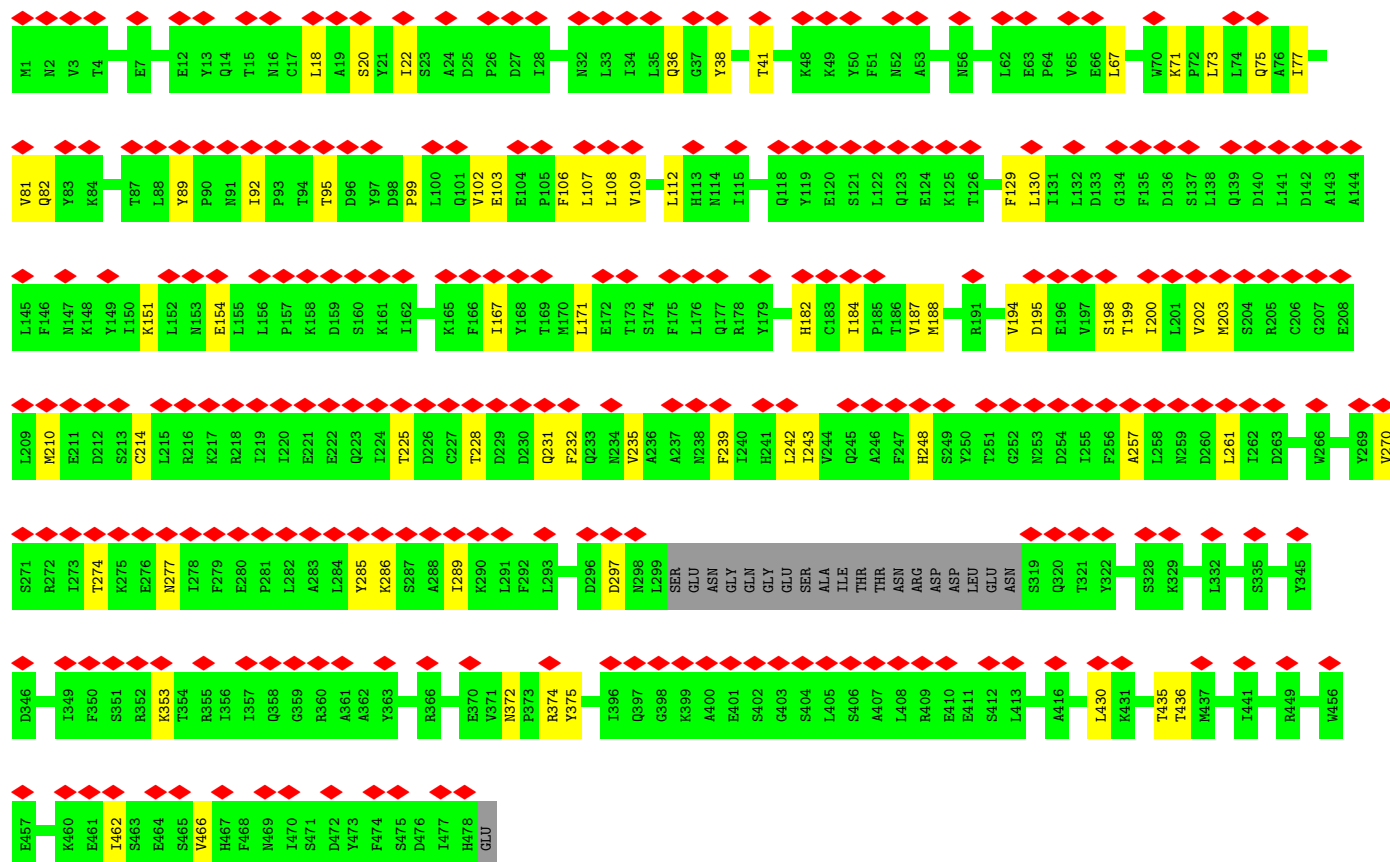
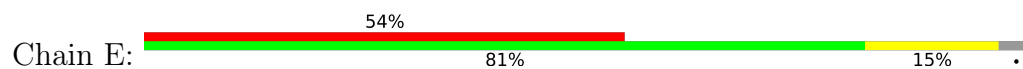


• Molecule 4: Origin recognition complex subunit 4

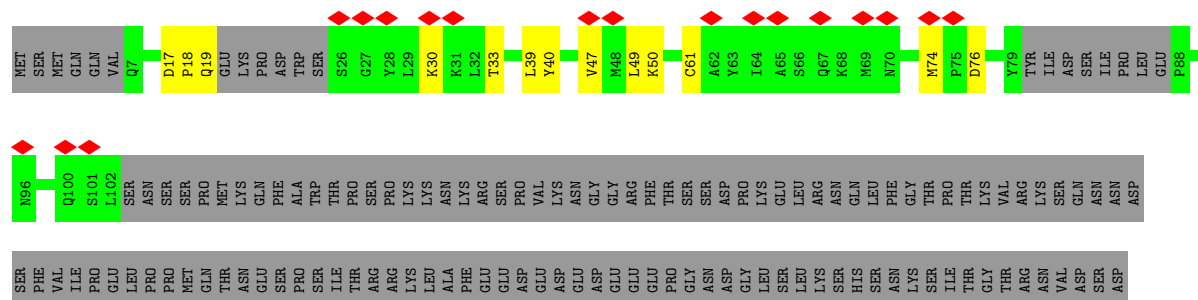
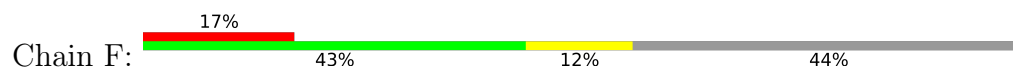


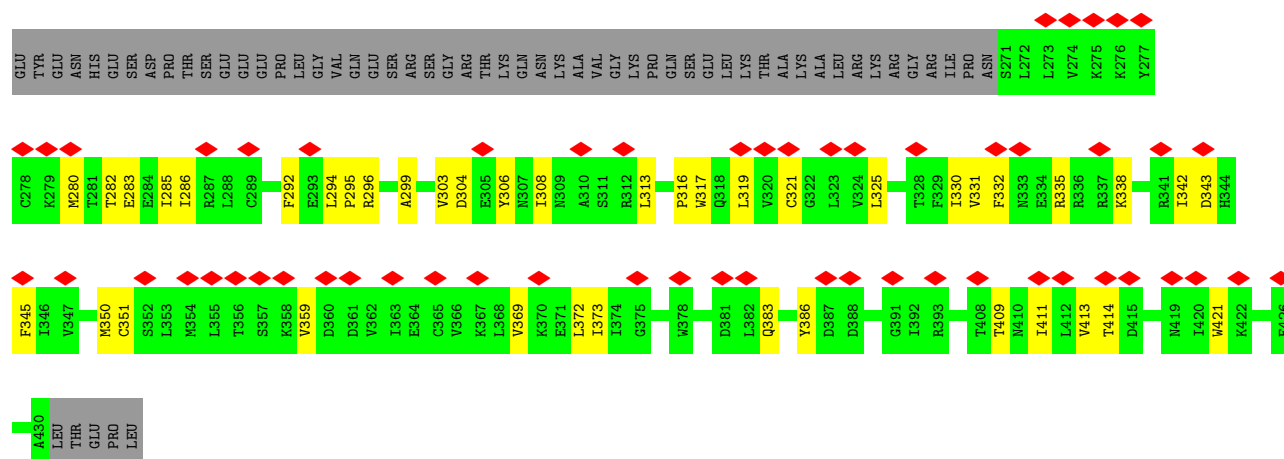


• Molecule 5: Origin recognition complex subunit 5

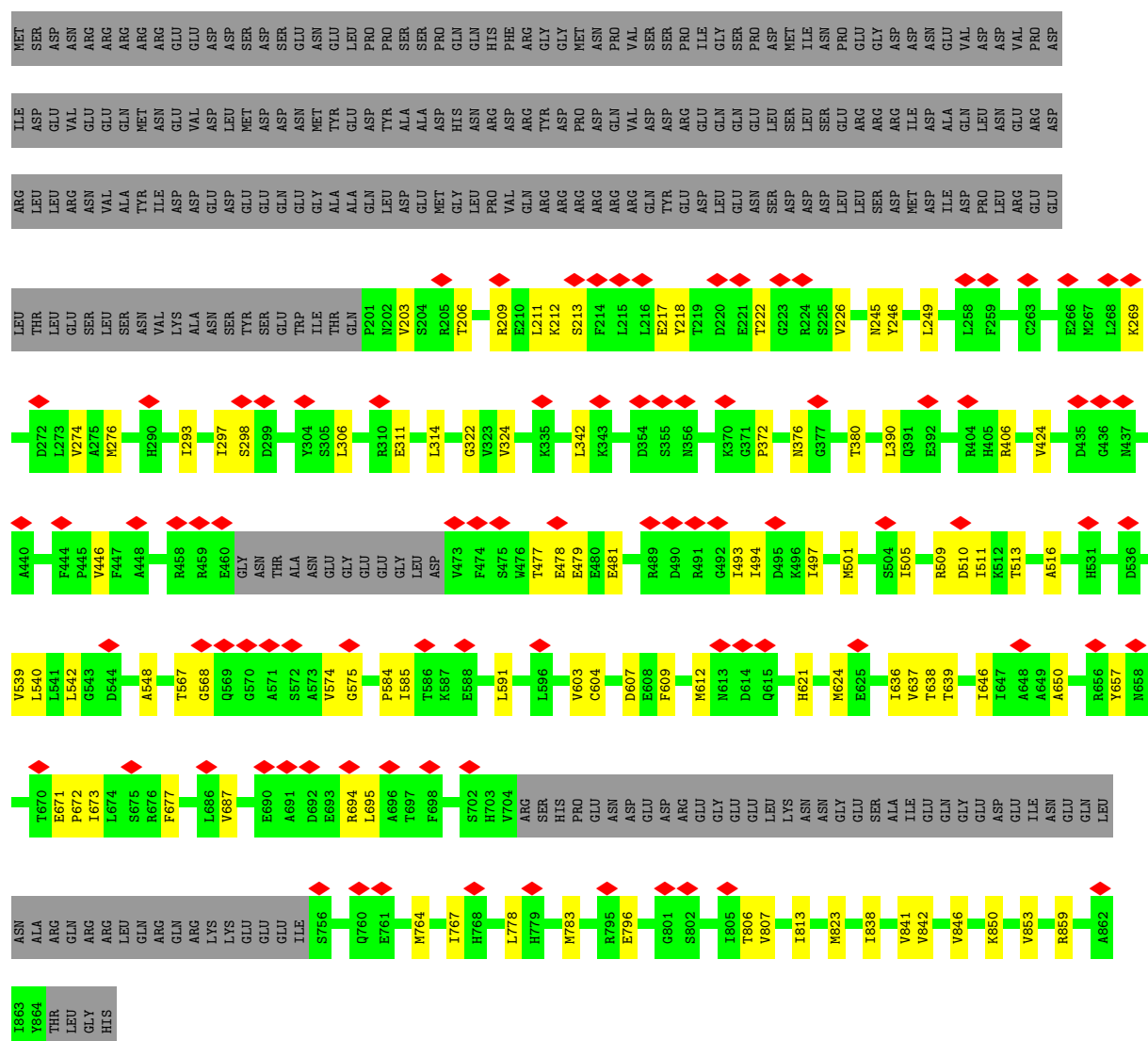


• Molecule 6: Origin recognition complex subunit 6

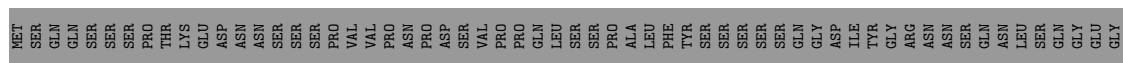




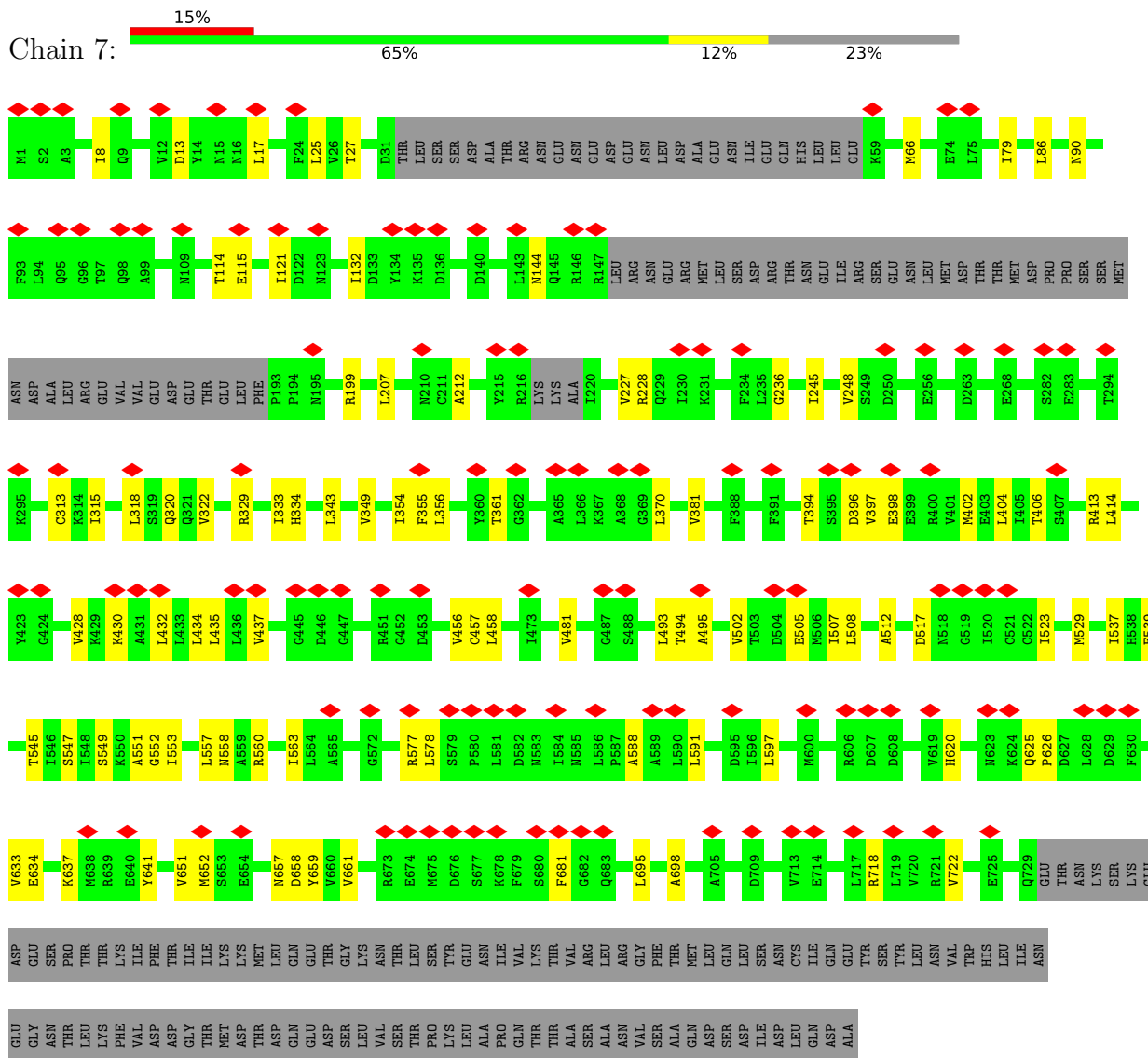
• Molecule 7: DNA replication licensing factor MCM2



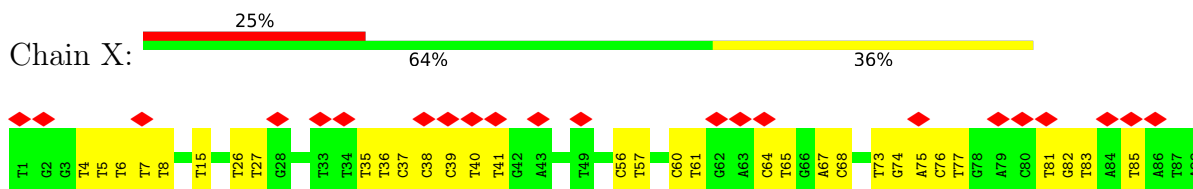
Chain 3:



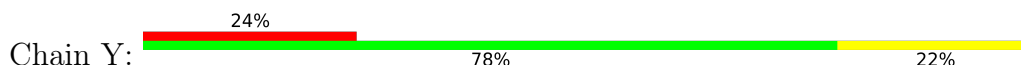
- Molecule 12: DNA replication licensing factor MCM7

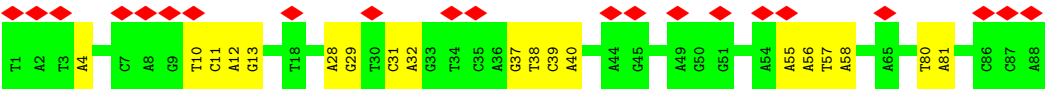


- Molecule 13: DNA (88-MER)



- Molecule 14: DNA (88-MER)





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	177687	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50.4	Depositor
Minimum defocus (nm)	2700	Depositor
Maximum defocus (nm)	4200	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.070	Depositor
Minimum map value	-0.037	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.016	Depositor
Map size (Å)	469.2, 469.2, 469.2	wwPDB
Map dimensions	340, 340, 340	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.38, 1.38, 1.38	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: ZN, MG, ATP, ADP

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	0/4020	0.33	0/5407
2	B	0.15	0/2446	0.38	0/3307
3	C	0.13	0/4918	0.35	0/6641
4	D	0.12	0/3628	0.35	0/4901
5	E	0.13	0/3835	0.34	0/5205
6	F	0.15	0/2026	0.39	0/2722
7	2	0.14	0/4835	0.36	0/6529
8	3	0.14	0/4775	0.37	0/6473
9	4	0.14	0/5185	0.36	0/7009
10	5	0.15	0/4796	0.37	0/6482
11	6	0.14	0/4954	0.37	0/6683
12	7	0.13	0/5219	0.34	0/7053
13	X	0.21	0/2006	0.49	0/3097
14	Y	0.21	0/2037	0.45	0/3139
All	All	0.14	0/54680	0.37	0/74648

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	3958	4070	4071	53	0
2	B	2387	2018	2379	67	0
3	C	4813	4766	4765	86	0
4	D	3566	3627	3628	57	0
5	E	3749	3780	3781	49	0
6	F	1995	2051	2051	50	0
7	2	4753	4820	4820	67	0
8	3	4696	4752	4752	66	0
9	4	5110	5172	5170	78	0
10	5	4730	4802	4800	84	0
11	6	4874	4908	4907	60	0
12	7	5138	5215	5213	76	0
13	X	1796	1010	1010	30	0
14	Y	1809	987	987	12	0
15	A	31	12	12	0	0
15	D	31	12	12	0	0
15	E	31	12	12	0	0
16	A	1	0	0	0	0
16	D	1	0	0	0	0
16	E	1	0	0	0	0
17	2	27	12	12	0	0
17	3	27	12	12	0	0
17	5	27	12	12	0	0
17	7	27	12	12	0	0
18	2	1	0	0	0	0
18	4	1	0	0	0	0
18	5	1	0	0	0	0
18	6	1	0	0	0	0
18	7	1	0	0	0	0
All	All	53583	52062	52418	763	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 (763) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:3:176:LEU:HD11	10:5:250:PHE:HB2	1.63	0.80
10:5:282:LEU:HD11	10:5:333:ILE:HG22	1.65	0.79
12:7:315:ILE:HD11	12:7:333:ILE:HD11	1.64	0.78
10:5:497:MET:HE2	10:5:519:VAL:HG11	1.67	0.76
12:7:434:LEU:HD23	12:7:695:LEU:HD22	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:5:339:THR:HG22	10:5:340:SER:H	1.52	0.75
2:B:221:VAL:HG22	6:F:283:GLU:CD	2.15	0.72
5:E:194:VAL:HG12	5:E:248:HIS:CE1	2.25	0.72
11:6:362:GLN:HA	11:6:376:THR:HG22	1.71	0.72
6:F:74:MET:N	6:F:74:MET:HE2	2.05	0.71
9:4:801:MET:HE2	9:4:822:VAL:HG13	1.72	0.71
12:7:591:LEU:HD22	12:7:681:PHE:CZ	2.26	0.71
2:B:220:PRO:HB3	6:F:19:GLN:HA	1.71	0.71
10:5:279:ASP:O	10:5:283:THR:HG23	1.91	0.71
10:5:655:ALA:O	10:5:659:ILE:HD12	1.91	0.71
3:C:361:MET:HE1	6:F:421:TRP:CZ2	2.26	0.70
8:3:702:LEU:O	8:3:706:ILE:HD12	1.91	0.70
9:4:271:ILE:O	9:4:275:THR:HG23	1.93	0.69
3:C:374:ALA:HB1	3:C:377:ILE:HD11	1.73	0.69
7:2:624:MET:HE1	7:2:677:PHE:CE1	2.28	0.68
3:C:513:LEU:HD22	3:C:556:PHE:CD2	2.29	0.68
2:B:192:GLN:CD	2:B:193:ILE:HG23	2.19	0.68
13:X:64:DC:H2'	13:X:65:DT:H72	1.76	0.68
1:A:809:LEU:HD11	1:A:901:ALA:CB	2.25	0.67
2:B:221:VAL:HG22	6:F:283:GLU:OE2	1.94	0.67
9:4:648:VAL:O	9:4:652:GLN:NE2	2.28	0.67
12:7:591:LEU:HD23	12:7:591:LEU:O	1.94	0.67
6:F:282:THR:HG22	6:F:286:ILE:HD12	1.76	0.67
12:7:394:THR:HG22	12:7:397:VAL:HG22	1.74	0.67
12:7:354:ILE:HD12	12:7:356:LEU:HD21	1.75	0.67
2:B:285:PHE:CE1	3:C:513:LEU:HD23	2.30	0.67
3:C:288:VAL:HG12	3:C:288:VAL:O	1.94	0.67
9:4:775:VAL:HG21	11:6:725:THR:HG22	1.77	0.67
3:C:513:LEU:HD11	3:C:517:MET:HE3	1.76	0.66
10:5:375:ALA:HB1	10:5:378:ILE:HD13	1.77	0.66
4:D:470:THR:HG22	4:D:471:GLU:H	1.59	0.66
9:4:602:THR:OG1	9:4:656:ILE:HD11	1.95	0.66
2:B:221:VAL:HB	6:F:76:ASP:OD2	1.96	0.66
3:C:513:LEU:HD11	3:C:517:MET:CE	2.26	0.65
9:4:708:VAL:HG12	9:4:708:VAL:O	1.96	0.65
3:C:410:ALA:HB2	3:C:424:LEU:HD21	1.76	0.65
5:E:182:HIS:HB2	5:E:184:ILE:HD11	1.77	0.65
8:3:443:THR:HG21	8:3:457:LEU:HD22	1.79	0.65
12:7:207:LEU:HD13	12:7:212:ALA:HB2	1.79	0.65
7:2:276:MET:HE2	7:2:293:ILE:HD11	1.79	0.65
10:5:417:ASP:HB3	10:5:418:PRO:HD2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:299:VAL:HG23	3:C:425:ILE:HG23	1.79	0.65
3:C:141:MET:HE1	13:X:15:DT:OP1	1.96	0.64
1:A:459:ILE:HD11	1:A:492:VAL:HG21	1.79	0.64
2:B:219:THR:HG23	6:F:296:ARG:HG2	1.80	0.64
13:X:76:DC:H2''	13:X:77:DT:H72	1.80	0.63
7:2:567:THR:HG22	7:2:568:GLY:H	1.62	0.63
11:6:374:PRO:O	11:6:376:THR:HG23	1.98	0.63
9:4:585:THR:HG21	9:4:628:VAL:HB	1.80	0.63
10:5:366:LEU:HD21	10:5:666:LEU:HD22	1.80	0.63
11:6:124:VAL:HG13	11:6:124:VAL:O	1.98	0.63
3:C:515:LYS:HG3	3:C:516:ILE:HG23	1.81	0.63
7:2:222:THR:HG22	7:2:222:THR:O	1.98	0.63
3:C:242:THR:CG2	5:E:430:LEU:HD22	2.29	0.63
2:B:207:PRO:HA	6:F:345:PHE:CE1	2.34	0.63
10:5:677:VAL:O	10:5:681:ILE:HD12	1.98	0.63
2:B:220:PRO:HD2	6:F:296:ARG:HD3	1.80	0.62
2:B:315:LEU:HD21	2:B:317:TYR:CD2	2.35	0.62
2:B:362:ILE:HD11	3:C:23:TYR:CZ	2.34	0.62
5:E:239:PHE:CE2	5:E:243:ILE:HD11	2.35	0.62
10:5:35:ILE:HG22	10:5:36:LEU:HD12	1.81	0.62
2:B:277:VAL:HG22	3:C:566:LEU:HD13	1.81	0.62
3:C:142:MET:HE3	3:C:219:PHE:CG	2.34	0.62
3:C:547:THR:HG22	3:C:547:THR:O	1.99	0.62
2:B:315:LEU:HD22	2:B:477:PHE:HA	1.82	0.61
12:7:591:LEU:HD22	12:7:681:PHE:HZ	1.64	0.61
13:X:60:DC:H2'	13:X:61:DT:H72	1.83	0.61
2:B:210:LEU:HD13	6:F:345:PHE:CE2	2.35	0.61
7:2:813:ILE:HG13	7:2:841:VAL:HG21	1.81	0.61
9:4:680:SER:HB3	12:7:588:ALA:HB3	1.83	0.61
9:4:714:GLU:O	9:4:715:LYS:HG2	2.01	0.61
2:B:210:LEU:HD13	6:F:345:PHE:CD2	2.36	0.61
13:X:5:DT:H2'	13:X:6:DT:H72	1.81	0.61
3:C:524:LEU:HD23	3:C:544:PHE:CE1	2.36	0.61
14:Y:37:DG:H2'	14:Y:38:DT:H72	1.82	0.61
4:D:313:LEU:HD12	4:D:321:ASN:HD21	1.67	0.60
10:5:540:ILE:HG22	10:5:541:ASP:H	1.66	0.60
2:B:212:LEU:HD23	6:F:335:ARG:NH2	2.16	0.60
10:5:437:VAL:CG1	10:5:471:LEU:HD22	2.32	0.60
10:5:497:MET:CE	10:5:519:VAL:HG11	2.32	0.60
7:2:591:LEU:HD21	7:2:636:ILE:HD13	1.81	0.60
4:D:461:GLU:HA	4:D:464:VAL:HG22	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:468:MET:HE3	3:C:113:THR:HG21	1.83	0.59
1:A:456:PHE:HD1	1:A:492:VAL:HG22	1.66	0.59
4:D:253:LYS:O	4:D:256:ILE:HD11	2.01	0.59
9:4:750:TYR:HE2	9:4:813:LEU:HD21	1.64	0.59
4:D:114:TYR:CZ	4:D:118:LEU:HD11	2.37	0.59
2:B:212:LEU:HD13	6:F:332:PHE:HE2	1.65	0.59
7:2:584:PRO:HG2	7:2:585:ILE:HD12	1.84	0.59
12:7:313:CYS:SG	12:7:333:ILE:HD12	2.43	0.59
6:F:351:CYS:SG	6:F:359:VAL:HG12	2.43	0.59
5:E:242:LEU:HD13	5:E:285:TYR:CD1	2.38	0.58
8:3:333:SER:OG	10:5:512:VAL:HG23	2.03	0.58
8:3:386:MET:HE1	8:3:663:ALA:HB1	1.85	0.58
7:2:624:MET:HE2	7:2:646:ILE:HB	1.85	0.58
1:A:809:LEU:HD11	1:A:901:ALA:HB3	1.86	0.58
4:D:305:TRP:O	4:D:309:LEU:HD23	2.04	0.58
10:5:556:VAL:O	10:5:556:VAL:HG13	2.04	0.58
11:6:533:ILE:HG22	11:6:533:ILE:O	2.04	0.58
6:F:40:TYR:HB3	7:2:269:LYS:HZ1	1.68	0.57
2:B:193:ILE:HG13	2:B:194:ILE:N	2.19	0.57
7:2:838:ILE:O	7:2:842:VAL:HG23	2.04	0.57
3:C:214:ASN:HD22	4:D:495:THR:HG23	1.69	0.57
2:B:263:MET:HE1	3:C:605:TYR:O	2.05	0.57
2:B:420:LEU:O	2:B:450:ILE:HG22	2.04	0.57
3:C:212:ASN:HD21	4:D:495:THR:HG21	1.69	0.57
12:7:315:ILE:CD1	12:7:333:ILE:HD11	2.34	0.57
3:C:130:GLU:C	3:C:131:LEU:HD22	2.30	0.57
11:6:135:VAL:HG12	11:6:135:VAL:O	2.03	0.57
1:A:456:PHE:CD1	1:A:492:VAL:HG22	2.40	0.57
12:7:245:ILE:HD11	12:7:349:VAL:HG21	1.86	0.57
1:A:809:LEU:HD11	1:A:901:ALA:HB1	1.87	0.57
8:3:729:ALA:O	8:3:733:LEU:HD23	2.05	0.56
2:B:190:LEU:HD11	3:C:362:PHE:HB2	1.87	0.56
9:4:224:LEU:HD22	9:4:227:ILE:HG13	1.87	0.56
9:4:763:THR:HG21	9:4:817:VAL:O	2.04	0.56
11:6:762:LYS:HD3	11:6:762:LYS:O	2.04	0.56
5:E:95:THR:HG23	5:E:95:THR:O	2.06	0.56
7:2:276:MET:HE2	7:2:293:ILE:CD1	2.35	0.56
10:5:425:LEU:O	10:5:429:VAL:HG23	2.05	0.56
1:A:875:ASP:O	4:D:253:LYS:HD2	2.06	0.56
3:C:381:LEU:HD23	3:C:382:THR:N	2.21	0.56
12:7:557:LEU:HD23	12:7:558:ASN:N	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:459:VAL:O	4:D:463:LEU:HD23	2.06	0.56
2:B:242:THR:HG22	2:B:243:PHE:N	2.21	0.56
11:6:710:ASP:C	11:6:711:LEU:HD22	2.30	0.56
3:C:214:ASN:ND2	4:D:495:THR:HG23	2.21	0.56
12:7:435:LEU:HD13	12:7:456:VAL:HG23	1.88	0.56
1:A:870:LEU:O	1:A:874:LEU:HD23	2.06	0.56
5:E:200:ILE:HD13	5:E:203:MET:HE2	1.88	0.56
1:A:564:LEU:HD12	1:A:596:ILE:HG23	1.88	0.55
3:C:242:THR:HG23	5:E:430:LEU:HD22	1.86	0.55
7:2:212:LYS:HD3	7:2:274:VAL:HG13	1.87	0.55
10:5:387:ALA:HB1	10:5:552:MET:SD	2.46	0.55
9:4:282:SER:O	9:4:285:VAL:HG12	2.07	0.55
10:5:278:CYS:SG	10:5:283:THR:HG22	2.46	0.55
7:2:477:THR:O	7:2:481:GLU:OE1	2.24	0.55
9:4:727:LEU:HD21	9:4:736:ILE:HD12	1.88	0.55
4:D:294:VAL:HG11	4:D:302:VAL:HG13	1.88	0.55
10:5:366:LEU:CD2	10:5:666:LEU:HD22	2.36	0.55
2:B:315:LEU:HD21	2:B:317:TYR:HD2	1.71	0.55
7:2:497:ILE:HG21	7:2:823:MET:HE3	1.88	0.55
6:F:372:LEU:O	6:F:372:LEU:HD23	2.07	0.55
9:4:769:GLU:HB2	9:4:819:LEU:HD11	1.89	0.55
8:3:709:ALA:HB1	8:3:721:VAL:HG11	1.89	0.55
13:X:7:DT:H2''	13:X:8:DT:H71	1.88	0.55
7:2:446:VAL:HG13	11:6:302:PRO:O	2.07	0.54
2:B:192:GLN:HA	2:B:195:MET:HE2	1.89	0.54
11:6:671:THR:C	11:6:672:LEU:HD12	2.31	0.54
9:4:227:ILE:HG22	9:4:231:ASN:OD1	2.08	0.54
1:A:822:LEU:HD22	1:A:887:ASP:O	2.07	0.54
10:5:622:LEU:HD21	10:5:681:ILE:HD11	1.89	0.54
13:X:4:DT:H2'	13:X:5:DT:H72	1.88	0.54
8:3:463:VAL:HG21	8:3:506:LEU:HD13	1.90	0.54
8:3:274:ILE:HD12	8:3:321:ILE:HD12	1.90	0.54
10:5:167:ILE:HD11	10:5:259:GLN:OE1	2.08	0.54
11:6:407:VAL:HG12	11:6:407:VAL:O	2.08	0.54
6:F:369:VAL:O	6:F:373:ILE:HG22	2.07	0.54
7:2:813:ILE:CG1	7:2:841:VAL:HG21	2.38	0.53
8:3:357:LYS:HD3	8:3:357:LYS:N	2.23	0.53
12:7:394:THR:HG23	12:7:396:ASP:H	1.73	0.53
1:A:699:VAL:HG13	1:A:700:SER:N	2.23	0.53
9:4:721:ALA:HB1	12:7:661:VAL:HG22	1.90	0.53
9:4:727:LEU:CD2	9:4:736:ILE:HD12	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:849:LEU:HD21	1:A:863:ILE:HB	1.89	0.53
3:C:241:ASN:O	3:C:241:ASN:ND2	2.41	0.53
7:2:778:LEU:HD23	10:5:577:THR:OG1	2.07	0.53
9:4:419:VAL:HG12	9:4:463:VAL:HG11	1.90	0.53
3:C:143:LEU:HD11	3:C:222:LEU:HD13	1.90	0.53
5:E:435:THR:HG22	5:E:436:THR:N	2.24	0.53
12:7:437:VAL:HG11	12:7:698:ALA:HB1	1.89	0.53
13:X:26:DT:H2'	13:X:27:DT:H71	1.90	0.53
7:2:493:ILE:HD11	7:2:823:MET:O	2.08	0.53
1:A:863:ILE:O	1:A:866:TRP:HD1	1.91	0.53
4:D:70:ILE:HD11	4:D:290:ASN:HB3	1.91	0.53
1:A:572:VAL:HB	1:A:576:GLN:HA	1.91	0.53
3:C:346:LEU:HD21	3:C:351:LEU:HD21	1.91	0.53
3:C:517:MET:HE2	3:C:549:PRO:HG2	1.91	0.53
5:E:194:VAL:HG23	5:E:195:ASP:N	2.24	0.53
10:5:469:MET:HE3	10:5:469:MET:HA	1.90	0.53
12:7:394:THR:O	12:7:398:GLU:OE1	2.27	0.53
2:B:473:TYR:CE2	2:B:475:PHE:CE1	2.97	0.52
6:F:331:VAL:HG13	6:F:386:TYR:CE1	2.43	0.52
10:5:341:SER:HB2	10:5:435:ILE:HG22	1.91	0.52
3:C:569:LEU:O	3:C:575:ALA:HB2	2.09	0.52
8:3:166:LEU:HD12	8:3:170:THR:OG1	2.10	0.52
3:C:513:LEU:HD22	3:C:556:PHE:CG	2.44	0.52
5:E:129:PHE:HD2	5:E:167:ILE:HD11	1.74	0.52
6:F:292:PHE:CZ	6:F:350:MET:HE3	2.44	0.52
7:2:806:THR:HG22	7:2:807:VAL:N	2.25	0.52
8:3:695:SER:HB2	8:3:737:LEU:HD23	1.91	0.52
9:4:562:ILE:HG13	9:4:562:ILE:O	2.09	0.52
12:7:537:ILE:HG23	12:7:563:ILE:HD13	1.91	0.52
2:B:207:PRO:HA	6:F:345:PHE:HE1	1.74	0.52
3:C:361:MET:HE1	6:F:421:TRP:CH2	2.44	0.52
11:6:304:LEU:HD11	11:6:307:ALA:HB2	1.92	0.52
2:B:190:LEU:O	2:B:194:ILE:HG22	2.10	0.52
7:2:609:PHE:O	7:2:612:MET:HE2	2.10	0.52
12:7:591:LEU:HD11	12:7:597:LEU:CD1	2.40	0.52
8:3:209:PHE:HB2	12:7:8:ILE:HG23	1.92	0.52
12:7:404:LEU:HD22	12:7:641:TYR:CE2	2.45	0.52
3:C:139:VAL:HA	3:C:142:MET:HE2	1.91	0.52
8:3:425:THR:HB	8:3:656:LEU:HD21	1.92	0.52
13:X:5:DT:C2'	13:X:6:DT:H72	2.40	0.52
3:C:587:ALA:O	3:C:591:PHE:HD1	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:67:LEU:HD23	5:E:73:LEU:HD22	1.93	0.51
5:E:199:THR:HA	5:E:202:VAL:HG22	1.91	0.51
7:2:850:LYS:O	7:2:853:VAL:HG12	2.10	0.51
8:3:700:ARG:O	8:3:704:THR:HG23	2.10	0.51
1:A:699:VAL:HG13	1:A:700:SER:H	1.74	0.51
2:B:446:VAL:HG12	2:B:446:VAL:O	2.08	0.51
5:E:462:ILE:O	5:E:466:VAL:HG22	2.09	0.51
2:B:362:ILE:HD11	3:C:23:TYR:CE2	2.45	0.51
11:6:132:VAL:HG23	11:6:132:VAL:O	2.10	0.51
5:E:228:THR:HG23	5:E:231:GLN:H	1.75	0.51
1:A:475:ILE:HG23	1:A:596:ILE:HG13	1.93	0.51
2:B:383:LEU:HD11	2:B:411:TYR:OH	2.10	0.51
4:D:375:ASN:C	4:D:376:LEU:HD22	2.35	0.51
8:3:443:THR:CG2	8:3:457:LEU:HD22	2.40	0.51
1:A:852:GLN:NE2	4:D:326:MET:SD	2.84	0.51
2:B:190:LEU:HG	3:C:362:PHE:CD1	2.45	0.51
5:E:99:PRO:O	5:E:102:VAL:HG12	2.11	0.51
9:4:528:PRO:O	9:4:736:ILE:HD11	2.11	0.51
1:A:906:ASN:O	1:A:906:ASN:ND2	2.42	0.51
8:3:171:LEU:HD11	8:3:180:VAL:HG21	1.93	0.51
13:X:35:DT:H2"	13:X:36:DT:H71	1.93	0.51
5:E:38:TYR:O	5:E:41:THR:HG22	2.11	0.50
8:3:389:VAL:HG23	8:3:668:ILE:HG23	1.93	0.50
5:E:194:VAL:HG12	5:E:248:HIS:NE2	2.26	0.50
7:2:342:LEU:HD12	7:2:372:PRO:HB2	1.93	0.50
9:4:801:MET:HE1	9:4:826:VAL:HB	1.94	0.50
10:5:389:VAL:CG1	10:5:666:LEU:HD21	2.42	0.50
5:E:71:LYS:O	5:E:75:GLN:NE2	2.44	0.50
10:5:540:ILE:HG22	10:5:541:ASP:N	2.25	0.50
1:A:811:LEU:HD13	1:A:820:GLN:HG2	1.93	0.50
2:B:274:PHE:CD1	3:C:582:ILE:HG23	2.47	0.50
2:B:288:ARG:HB3	2:B:289:PRO:HD3	1.94	0.50
7:2:322:GLY:HA3	7:2:390:LEU:HD11	1.94	0.50
11:6:522:ASP:HB2	11:6:525:ILE:HG23	1.94	0.50
12:7:633:VAL:HG12	12:7:634:GLU:N	2.26	0.50
1:A:548:PHE:CE1	1:A:552:ARG:HG2	2.47	0.50
12:7:633:VAL:HG11	12:7:637:LYS:HD2	1.94	0.50
13:X:6:DT:H2"	13:X:7:DT:H71	1.93	0.50
2:B:319:VAL:HG21	3:C:491:LEU:HD21	1.94	0.50
11:6:679:LEU:C	11:6:679:LEU:HD23	2.36	0.50
12:7:494:THR:HG22	12:7:495:ALA:N	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:LEU:O	1:A:449:LEU:HD12	2.12	0.49
2:B:212:LEU:HD13	6:F:332:PHE:CE2	2.47	0.49
4:D:119:LEU:HD23	4:D:127:PHE:CE2	2.47	0.49
8:3:100:LEU:HD11	8:3:161:PHE:CE1	2.47	0.49
2:B:230:THR:O	2:B:230:THR:HG23	2.13	0.49
4:D:410:LEU:HD22	5:E:375:TYR:HB3	1.94	0.49
7:2:211:LEU:HD11	7:2:274:VAL:HG21	1.93	0.49
4:D:346:ALA:HB1	5:E:20:SER:HB3	1.94	0.49
8:3:486:ILE:HA	8:3:489:VAL:HG12	1.95	0.49
11:6:630:LEU:O	11:6:631:ALA:HB3	2.13	0.49
3:C:479:ILE:HG22	3:C:480:PHE:CD1	2.48	0.49
7:2:539:VAL:HG12	7:2:540:LEU:N	2.27	0.49
10:5:453:VAL:HG11	10:5:509:ILE:HD13	1.94	0.49
14:Y:37:DG:C2'	14:Y:38:DT:H72	2.42	0.49
7:2:212:LYS:CD	7:2:274:VAL:HG13	2.43	0.49
8:3:294:VAL:HG12	8:3:326:VAL:HG22	1.94	0.49
5:E:81:VAL:HG13	5:E:82:GLN:N	2.28	0.49
9:4:521:LEU:HD21	9:4:742:LEU:HD11	1.93	0.49
9:4:651:GLN:O	9:4:653:THR:HG23	2.12	0.49
11:6:548:LEU:HD23	11:6:548:LEU:C	2.37	0.49
6:F:285:ILE:HD11	6:F:303:VAL:HG22	1.93	0.49
9:4:609:VAL:HG21	12:7:505:GLU:N	2.28	0.49
12:7:86:LEU:HD23	12:7:90:ASN:OD1	2.13	0.49
14:Y:11:DC:H2''	14:Y:12:DA:C8	2.48	0.49
8:3:653:ILE:N	8:3:654:PRO:HD2	2.28	0.49
9:4:342:MET:HE3	9:4:342:MET:HA	1.94	0.48
3:C:491:LEU:HD22	3:C:521:LEU:HD21	1.95	0.48
4:D:238:VAL:HG22	4:D:245:VAL:HB	1.95	0.48
11:6:386:VAL:HG12	11:6:387:GLU:N	2.28	0.48
1:A:488:THR:O	1:A:492:VAL:HG23	2.13	0.48
5:E:194:VAL:HG23	5:E:195:ASP:H	1.78	0.48
1:A:811:LEU:HD22	1:A:820:GLN:NE2	2.28	0.48
7:2:324:VAL:HG23	7:2:424:VAL:HG21	1.94	0.48
8:3:568:THR:HG22	8:3:568:THR:O	2.12	0.48
9:4:311:CYS:SG	9:4:326:ILE:HD11	2.53	0.48
9:4:527:ALA:HB1	9:4:530:ILE:HD12	1.95	0.48
9:4:747:LEU:HD23	9:4:751:ILE:HD12	1.95	0.48
10:5:170:SER:O	10:5:171:VAL:HG23	2.13	0.48
1:A:866:TRP:HE3	1:A:870:LEU:HD12	1.79	0.48
7:2:306:LEU:HD11	7:2:406:ARG:CG	2.44	0.48
7:2:567:THR:HG23	7:2:607:ASP:OD1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:X:7:DT:C2'	13:X:8:DT:H71	2.43	0.48
2:B:269:VAL:HG11	2:B:273:GLU:OE2	2.13	0.48
7:2:510:ASP:OD2	7:2:511:ILE:N	2.47	0.48
7:2:764:MET:O	7:2:767:ILE:HG22	2.13	0.48
10:5:663:LEU:O	10:5:663:LEU:HD23	2.14	0.48
2:B:461:ALA:N	2:B:462:PRO:CD	2.77	0.48
3:C:219:PHE:O	3:C:223:LEU:HD13	2.12	0.48
12:7:343:LEU:HD21	12:7:381:VAL:HG21	1.95	0.48
1:A:388:LEU:HD12	1:A:394:PHE:CD1	2.49	0.48
4:D:52:LEU:HD23	4:D:56:LEU:HD23	1.96	0.48
4:D:117:SER:O	4:D:120:GLN:HG2	2.14	0.48
4:D:497:ILE:HD11	4:D:502:ARG:HB3	1.96	0.48
4:D:90:ILE:HD13	4:D:211:THR:HG21	1.94	0.47
4:D:230:LEU:O	4:D:234:LEU:HD13	2.14	0.47
8:3:363:LEU:O	8:3:367:LEU:HG	2.14	0.47
7:2:624:MET:HE2	7:2:646:ILE:CB	2.44	0.47
14:Y:57:DT:C2	14:Y:58:DA:N7	2.83	0.47
1:A:896:ILE:HG13	1:A:897:SER:H	1.78	0.47
3:C:517:MET:HE1	3:C:553:ILE:HB	1.95	0.47
8:3:474:GLU:OE2	10:5:491:VAL:HG13	2.14	0.47
2:B:361:LEU:HD23	2:B:361:LEU:H	1.79	0.47
3:C:92:TYR:OH	3:C:118:LEU:HD12	2.14	0.47
7:2:297:ILE:HG22	7:2:298:SER:N	2.29	0.47
9:4:820:GLU:O	9:4:824:GLU:OE1	2.32	0.47
12:7:591:LEU:HD21	12:7:597:LEU:HD13	1.95	0.47
3:C:130:GLU:N	3:C:130:GLU:OE1	2.48	0.47
5:E:89:TYR:O	5:E:92:ILE:HG22	2.14	0.47
10:5:437:VAL:HG13	10:5:472:ALA:HB2	1.97	0.47
6:F:49:LEU:O	6:F:50:LYS:HE2	2.14	0.47
7:2:501:MET:HE3	7:2:516:ALA:HB2	1.96	0.47
7:2:567:THR:HG22	7:2:568:GLY:N	2.29	0.47
8:3:214:TYR:CE2	8:3:229:ALA:HB3	2.50	0.47
8:3:279:ASP:OD2	8:3:280:ASP:N	2.47	0.47
9:4:818:GLU:N	9:4:818:GLU:OE1	2.48	0.47
3:C:205:PHE:HB2	3:C:208:VAL:HG22	1.96	0.47
3:C:242:THR:HG22	5:E:430:LEU:HD22	1.96	0.47
10:5:339:THR:HG22	10:5:340:SER:N	2.27	0.47
12:7:361:THR:HG23	12:7:361:THR:O	2.14	0.47
12:7:651:VAL:HG12	12:7:652:MET:N	2.30	0.47
3:C:513:LEU:HD13	3:C:556:PHE:HB2	1.96	0.47
8:3:411:PRO:HG2	10:5:544:THR:HG23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:7:428:VAL:O	12:7:432:LEU:HD23	2.14	0.47
12:7:493:LEU:HD22	12:7:512:ALA:HB3	1.96	0.47
6:F:39:LEU:C	6:F:39:LEU:HD23	2.40	0.47
8:3:307:ASN:HB3	10:5:207:LEU:HD12	1.97	0.47
8:3:197:ILE:HD11	8:3:251:ILE:HB	1.97	0.46
9:4:333:LEU:HD11	9:4:400:GLN:OE1	2.14	0.46
12:7:114:THR:HG23	12:7:115:GLU:N	2.31	0.46
3:C:409:VAL:HG22	3:C:453:LEU:HD11	1.97	0.46
2:B:207:PRO:HA	6:F:345:PHE:CZ	2.50	0.46
5:E:18:LEU:CD2	5:E:187:VAL:HG21	2.46	0.46
7:2:687:VAL:HG12	7:2:687:VAL:O	2.16	0.46
14:Y:80:DT:H2''	14:Y:81:DA:N7	2.29	0.46
4:D:119:LEU:HD23	4:D:127:PHE:CD2	2.50	0.46
7:2:778:LEU:HD12	7:2:783:MET:CE	2.45	0.46
8:3:383:LEU:HD12	8:3:711:ALA:HB3	1.98	0.46
6:F:338:LYS:O	6:F:338:LYS:HG2	2.15	0.46
7:2:783:MET:SD	10:5:573:ILE:CG2	3.04	0.46
8:3:201:HIS:HB3	8:3:241:LEU:HD22	1.97	0.46
8:3:703:GLU:OE2	12:7:620:HIS:CD2	2.69	0.46
13:X:36:DT:H2''	13:X:37:DC:C6	2.50	0.46
1:A:709:VAL:HG21	1:A:782:LEU:HD13	1.98	0.46
5:E:22:ILE:HD12	5:E:129:PHE:HZ	1.80	0.46
6:F:306:TYR:HD1	6:F:319:LEU:HD11	1.81	0.46
12:7:27:THR:O	12:7:27:THR:HG22	2.15	0.46
12:7:245:ILE:CD1	12:7:349:VAL:HG11	2.45	0.46
13:X:37:DC:H2''	13:X:38:DC:C5	2.51	0.46
13:X:82:DG:H2'	13:X:83:DT:H72	1.98	0.46
4:D:342:ILE:N	4:D:343:PRO:HD2	2.30	0.46
8:3:367:LEU:HD12	8:3:378:LYS:HB3	1.98	0.46
9:4:725:THR:HG21	12:7:657:ASN:CG	2.41	0.46
10:5:46:TYR:O	10:5:50:LEU:HD23	2.16	0.46
11:6:790:ARG:O	11:6:835:ILE:HD12	2.15	0.46
12:7:66:MET:HE1	12:7:79:ILE:HG22	1.98	0.46
3:C:334:PHE:CE2	3:C:465:LEU:HD13	2.51	0.46
5:E:198:SER:O	5:E:202:VAL:HG13	2.16	0.46
10:5:35:ILE:CG2	10:5:36:LEU:HD12	2.46	0.46
10:5:339:THR:O	10:5:340:SER:C	2.59	0.46
10:5:569:ALA:O	10:5:573:ILE:HG13	2.16	0.46
3:C:265:LYS:C	3:C:266:LEU:HD22	2.40	0.46
3:C:382:THR:HG22	3:C:388:LEU:HD22	1.98	0.46
3:C:479:ILE:HG22	3:C:480:PHE:CE1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:380:LEU:HA	4:D:383:LEU:HD23	1.97	0.46
7:2:796:GLU:HB3	7:2:859:ARG:NE	2.31	0.46
10:5:282:LEU:HD21	10:5:330:ILE:HG22	1.98	0.46
10:5:301:TYR:HD1	10:5:327:TYR:CE1	2.34	0.46
11:6:326:LYS:HE3	11:6:326:LYS:HA	1.98	0.46
2:B:422:LEU:HD23	2:B:422:LEU:H	1.81	0.46
4:D:361:CYS:HB3	4:D:364:LEU:HD12	1.98	0.46
4:D:497:ILE:HB	4:D:498:PRO:HD2	1.98	0.46
5:E:108:LEU:O	5:E:112:LEU:HD13	2.16	0.46
5:E:214:CYS:SG	5:E:270:VAL:HG11	2.56	0.46
8:3:100:LEU:HD11	8:3:161:PHE:CD1	2.51	0.46
2:B:269:VAL:HG12	2:B:270:THR:N	2.31	0.45
11:6:153:ILE:HG21	11:6:267:PHE:CE1	2.51	0.45
2:B:210:LEU:HD21	2:B:212:LEU:CD2	2.46	0.45
4:D:294:VAL:HG11	4:D:302:VAL:CG1	2.46	0.45
9:4:224:LEU:HD22	9:4:227:ILE:CG1	2.46	0.45
10:5:412:VAL:HG12	10:5:413:LEU:N	2.30	0.45
11:6:134:LYS:N	11:6:134:LYS:CD	2.79	0.45
14:Y:10:DT:H2''	14:Y:11:DC:C5	2.51	0.45
2:B:476:VAL:HG23	2:B:476:VAL:O	2.15	0.45
5:E:286:LYS:O	5:E:289:ILE:HG22	2.16	0.45
8:3:382:LEU:C	8:3:382:LEU:HD23	2.42	0.45
9:4:318:ASN:HB2	9:4:319:PRO:HD2	1.97	0.45
9:4:602:THR:CG2	9:4:656:ILE:HD11	2.47	0.45
8:3:538:SER:O	8:3:542:ARG:HG2	2.17	0.45
9:4:199:MET:HE2	9:4:227:ILE:HD11	1.99	0.45
9:4:521:LEU:CD2	9:4:742:LEU:HD11	2.47	0.45
9:4:522:LEU:HB3	9:4:541:LEU:HD13	1.97	0.45
12:7:494:THR:HG22	12:7:495:ALA:H	1.81	0.45
12:7:507:ILE:C	12:7:508:LEU:HD12	2.42	0.45
13:X:4:DT:C2'	13:X:5:DT:H72	2.47	0.45
1:A:811:LEU:HD11	1:A:825:ILE:HA	1.99	0.45
1:A:896:ILE:HG13	1:A:897:SER:N	2.32	0.45
2:B:227:LEU:HG	2:B:228:TYR:CD1	2.52	0.45
4:D:294:VAL:HG12	4:D:296:SER:H	1.81	0.45
10:5:437:VAL:HG11	10:5:471:LEU:HD22	1.99	0.45
10:5:513:LEU:HD23	10:5:513:LEU:H	1.82	0.45
11:6:304:LEU:HD12	11:6:353:PHE:HE1	1.82	0.45
12:7:402:MET:O	12:7:406:THR:HG23	2.16	0.45
1:A:794:MET:HE2	1:A:904:ALA:HB2	1.99	0.45
3:C:246:ASN:C	3:C:246:ASN:HD22	2.24	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:2:494:ILE:HG23	7:2:823:MET:CE	2.46	0.45
9:4:338:VAL:HA	9:4:395:GLN:OE1	2.15	0.45
9:4:361:ASP:OD1	9:4:361:ASP:O	2.34	0.45
11:6:166:LEU:HG	11:6:170:ILE:HD12	1.97	0.45
14:Y:39:DC:H2''	14:Y:40:DA:C8	2.51	0.45
7:2:203:VAL:HG12	7:2:206:THR:OG1	2.17	0.45
10:5:136:GLN:HE22	10:5:282:LEU:HB2	1.82	0.45
13:X:6:DT:C2'	13:X:7:DT:H71	2.46	0.45
2:B:221:VAL:HA	6:F:283:GLU:OE2	2.17	0.45
1:A:836:ASN:O	1:A:842:VAL:HG21	2.17	0.45
8:3:432:THR:HG22	8:3:434:GLY:H	1.82	0.45
9:4:451:ARG:NH2	11:6:445:VAL:HG21	2.32	0.45
10:5:653:LEU:HD23	10:5:653:LEU:C	2.42	0.45
6:F:30:LYS:O	6:F:33:THR:HG22	2.17	0.45
7:2:638:THR:HG22	7:2:639:THR:N	2.31	0.45
9:4:553:THR:HG23	9:4:554:LYS:N	2.32	0.45
10:5:472:ALA:O	10:5:517:THR:HG22	2.17	0.45
11:6:178:LEU:HD11	11:6:265:ILE:HD11	1.97	0.45
1:A:839:ASN:HB3	1:A:842:VAL:HG22	1.99	0.44
3:C:405:HIS:HA	3:C:408:PHE:CE2	2.53	0.44
3:C:520:VAL:O	3:C:524:LEU:HG	2.17	0.44
4:D:105:GLN:HB3	4:D:335:PRO:HG2	1.99	0.44
5:E:77:ILE:CD1	5:E:130:LEU:HD11	2.47	0.44
7:2:574:VAL:HG12	7:2:575:GLY:N	2.32	0.44
11:6:125:GLN:O	11:6:126:SER:C	2.60	0.44
11:6:701:MET:HE3	11:6:705:ILE:HG13	1.99	0.44
12:7:508:LEU:HD21	12:7:553:ILE:HG13	1.99	0.44
1:A:625:THR:HG22	1:A:626:GLY:N	2.33	0.44
10:5:420:THR:OG1	10:5:556:VAL:HG22	2.18	0.44
12:7:333:ILE:HG22	12:7:334:HIS:N	2.32	0.44
13:X:74:DG:H2''	13:X:75:DA:C8	2.53	0.44
4:D:281:LEU:HD23	4:D:285:VAL:HG23	1.99	0.44
5:E:81:VAL:HG13	5:E:82:GLN:H	1.81	0.44
7:2:478:GLU:HG2	7:2:479:GLU:N	2.32	0.44
9:4:298:THR:HG22	9:4:298:THR:O	2.17	0.44
12:7:248:VAL:HG12	12:7:313:CYS:CB	2.48	0.44
2:B:365:GLY:HA3	2:B:432:ILE:HD13	1.99	0.44
4:D:286:ASP:OD1	4:D:287:ALA:N	2.51	0.44
4:D:459:VAL:HG12	4:D:463:LEU:HD23	1.99	0.44
5:E:151:LYS:HA	5:E:154:GLU:OE1	2.17	0.44
6:F:47:VAL:HG21	11:6:349:THR:HB	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:321:CYS:O	6:F:325:LEU:HD13	2.18	0.44
8:3:415:LYS:HG3	8:3:416:SER:N	2.33	0.44
9:4:332:VAL:HG11	9:4:428:ARG:C	2.43	0.44
10:5:69:ILE:HG23	10:5:70:GLY:N	2.33	0.44
12:7:13:ASP:O	12:7:17:LEU:HD13	2.18	0.44
4:D:329:GLU:OE2	4:D:330:THR:HG23	2.17	0.44
4:D:509:ILE:HG22	4:D:510:LEU:N	2.33	0.44
5:E:200:ILE:HD13	5:E:203:MET:CE	2.47	0.44
7:2:591:LEU:HD21	7:2:636:ILE:CD1	2.46	0.44
8:3:448:THR:HG23	8:3:448:THR:O	2.17	0.44
13:X:38:DC:H2"	13:X:39:DC:C6	2.53	0.44
6:F:292:PHE:HB3	6:F:294:LEU:CD1	2.48	0.44
7:2:246:TYR:HA	7:2:249:LEU:HD21	1.98	0.44
9:4:553:THR:HG23	9:4:554:LYS:H	1.82	0.44
11:6:513:ILE:HG23	11:6:514:ASN:N	2.32	0.44
12:7:430:LYS:O	12:7:434:LEU:HD13	2.17	0.44
6:F:342:ILE:O	6:F:343:ASP:C	2.60	0.44
7:2:213:SER:O	7:2:217:GLU:HG3	2.18	0.44
8:3:36:THR:HG22	8:3:36:THR:O	2.18	0.44
10:5:261:ILE:HG22	10:5:263:GLU:H	1.82	0.44
10:5:279:ASP:OD1	10:5:279:ASP:N	2.51	0.44
1:A:875:ASP:O	4:D:253:LYS:HA	2.17	0.44
1:A:899:GLU:O	1:A:903:ARG:HG2	2.18	0.44
2:B:285:PHE:HE1	3:C:513:LEU:HD23	1.81	0.44
3:C:545:ARG:HD3	3:C:576:PHE:CE2	2.53	0.44
3:C:611:CYS:SG	3:C:612:VAL:HG13	2.58	0.44
6:F:372:LEU:HD23	6:F:372:LEU:C	2.43	0.44
9:4:720:LEU:C	9:4:720:LEU:HD23	2.43	0.44
10:5:66:GLU:O	10:5:69:ILE:HG22	2.17	0.44
11:6:290:ILE:HG22	11:6:291:SER:N	2.32	0.44
10:5:170:SER:C	10:5:171:VAL:HG23	2.43	0.44
1:A:863:ILE:O	1:A:866:TRP:CD1	2.71	0.43
12:7:553:ILE:HD12	12:7:553:ILE:H	1.82	0.43
3:C:142:MET:HE3	3:C:219:PHE:CD1	2.52	0.43
11:6:762:LYS:NZ	11:6:764:ILE:HD11	2.33	0.43
1:A:811:LEU:HD21	1:A:828:GLU:HB2	1.99	0.43
9:4:451:ARG:HE	11:6:445:VAL:HG11	1.82	0.43
11:6:571:ILE:HG22	11:6:572:CYS:N	2.33	0.43
3:C:351:LEU:HD22	3:C:388:LEU:HD23	2.00	0.43
4:D:56:LEU:HD11	4:D:345:VAL:HG21	2.00	0.43
9:4:314:MET:SD	9:4:415:ILE:HD11	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:4:351:VAL:O	9:4:352:CYS:SG	2.76	0.43
10:5:391:LEU:HD23	10:5:391:LEU:C	2.44	0.43
10:5:458:MET:O	10:5:459:THR:OG1	2.34	0.43
11:6:533:ILE:O	11:6:533:ILE:CG2	2.66	0.43
11:6:795:ILE:HG22	11:6:796:THR:N	2.32	0.43
1:A:866:TRP:CE3	1:A:870:LEU:HD12	2.53	0.43
5:E:106:PHE:O	5:E:109:VAL:HG12	2.18	0.43
9:4:342:MET:HE1	9:4:389:CYS:CB	2.49	0.43
10:5:321:VAL:HG23	10:5:322:ALA:N	2.34	0.43
14:Y:28:DA:H2"	14:Y:29:DG:C8	2.54	0.43
1:A:516:LEU:HG	1:A:571:MET:HE1	2.01	0.43
2:B:469:LYS:O	2:B:473:TYR:HB2	2.18	0.43
3:C:104:THR:HA	3:C:236:LEU:O	2.18	0.43
4:D:86:ILE:O	4:D:90:ILE:HG12	2.18	0.43
9:4:650:GLU:HG2	9:4:701:ARG:NH2	2.33	0.43
11:6:293:THR:HG22	11:6:294:VAL:N	2.33	0.43
2:B:210:LEU:HD21	2:B:212:LEU:HD23	2.01	0.43
4:D:470:THR:HG22	4:D:471:GLU:N	2.30	0.43
5:E:372:ASN:O	5:E:374:ARG:N	2.51	0.43
6:F:280:MET:HE1	6:F:317:TRP:CZ2	2.54	0.43
6:F:299:ALA:O	6:F:303:VAL:HG23	2.18	0.43
7:2:509:ARG:O	7:2:513:THR:HG23	2.19	0.43
7:2:842:VAL:O	7:2:846:VAL:HG23	2.17	0.43
8:3:555:GLU:N	8:3:555:GLU:OE1	2.52	0.43
9:4:678:ILE:HG22	9:4:691:ASN:O	2.19	0.43
9:4:770:LEU:HD21	9:4:802:ILE:HG22	2.00	0.43
10:5:453:VAL:HG11	10:5:509:ILE:CD1	2.47	0.43
2:B:216:PHE:HE1	6:F:295:PRO:HA	1.82	0.43
2:B:250:ARG:HA	3:C:533:MET:HB2	2.00	0.43
3:C:265:LYS:O	3:C:266:LEU:HD22	2.19	0.43
5:E:36:GLN:OE1	5:E:188:MET:SD	2.77	0.43
5:E:38:TYR:HA	5:E:171:LEU:HD13	2.01	0.43
7:2:542:LEU:HD13	7:2:650:ALA:HB3	2.01	0.43
8:3:672:THR:HG22	8:3:673:GLN:N	2.33	0.43
9:4:451:ARG:HH21	11:6:445:VAL:HG21	1.84	0.43
9:4:530:ILE:HG22	9:4:531:TYR:N	2.33	0.43
10:5:286:VAL:HG21	10:5:333:ILE:HD11	2.01	0.43
10:5:365:LYS:HG3	10:5:369:ILE:HD11	2.01	0.43
10:5:539:ASN:O	10:5:540:ILE:HD13	2.19	0.43
12:7:413:ARG:HH22	12:7:414:LEU:HD23	1.84	0.43
12:7:434:LEU:HD23	12:7:695:LEU:CD2	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:628:THR:HG23	1:A:631:GLU:H	1.84	0.43
1:A:684:MET:SD	1:A:684:MET:C	3.02	0.43
4:D:322:ARG:O	4:D:326:MET:HG2	2.18	0.43
5:E:257:ALA:O	5:E:261:LEU:HD23	2.18	0.43
7:2:603:VAL:HG22	7:2:604:CYS:N	2.34	0.43
8:3:235:ASP:OD2	8:3:241:LEU:HD21	2.19	0.43
8:3:421:PHE:O	8:3:425:THR:HG23	2.19	0.43
9:4:591:THR:HA	12:7:547:SER:HB3	2.01	0.43
12:7:718:ARG:O	12:7:722:VAL:HG23	2.19	0.43
13:X:81:DT:H2"	13:X:82:DG:C8	2.54	0.43
3:C:64:LEU:HD21	3:C:288:VAL:CG2	2.48	0.43
4:D:177:LEU:O	4:D:180:VAL:HG12	2.19	0.43
5:E:242:LEU:HB3	5:E:285:TYR:CE1	2.53	0.43
8:3:358:ASP:OD1	8:3:358:ASP:O	2.37	0.43
9:4:384:LEU:HD23	9:4:385:ILE:N	2.34	0.43
12:7:577:ARG:O	12:7:578:LEU:HD22	2.19	0.43
2:B:210:LEU:C	2:B:210:LEU:HD23	2.44	0.42
2:B:315:LEU:HD13	2:B:477:PHE:CD2	2.53	0.42
3:C:130:GLU:O	3:C:131:LEU:HD22	2.19	0.42
3:C:240:ILE:HG22	3:C:241:ASN:N	2.33	0.42
5:E:232:PHE:O	5:E:235:VAL:HG12	2.18	0.42
8:3:486:ILE:CG2	8:3:490:MET:HE2	2.49	0.42
9:4:279:CYS:O	9:4:283:LEU:HD23	2.19	0.42
9:4:801:MET:HE3	9:4:825:ALA:HB3	2.01	0.42
11:6:537:VAL:HG12	11:6:538:PHE:N	2.34	0.42
11:6:663:ILE:CD1	11:6:672:LEU:HD13	2.49	0.42
12:7:517:ASP:OD1	12:7:560:ARG:HB3	2.19	0.42
13:X:67:DA:H2"	13:X:68:DC:OP2	2.19	0.42
13:X:85:DT:O2	14:Y:4:DA:N1	2.51	0.42
3:C:287:THR:HG23	3:C:288:VAL:HG23	2.01	0.42
4:D:208:THR:HG22	4:D:209:LYS:N	2.34	0.42
8:3:254:GLN:HB2	8:3:283:VAL:HG13	2.01	0.42
8:3:276:VAL:HG12	8:3:277:ILE:N	2.34	0.42
9:4:396:VAL:HG21	11:6:412:LEU:HD22	2.01	0.42
10:5:419:GLY:C	10:5:420:THR:HG22	2.44	0.42
11:6:710:ASP:O	11:6:711:LEU:HD22	2.19	0.42
12:7:199:ARG:O	12:7:199:ARG:HG3	2.19	0.42
12:7:502:VAL:O	12:7:502:VAL:HG22	2.19	0.42
2:B:342:GLN:HE22	2:B:358:ILE:HD11	1.84	0.42
3:C:375:ASP:OD1	3:C:376:GLU:N	2.52	0.42
4:D:55:ARG:O	4:D:58:GLN:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:3:116:VAL:HG22	8:3:178:LYS:HD3	2.01	0.42
8:3:223:THR:HG21	10:5:244:ILE:HA	2.01	0.42
9:4:534:GLU:HG2	9:4:535:ASP:N	2.34	0.42
10:5:287:ILE:HG23	10:5:288:PRO:HD2	2.00	0.42
1:A:477:VAL:HG22	1:A:478:ALA:N	2.33	0.42
1:A:803:LEU:HD21	1:A:833:ILE:HD11	2.01	0.42
2:B:469:LYS:O	2:B:473:TYR:CB	2.67	0.42
3:C:247:ILE:O	3:C:247:ILE:HG22	2.20	0.42
3:C:470:ASN:HB2	3:C:475:LEU:HD12	2.01	0.42
4:D:459:VAL:HG12	4:D:463:LEU:CD2	2.50	0.42
7:2:376:ASN:O	7:2:380:THR:HG23	2.19	0.42
11:6:153:ILE:HG21	11:6:267:PHE:HE1	1.84	0.42
6:F:306:TYR:CD1	6:F:319:LEU:HD11	2.54	0.42
7:2:446:VAL:HG22	11:6:303:GLU:HB2	2.00	0.42
7:2:671:GLU:N	7:2:672:PRO:HD2	2.34	0.42
8:3:386:MET:HE2	8:3:715:VAL:N	2.34	0.42
11:6:784:ASP:OD2	11:6:795:ILE:HB	2.20	0.42
12:7:228:ARG:HG2	12:7:329:ARG:HE	1.83	0.42
12:7:658:ASP:OD1	12:7:659:TYR:N	2.53	0.42
2:B:242:THR:HG22	2:B:243:PHE:H	1.84	0.42
2:B:437:PHE:CE2	2:B:441:LEU:HD11	2.54	0.42
3:C:141:MET:O	3:C:145:ARG:HD3	2.19	0.42
3:C:219:PHE:CZ	3:C:223:LEU:HD11	2.54	0.42
7:2:311:GLU:HA	7:2:314:LEU:HD23	2.00	0.42
8:3:379:LYS:O	8:3:383:LEU:HD23	2.19	0.42
8:3:667:VAL:HG11	8:3:714:LYS:HA	2.01	0.42
10:5:45:ILE:O	10:5:48:ASP:OD1	2.38	0.42
12:7:404:LEU:C	12:7:404:LEU:HD23	2.44	0.42
13:X:64:DC:C2'	13:X:65:DT:H72	2.46	0.42
1:A:827:ASP:HA	1:A:830:LYS:HG2	2.02	0.42
2:B:241:ASP:OD2	2:B:241:ASP:O	2.38	0.42
3:C:425:ILE:HD12	3:C:425:ILE:H	1.84	0.42
6:F:304:ASP:OD1	6:F:308:ILE:HD12	2.19	0.42
7:2:657:TYR:O	7:2:657:TYR:CG	2.72	0.42
8:3:444:ALA:O	8:3:446:VAL:HG23	2.20	0.42
10:5:415:LEU:HA	10:5:523:ALA:HB3	2.02	0.42
11:6:642:ASP:HB2	11:6:683:ASN:HB3	2.01	0.42
11:6:762:LYS:O	11:6:762:LYS:CD	2.65	0.42
1:A:809:LEU:HD13	1:A:902:LYS:CD	2.50	0.42
1:A:850:PHE:HE1	1:A:863:ILE:HG22	1.84	0.42
3:C:374:ALA:HB1	3:C:377:ILE:CD1	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:296:SER:O	4:D:298:ILE:HD12	2.19	0.42
9:4:183:THR:HG22	9:4:184:ASN:N	2.35	0.42
10:5:25:THR:O	10:5:28:ILE:HG22	2.19	0.42
12:7:456:VAL:HG12	12:7:457:CYS:N	2.33	0.42
2:B:447:ILE:HG22	2:B:449:GLN:H	1.85	0.42
4:D:509:ILE:HG22	4:D:510:LEU:H	1.85	0.42
8:3:403:ILE:HG13	8:3:707:ARG:HD3	2.02	0.42
10:5:436:ALA:HB1	10:5:476:VAL:O	2.20	0.42
11:6:526:TYR:O	11:6:530:VAL:HG23	2.19	0.42
12:7:207:LEU:HD12	12:7:207:LEU:C	2.45	0.42
12:7:523:ILE:HD13	12:7:529:MET:HE1	2.01	0.42
13:X:56:DC:H2"	13:X:57:DT:H72	2.01	0.42
13:X:75:DA:H2"	13:X:76:DC:C6	2.55	0.42
2:B:212:LEU:HD23	2:B:212:LEU:HA	1.88	0.42
2:B:444:LEU:C	2:B:444:LEU:HD23	2.45	0.42
4:D:71:ILE:HG22	4:D:76:GLN:OE1	2.19	0.42
5:E:261:LEU:CD1	5:E:297:ASP:OD2	2.67	0.42
6:F:39:LEU:HD22	6:F:61:CYS:HB3	2.02	0.42
7:2:505:ILE:HD13	7:2:548:ALA:HB3	2.01	0.42
10:5:83:PRO:O	10:5:87:ILE:HD12	2.20	0.42
11:6:134:LYS:NZ	11:6:160:MET:HE1	2.34	0.42
12:7:481:VAL:HG11	12:7:512:ALA:HB1	2.02	0.42
4:D:253:LYS:HE2	4:D:255:ASN:OD1	2.20	0.41
6:F:330:ILE:HD11	6:F:383:GLN:HA	2.01	0.41
9:4:600:GLY:HA3	12:7:549:SER:O	2.20	0.41
9:4:604:TYR:HA	9:4:658:LYS:HE3	2.02	0.41
9:4:817:VAL:HG12	9:4:818:GLU:N	2.34	0.41
10:5:549:ARG:HG3	10:5:651:ARG:HH12	1.85	0.41
11:6:538:PHE:CE2	11:6:727:LEU:HD11	2.55	0.41
12:7:318:LEU:HD23	12:7:320:GLN:H	1.85	0.41
12:7:551:ALA:O	12:7:552:GLY:C	2.63	0.41
1:A:705:ARG:HA	1:A:708:LYS:HG2	2.02	0.41
4:D:289:ARG:CZ	4:D:313:LEU:HD23	2.50	0.41
7:2:671:GLU:N	7:2:672:PRO:CD	2.83	0.41
7:2:694:ARG:HG3	7:2:695:LEU:N	2.35	0.41
10:5:377:SER:C	10:5:378:ILE:HD12	2.45	0.41
11:6:179:PRO:O	11:6:182:GLN:HG2	2.20	0.41
1:A:602:MET:HE2	1:A:602:MET:HA	2.00	0.41
1:A:874:LEU:HD22	1:A:879:LEU:HB3	2.02	0.41
6:F:413:VAL:HG12	6:F:414:THR:N	2.35	0.41
11:6:646:ILE:HD12	13:X:73:DT:OP2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:299:VAL:HG23	3:C:425:ILE:CG2	2.48	0.41
4:D:283:ASP:O	4:D:286:ASP:OD1	2.38	0.41
5:E:77:ILE:O	5:E:81:VAL:HG12	2.21	0.41
6:F:313:LEU:HD21	6:F:319:LEU:HD21	2.00	0.41
7:2:206:THR:HA	7:2:209:ARG:HG2	2.01	0.41
7:2:494:ILE:HG23	7:2:823:MET:HE1	2.03	0.41
8:3:279:ASP:O	8:3:280:ASP:C	2.62	0.41
8:3:306:MET:HE1	10:5:179:LEU:CD1	2.51	0.41
12:7:539:GLU:OE1	12:7:545:THR:O	2.38	0.41
2:B:269:VAL:HG12	2:B:270:THR:H	1.85	0.41
3:C:580:ALA:O	3:C:583:LEU:HG	2.21	0.41
7:2:218:TYR:CE1	7:2:226:VAL:HG22	2.55	0.41
9:4:276:ILE:HG23	9:4:277:LYS:N	2.36	0.41
12:7:245:ILE:HD11	12:7:349:VAL:HG11	2.02	0.41
14:Y:12:DA:H2''	14:Y:13:DG:C8	2.55	0.41
3:C:244:LEU:HG	3:C:245:SER:N	2.35	0.41
4:D:284:MET:SD	4:D:324:ILE:HG23	2.60	0.41
4:D:450:LEU:HD23	4:D:450:LEU:H	1.86	0.41
4:D:493:GLN:O	4:D:497:ILE:HG12	2.21	0.41
5:E:210:MET:HE1	5:E:232:PHE:O	2.20	0.41
8:3:212:ARG:NH2	8:3:229:ALA:HB1	2.36	0.41
9:4:589:VAL:HG12	9:4:590:TYR:N	2.35	0.41
10:5:360:LEU:HD22	10:5:594:ILE:HD11	2.03	0.41
3:C:298:PHE:HE2	3:C:475:LEU:HD11	1.86	0.41
5:E:103:GLU:H	5:E:107:LEU:HD12	1.86	0.41
6:F:17:ASP:HB3	6:F:18:PRO:HD3	2.03	0.41
9:4:293:LEU:O	9:4:293:LEU:HD23	2.20	0.41
13:X:60:DC:H2''	13:X:61:DT:C6	2.55	0.41
14:Y:55:DA:C4	14:Y:56:DA:C8	3.09	0.41
1:A:685:SER:OG	1:A:688:ALA:HB3	2.21	0.41
2:B:244:GLU:OE1	5:E:353:LYS:HD3	2.21	0.41
2:B:301:PHE:N	2:B:302:PRO:HD2	2.35	0.41
8:3:414:ALA:O	8:3:415:LYS:C	2.63	0.41
11:6:448:LEU:O	11:6:448:LEU:HD12	2.21	0.41
12:7:236:GLY:N	12:7:355:PHE:O	2.54	0.41
14:Y:31:DC:H4'	14:Y:32:DA:OP1	2.20	0.41
3:C:328:TYR:HD1	3:C:468:LEU:HD21	1.86	0.41
5:E:225:THR:O	5:E:225:THR:HG23	2.21	0.41
5:E:274:THR:H	5:E:277:ASN:HB2	1.85	0.41
8:3:443:THR:O	8:3:444:ALA:HB3	2.21	0.41
8:3:523:TYR:CD2	8:3:548:VAL:HG11	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:4:692:ILE:O	9:4:692:ILE:HG22	2.20	0.41
10:5:136:GLN:NE2	10:5:279:ASP:OD1	2.54	0.41
12:7:25:LEU:HD13	12:7:121:ILE:HD13	2.03	0.41
12:7:227:VAL:HG12	12:7:227:VAL:O	2.21	0.41
12:7:428:VAL:HG13	12:7:458:LEU:HD21	2.03	0.41
12:7:625:GLN:HB3	12:7:626:PRO:HD2	2.02	0.41
3:C:73:ASP:HA	3:C:76:ILE:HG22	2.02	0.41
3:C:346:LEU:HD21	3:C:351:LEU:CD2	2.51	0.41
4:D:219:ILE:HD11	4:D:231:LEU:HD22	2.03	0.41
6:F:411:ILE:H	6:F:411:ILE:HD12	1.86	0.41
9:4:306:TYR:HB3	9:4:465:HIS:ND1	2.36	0.41
9:4:543:GLN:OE1	9:4:562:ILE:O	2.39	0.41
11:6:297:THR:HG23	11:6:391:PRO:HG3	2.03	0.41
11:6:400:VAL:HG12	11:6:401:GLU:N	2.35	0.41
11:6:635:ILE:HG22	11:6:636:CYS:N	2.36	0.41
12:7:370:LEU:N	12:7:370:LEU:HD12	2.36	0.41
3:C:527:LEU:HA	3:C:530:GLU:HG2	2.03	0.40
4:D:143:ILE:HD11	4:D:181:PHE:HA	2.03	0.40
6:F:283:GLU:HA	6:F:286:ILE:HB	2.02	0.40
10:5:87:ILE:N	10:5:88:PRO:HD2	2.36	0.40
10:5:98:ALA:HB2	10:5:135:PHE:CE2	2.56	0.40
13:X:35:DT:C2'	13:X:36:DT:H71	2.51	0.40
3:C:42:VAL:HG12	3:C:43:LYS:N	2.36	0.40
3:C:487:ILE:O	3:C:491:LEU:HG	2.21	0.40
8:3:544:ASP:CG	8:3:704:THR:HG22	2.46	0.40
8:3:667:VAL:O	8:3:668:ILE:HD13	2.22	0.40
9:4:262:LEU:HD23	9:4:263:ASN:N	2.37	0.40
10:5:41:ASP:CG	10:5:41:ASP:O	2.64	0.40
10:5:526:ILE:HG13	10:5:526:ILE:O	2.20	0.40
10:5:559:ASP:O	10:5:560:HIS:C	2.64	0.40
11:6:568:ASP:OD2	11:6:677:SER:HB3	2.21	0.40
13:X:73:DT:H2''	13:X:74:DG:C8	2.56	0.40
3:C:116:ILE:HG22	3:C:117:GLU:N	2.37	0.40
6:F:316:PRO:O	6:F:319:LEU:HG	2.21	0.40
9:4:486:MET:C	9:4:486:MET:SD	3.04	0.40
9:4:531:TYR:CZ	9:4:532:GLU:OE1	2.74	0.40
10:5:430:GLU:OE1	10:5:436:ALA:HB3	2.21	0.40
11:6:322:GLU:HG3	11:6:326:LYS:O	2.21	0.40
12:7:228:ARG:NH2	12:7:322:VAL:HA	2.37	0.40
1:A:613:ILE:H	1:A:613:ILE:HD12	1.87	0.40
1:A:790:VAL:O	1:A:794:MET:HG2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:903:ARG:O	1:A:907:GLU:CD	2.64	0.40
2:B:206:SER:OG	2:B:209:LYS:HG2	2.21	0.40
7:2:621:HIS:HA	7:2:673:ILE:HD12	2.04	0.40
7:2:671:GLU:HG2	7:2:672:PRO:N	2.36	0.40
7:2:806:THR:CG2	7:2:807:VAL:N	2.85	0.40
8:3:221:LEU:HD12	8:3:221:LEU:H	1.86	0.40
8:3:703:GLU:O	8:3:707:ARG:HG3	2.21	0.40
9:4:446:ALA:HB3	9:4:452:VAL:HG23	2.04	0.40
10:5:400:LEU:HB3	10:5:401:PRO:HD2	2.04	0.40
10:5:549:ARG:HG3	10:5:651:ARG:NH1	2.36	0.40
11:6:785:ALA:O	11:6:786:GLN:HB3	2.22	0.40
6:F:409:THR:HG23	6:F:409:THR:O	2.22	0.40
7:2:245:ASN:OD1	7:2:245:ASN:O	2.39	0.40
7:2:637:VAL:O	7:2:637:VAL:HG23	2.21	0.40
9:4:342:MET:HG2	11:6:448:LEU:CD1	2.52	0.40
9:4:435:VAL:HG22	9:4:436:THR:N	2.37	0.40
10:5:340:SER:O	10:5:341:SER:C	2.65	0.40
10:5:470:VAL:HG21	10:5:513:LEU:HD12	2.04	0.40
12:7:132:ILE:HD13	12:7:144:ASN:OD1	2.22	0.40
13:X:40:DT:H2'	13:X:41:DT:H71	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	A	486/949 (51%)	464 (96%)	22 (4%)	0	100	100
2	B	283/620 (46%)	260 (92%)	23 (8%)	0	100	100
3	C	578/616 (94%)	535 (93%)	43 (7%)	0	100	100
4	D	429/529 (81%)	403 (94%)	26 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	E	455/479 (95%)	434 (95%)	21 (5%)	0	100	100
6	F	234/435 (54%)	217 (93%)	17 (7%)	0	100	100
7	2	595/868 (68%)	547 (92%)	48 (8%)	0	100	100
8	3	589/1006 (58%)	551 (94%)	38 (6%)	0	100	100
9	4	635/933 (68%)	577 (91%)	58 (9%)	0	100	100
10	5	590/775 (76%)	539 (91%)	51 (9%)	0	100	100
11	6	608/1017 (60%)	556 (91%)	52 (9%)	0	100	100
12	7	646/845 (76%)	604 (94%)	42 (6%)	0	100	100
All	All	6128/9072 (68%)	5687 (93%)	441 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	438/842 (52%)	437 (100%)	1 (0%)	87	85
2	B	271/573 (47%)	271 (100%)	0	100	100
3	C	543/576 (94%)	543 (100%)	0	100	100
4	D	404/488 (83%)	404 (100%)	0	100	100
5	E	423/440 (96%)	423 (100%)	0	100	100
6	F	228/406 (56%)	228 (100%)	0	100	100
7	2	526/770 (68%)	526 (100%)	0	100	100
8	3	518/864 (60%)	518 (100%)	0	100	100
9	4	582/848 (69%)	582 (100%)	0	100	100
10	5	539/688 (78%)	539 (100%)	0	100	100
11	6	537/886 (61%)	537 (100%)	0	100	100
12	7	575/753 (76%)	575 (100%)	0	100	100
All	All	5584/8134 (69%)	5583 (100%)	1 (0%)	100	100

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

Mol	Chain	Res	Type
1	A	906	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	364	ASN
1	A	820	GLN
1	A	872	GLN
2	B	256	ASN
2	B	279	ASN
2	B	286	GLN
2	B	458	HIS
3	C	67	HIS
3	C	278	ASN
3	C	308	ASN
3	C	477	GLN
4	D	53	GLN
4	D	491	GLN
5	E	248	HIS
5	E	253	ASN
5	E	422	GLN
7	2	356	ASN
7	2	405	HIS
7	2	508	HIS
7	2	531	HIS
7	2	653	ASN
7	2	779	HIS
7	2	849	GLN
8	3	46	GLN
8	3	175	HIS
9	4	242	ASN
9	4	287	ASN
9	4	320	ASN
9	4	386	HIS
9	4	450	GLN
9	4	676	ASN
9	4	735	HIS
9	4	797	GLN
9	4	815	ASN
10	5	136	GLN
10	5	155	HIS

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Mol	Chain	Res	Type
10	5	188	HIS
10	5	344	ASN
10	5	574	ASN
10	5	579	ASN
10	5	587	GLN
10	5	652	GLN
11	6	182	GLN
11	6	633	ASN
11	6	653	HIS
12	7	124	ASN
12	7	144	ASN
12	7	384	HIS
12	7	568	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 8 are monoatomic - leaving 7 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	$\# Z > 2$
17	ADP	5	801	-	28,29,29	1.39	4 (14%)	43,45,45	1.90	10 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
15	ATP	A	2001	16	32,33,33	0.28	0	48,52,52	0.70	0
17	ADP	2	901	-	28,29,29	1.39	4 (14%)	43,45,45	1.91	10 (23%)
17	ADP	3	2001	-	28,29,29	1.39	4 (14%)	43,45,45	1.90	11 (25%)
15	ATP	D	2001	16	32,33,33	0.33	0	48,52,52	0.73	0
17	ADP	7	901	-	28,29,29	1.37	4 (14%)	43,45,45	1.89	11 (25%)
15	ATP	E	2001	16	32,33,33	0.30	0	48,52,52	0.72	0

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
17	ADP	5	801	-	-	2/16/32/32	0/3/3/3
15	ATP	A	2001	16	-	4/22/38/38	0/3/3/3
17	ADP	2	901	-	-	3/16/32/32	0/3/3/3
17	ADP	3	2001	-	-	4/16/32/32	0/3/3/3
15	ATP	D	2001	16	-	2/22/38/38	0/3/3/3
17	ADP	7	901	-	-	4/16/32/32	0/3/3/3
15	ATP	E	2001	16	-	1/22/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	2	901	ADP	C5-C4	4.62	1.47	1.39
17	5	801	ADP	C5-C4	4.60	1.47	1.39
17	3	2001	ADP	C5-C4	4.51	1.47	1.39
17	7	901	ADP	C5-C4	4.49	1.47	1.39
17	3	2001	ADP	C5-C6	2.68	1.48	1.41
17	5	801	ADP	C5-C6	2.64	1.48	1.41
17	2	901	ADP	C5-C6	2.63	1.48	1.41
17	7	901	ADP	C5-C6	2.53	1.48	1.41
17	3	2001	ADP	C5-N7	-2.46	1.34	1.39
17	5	801	ADP	C5-N7	-2.44	1.34	1.39
17	7	901	ADP	C5-N7	-2.38	1.34	1.39
17	2	901	ADP	C5-N7	-2.35	1.34	1.39
17	2	901	ADP	C8-N7	2.30	1.36	1.31
17	3	2001	ADP	C8-N7	2.29	1.36	1.31
17	7	901	ADP	C8-N7	2.25	1.36	1.31
17	5	801	ADP	C8-N7	2.21	1.35	1.31

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	5	801	ADP	C5-C4-N3	-6.03	118.41	126.72
17	3	2001	ADP	C5-C4-N3	-5.86	118.65	126.72
17	2	901	ADP	C5-C4-N3	-5.80	118.73	126.72
17	7	901	ADP	C5-C4-N3	-5.65	118.94	126.72
17	5	801	ADP	N3-C4-N9	4.95	135.59	127.17
17	2	901	ADP	N3-C4-N9	4.80	135.34	127.17
17	7	901	ADP	N3-C4-N9	4.65	135.08	127.17
17	3	2001	ADP	N3-C4-N9	4.65	135.07	127.17
17	3	2001	ADP	C2-N3-C4	3.84	121.22	111.83
17	5	801	ADP	C2-N3-C4	3.82	121.16	111.83
17	2	901	ADP	C2-N3-C4	3.81	121.14	111.83
17	7	901	ADP	C2-N3-C4	3.75	120.99	111.83
17	2	901	ADP	N3-C2-N1	-3.64	123.07	128.58
17	7	901	ADP	N3-C2-N1	-3.64	123.08	128.58
17	3	2001	ADP	N3-C2-N1	-3.60	123.14	128.58
17	3	2001	ADP	C4-C5-N7	-3.54	106.53	110.58
17	5	801	ADP	N3-C2-N1	-3.53	123.23	128.58
17	5	801	ADP	C4-C5-N7	-3.40	106.69	110.58
17	7	901	ADP	C4-C5-N7	-3.35	106.75	110.58
17	2	901	ADP	C4-C5-N7	-3.33	106.78	110.58
17	7	901	ADP	C4-N9-C8	3.14	109.03	105.74
17	2	901	ADP	C4-N9-C8	3.10	109.00	105.74
17	3	2001	ADP	C4-N9-C8	2.91	108.79	105.74
17	5	801	ADP	C4-N9-C8	2.87	108.75	105.74
17	5	801	ADP	C5-N7-C8	2.80	107.84	103.45
17	3	2001	ADP	C5-N7-C8	2.78	107.82	103.45
17	7	901	ADP	C5-N7-C8	2.67	107.64	103.45
17	2	901	ADP	C5-N7-C8	2.63	107.59	103.45
17	2	901	ADP	C3'-C2'-C1'	2.52	106.23	101.46
17	5	801	ADP	C3'-C2'-C1'	2.41	106.02	101.46
17	7	901	ADP	C3'-C2'-C1'	2.38	105.96	101.46
17	3	2001	ADP	N9-C8-N7	-2.36	110.58	113.94
17	7	901	ADP	N9-C8-N7	-2.35	110.60	113.94
17	3	2001	ADP	C3'-C2'-C1'	2.32	105.86	101.46
17	2	901	ADP	N9-C8-N7	-2.28	110.70	113.94
17	5	801	ADP	N9-C8-N7	-2.28	110.71	113.94
17	3	2001	ADP	C6-C5-N7	2.12	136.18	132.09
17	2	901	ADP	C2-N1-C6	2.06	122.11	118.73
17	7	901	ADP	C2-N1-C6	2.06	122.11	118.73
17	7	901	ADP	C6-C5-N7	2.04	136.02	132.09
17	5	801	ADP	C2-N1-C6	2.03	122.06	118.73
17	3	2001	ADP	C2-N1-C6	2.01	122.04	118.73

There are no chirality outliers.

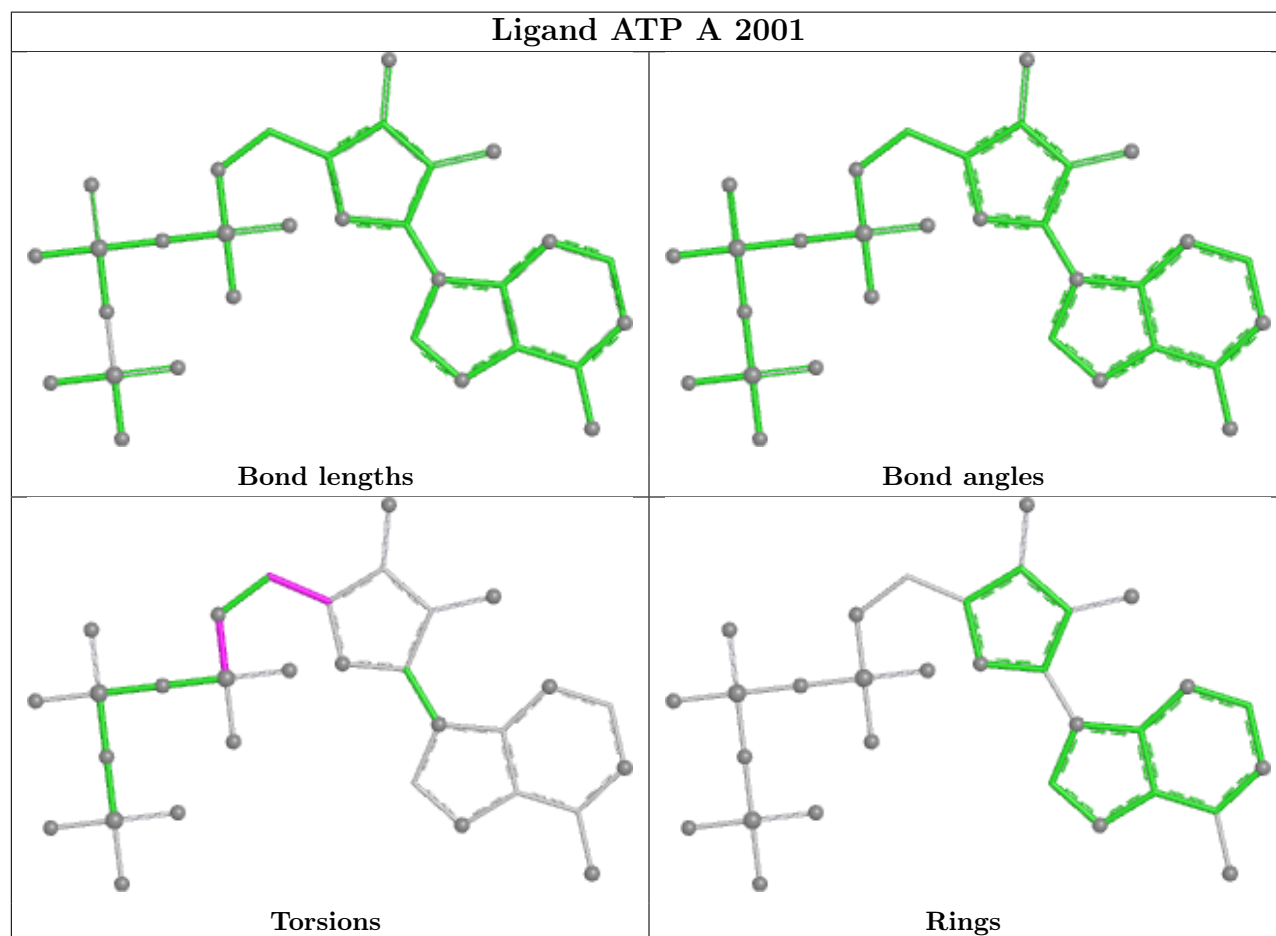
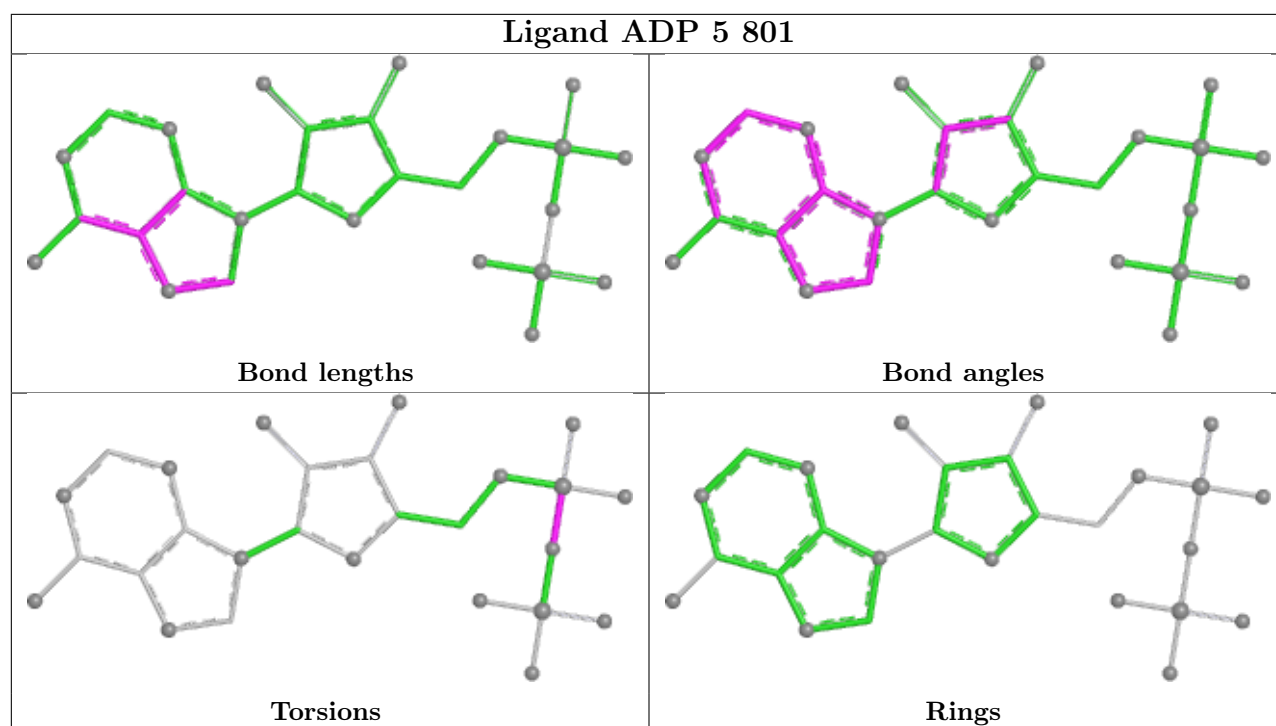
All (20) torsion outliers are listed below:

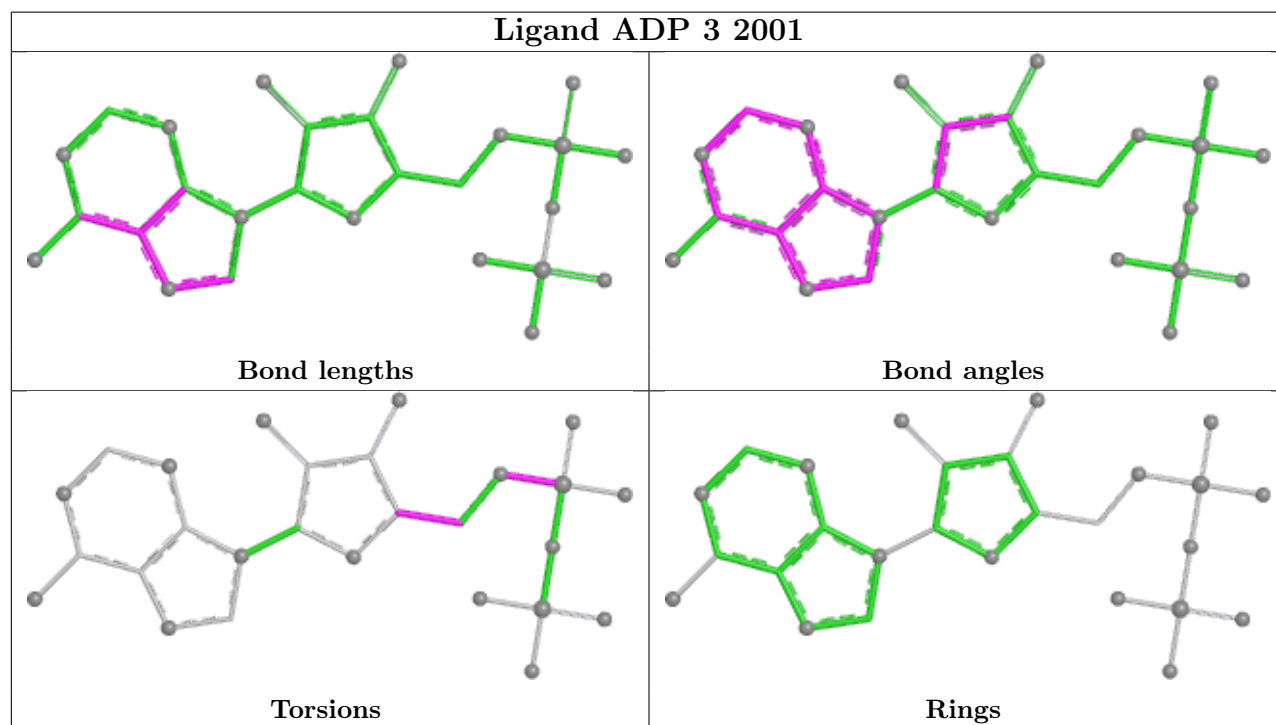
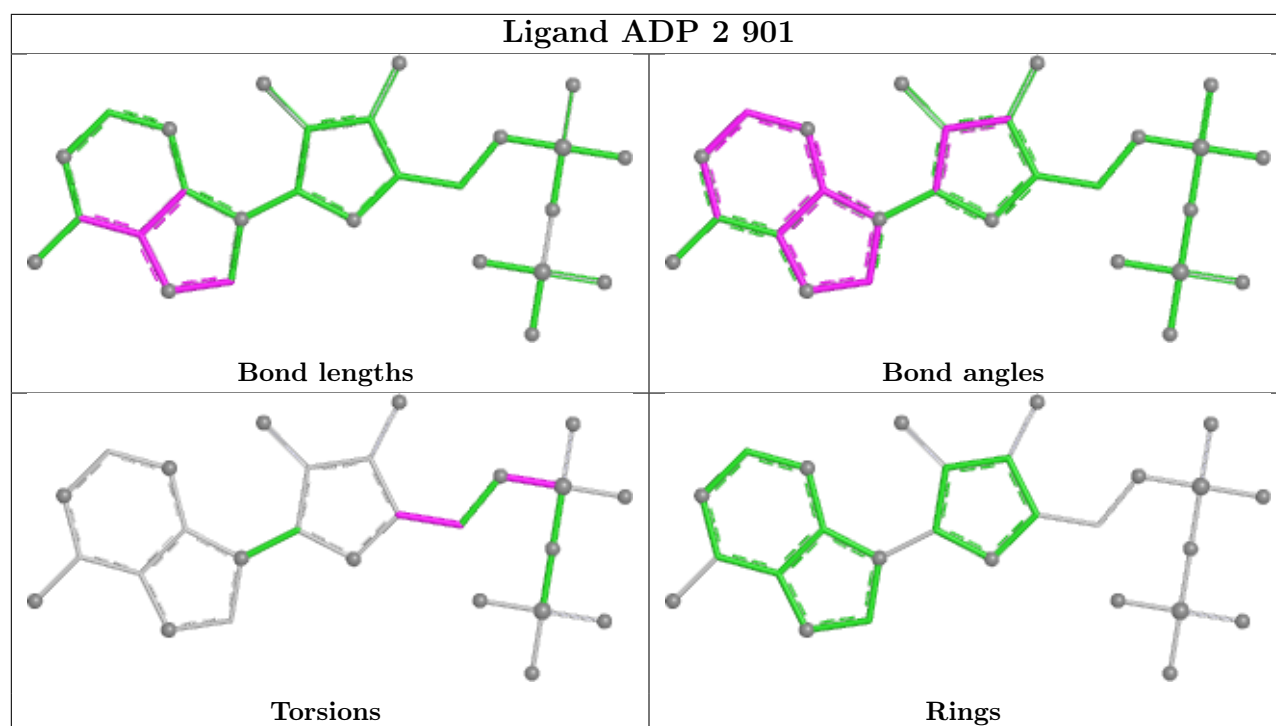
Mol	Chain	Res	Type	Atoms
15	A	2001	ATP	C5'-O5'-PA-O1A
15	A	2001	ATP	C5'-O5'-PA-O2A
15	A	2001	ATP	C5'-O5'-PA-O3A
17	2	901	ADP	C5'-O5'-PA-O2A
17	3	2001	ADP	C5'-O5'-PA-O2A
17	3	2001	ADP	C5'-O5'-PA-O3A
17	7	901	ADP	C5'-O5'-PA-O1A
17	7	901	ADP	C5'-O5'-PA-O2A
17	7	901	ADP	C5'-O5'-PA-O3A
17	3	2001	ADP	O4'-C4'-C5'-O5'
15	A	2001	ATP	O4'-C4'-C5'-O5'
15	D	2001	ATP	C3'-C4'-C5'-O5'
17	3	2001	ADP	C3'-C4'-C5'-O5'
17	5	801	ADP	PB-O3A-PA-O1A
17	7	901	ADP	O4'-C4'-C5'-O5'
17	2	901	ADP	C5'-O5'-PA-O3A
17	5	801	ADP	PB-O3A-PA-O2A
15	D	2001	ATP	O4'-C4'-C5'-O5'
15	E	2001	ATP	PB-O3A-PA-O2A
17	2	901	ADP	O4'-C4'-C5'-O5'

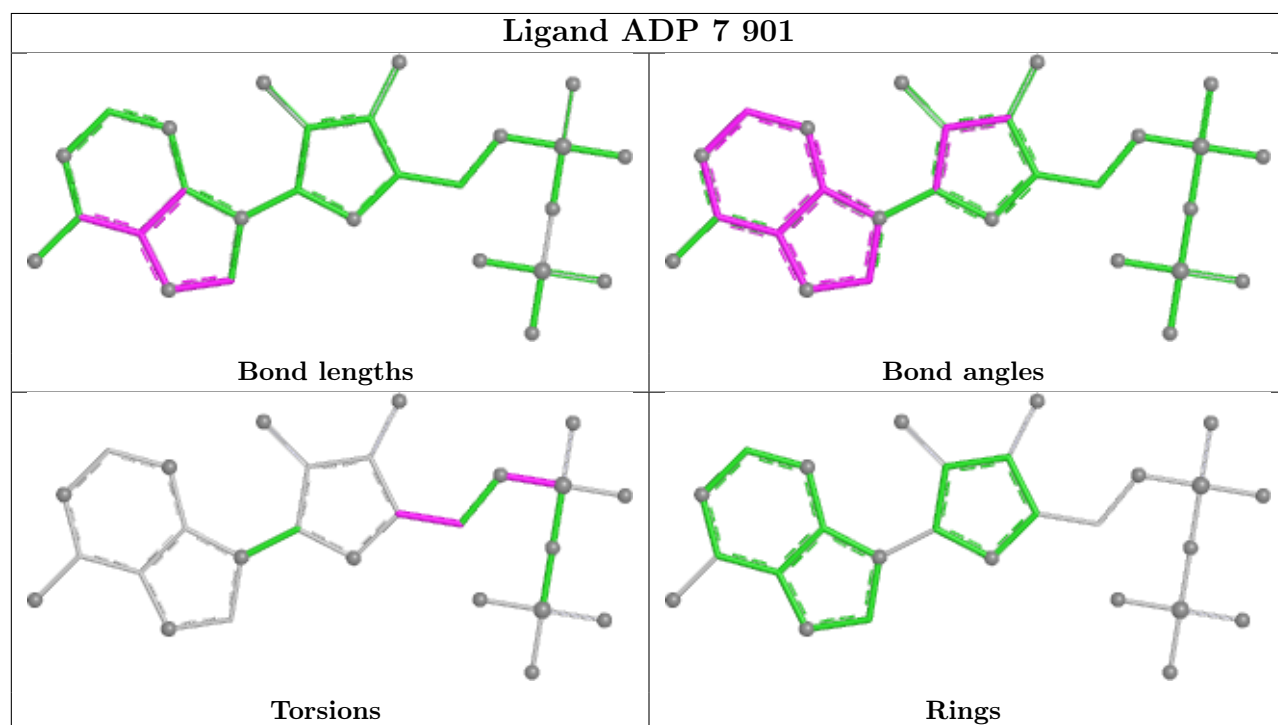
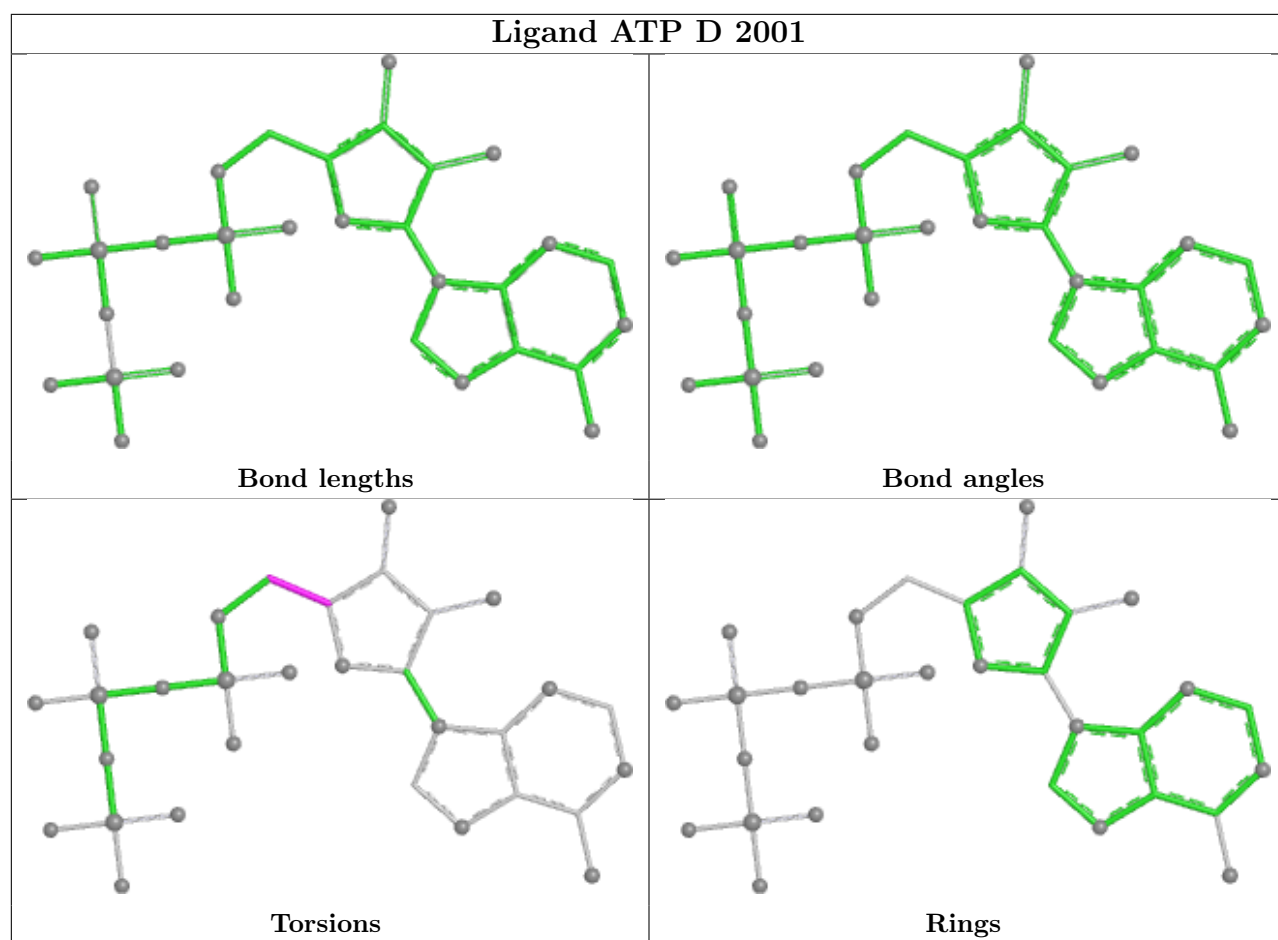
There are no ring outliers.

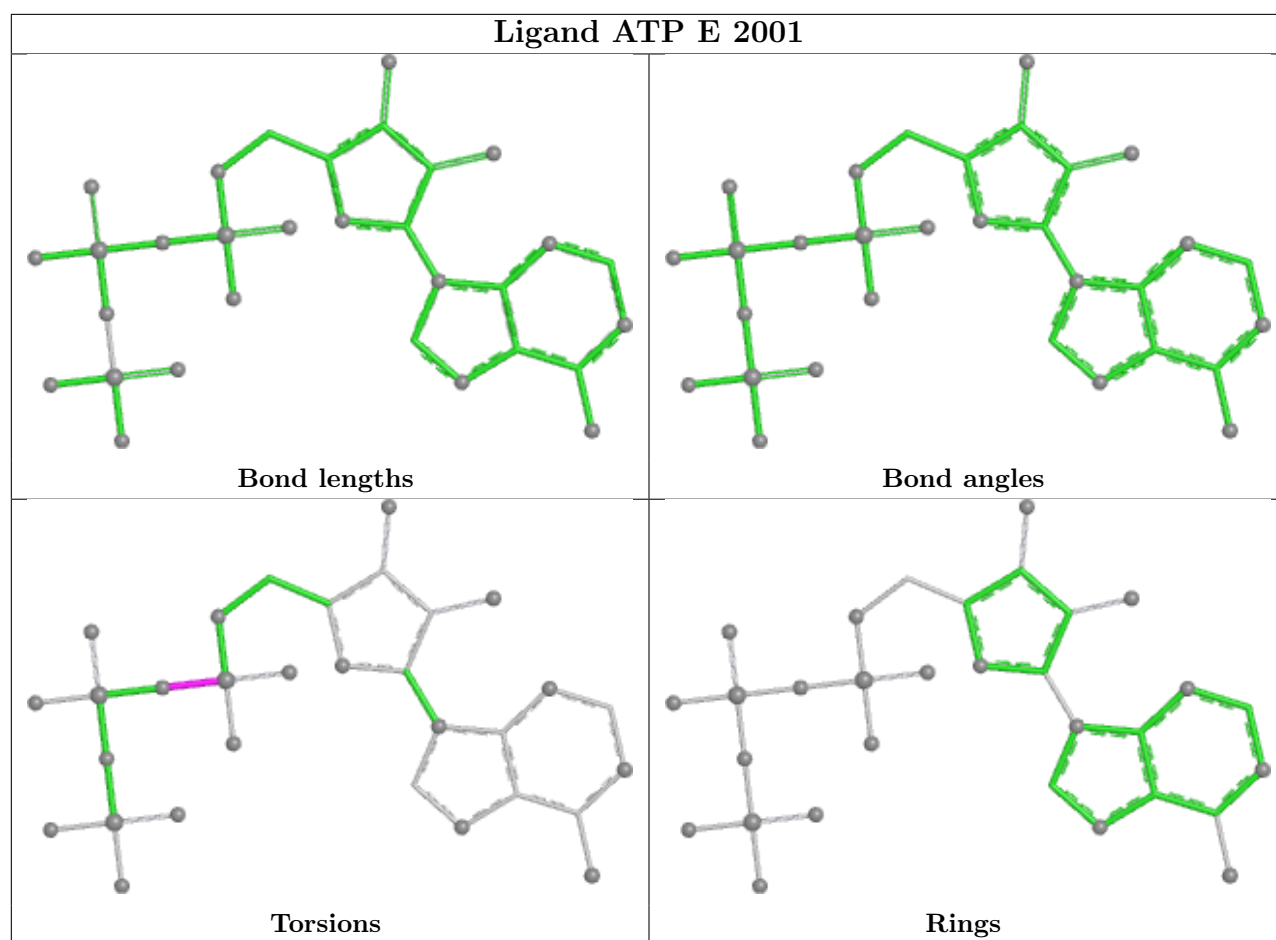
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









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.

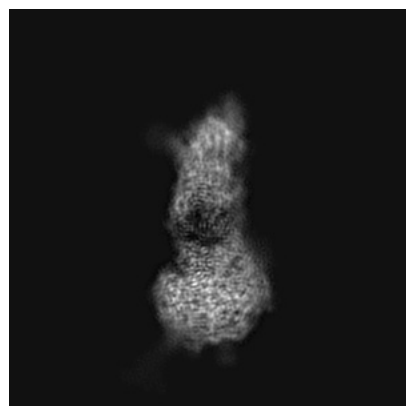
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4980. These allow visual inspection of the internal detail of the map and identification of artifacts.

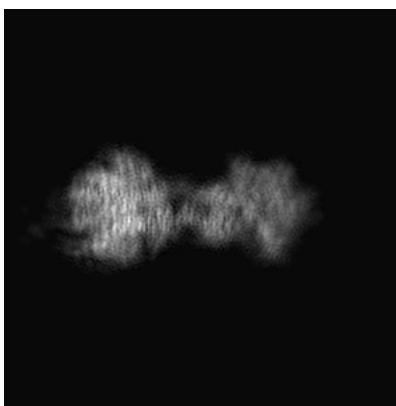
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

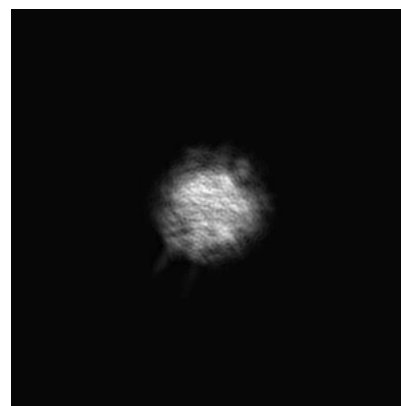
6.1.1 Primary map



X

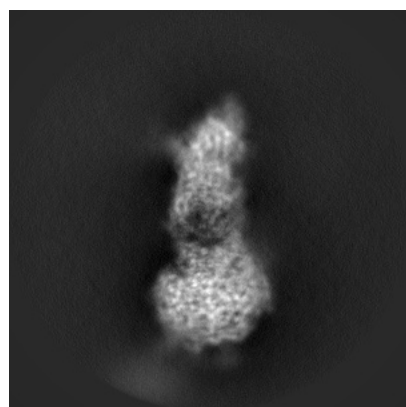


Y

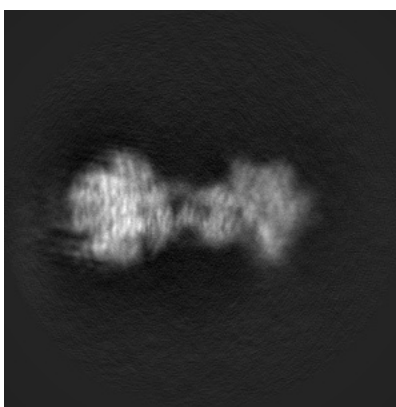


Z

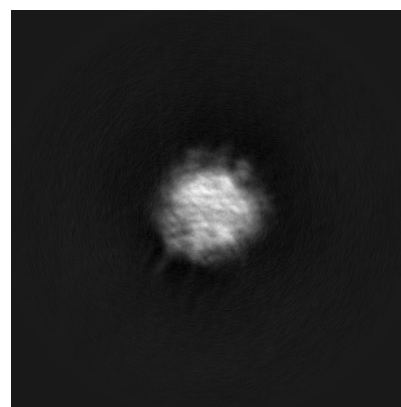
6.1.2 Raw map



X



Y

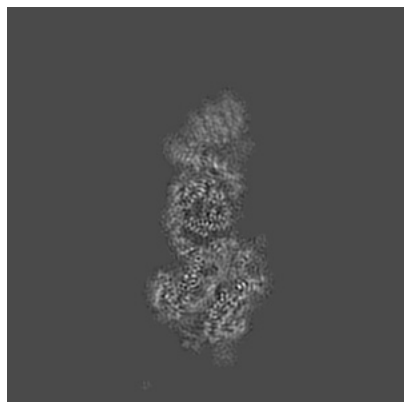


Z

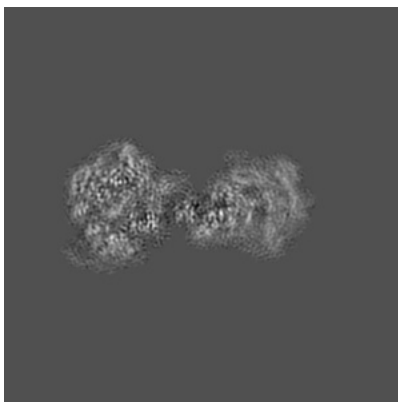
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

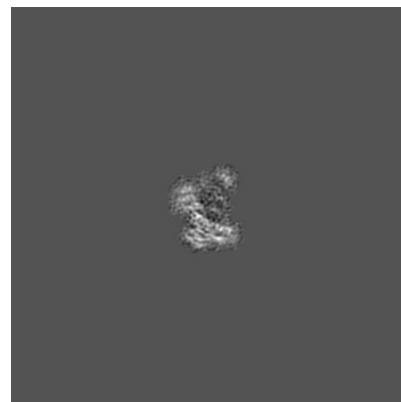
6.2.1 Primary map



X Index: 170

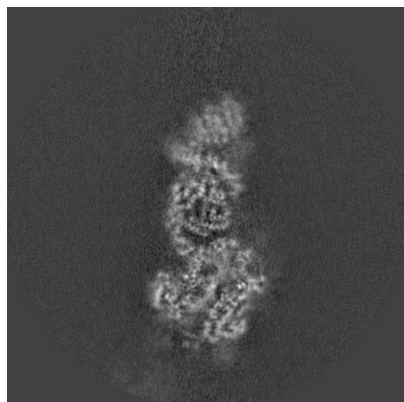


Y Index: 170

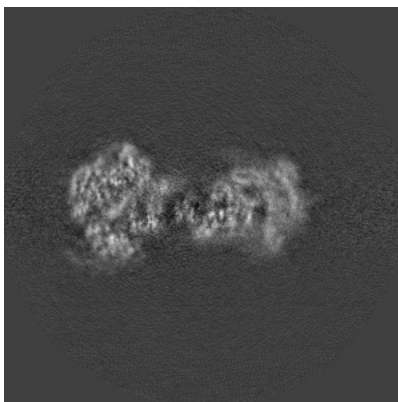


Z Index: 170

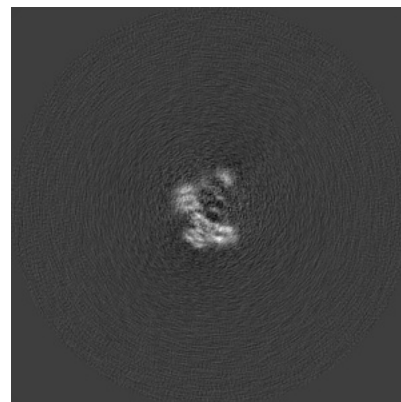
6.2.2 Raw map



X Index: 170



Y Index: 170

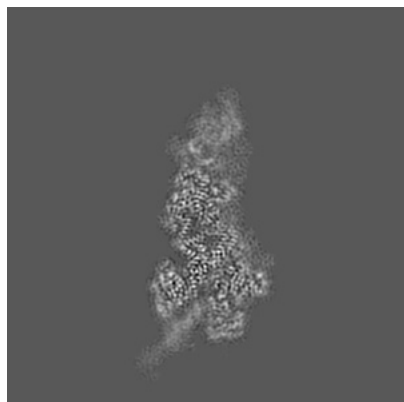


Z Index: 170

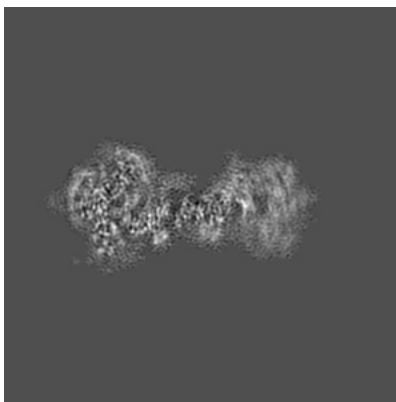
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

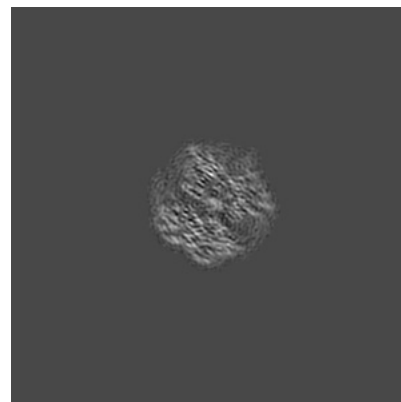
6.3.1 Primary map



X Index: 159

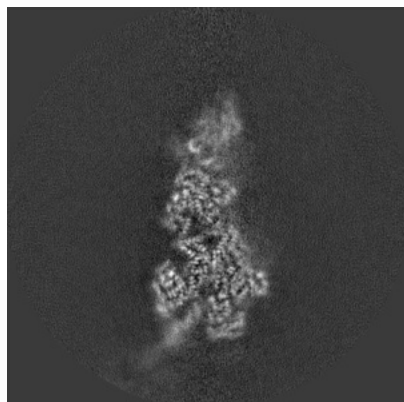


Y Index: 177

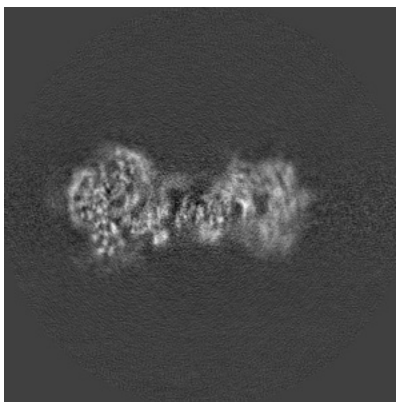


Z Index: 102

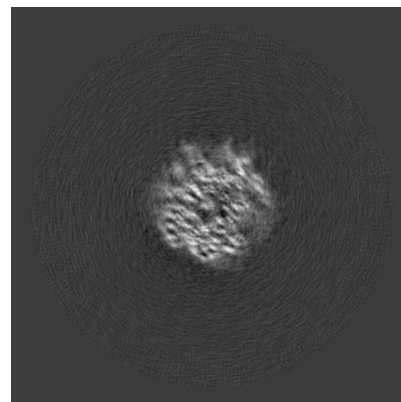
6.3.2 Raw map



X Index: 158



Y Index: 177

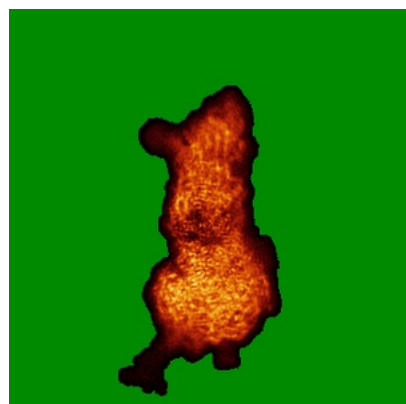


Z Index: 94

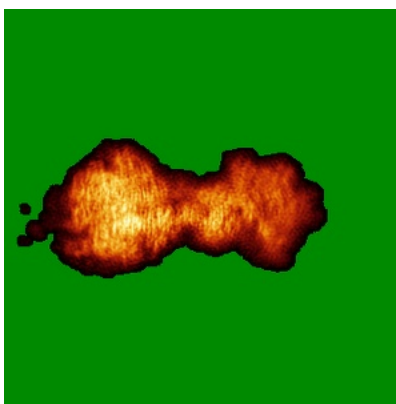
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

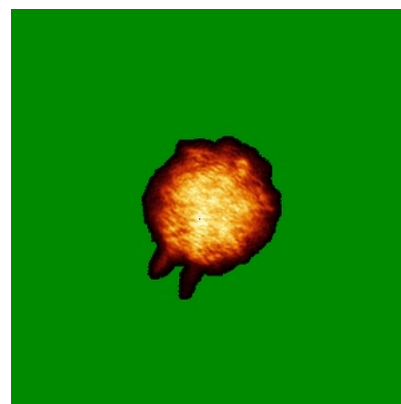
6.4.1 Primary map



X

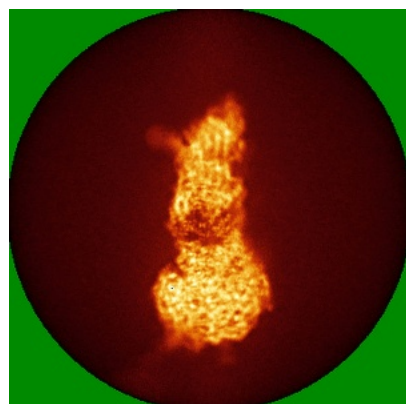


Y

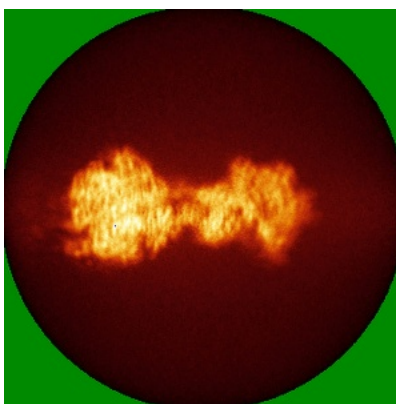


Z

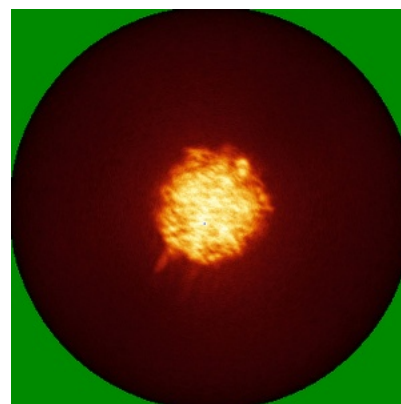
6.4.2 Raw map



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



X



Y



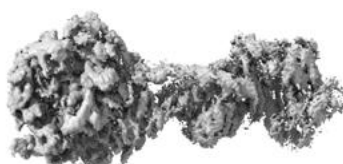
Z

The images above show the 3D surface view of the map at the recommended contour level 0.016. 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



X



Y



Z

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

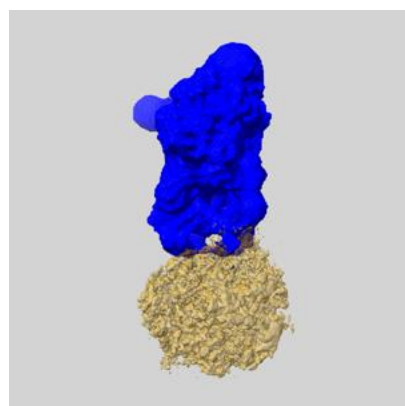
6.6 Mask visualisation [i](#)

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

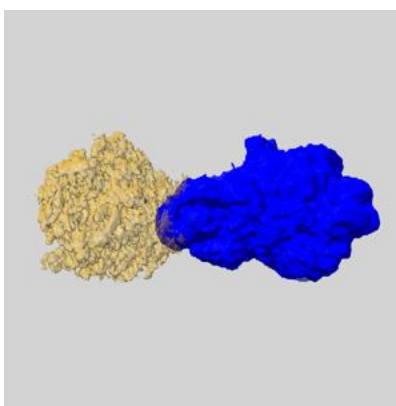
A mask typically either:

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

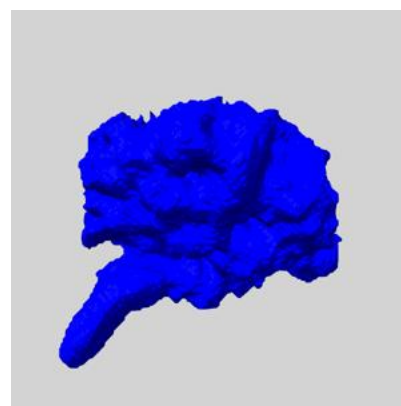
6.6.1 emd_4980_msk_2.map [i](#)



X

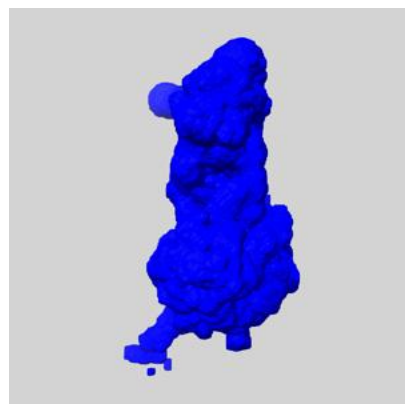


Y

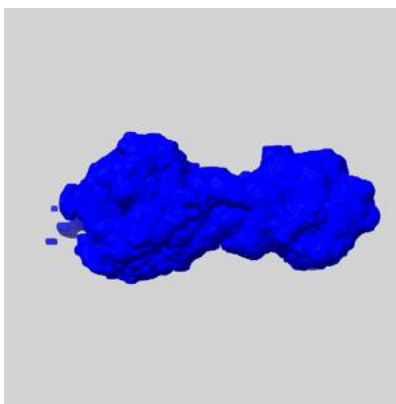


Z

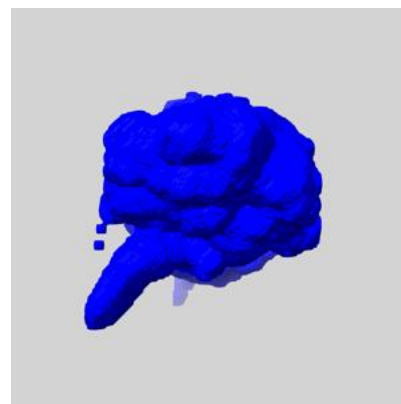
6.6.2 emd_4980_msk_3.map [i](#)



X

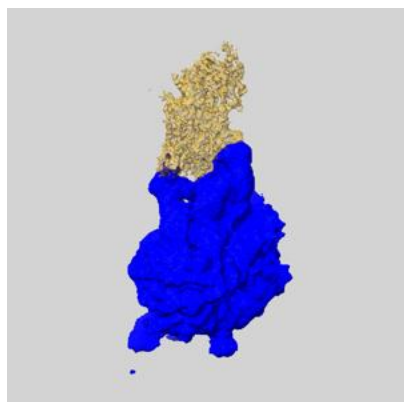


Y

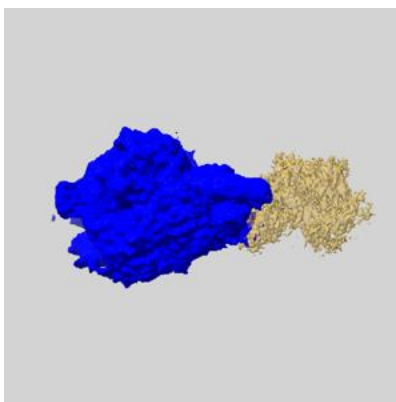


Z

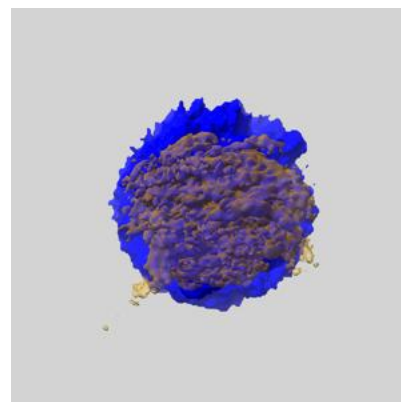
6.6.3 emd_4980_msk_1.map [i](#)



X



Y

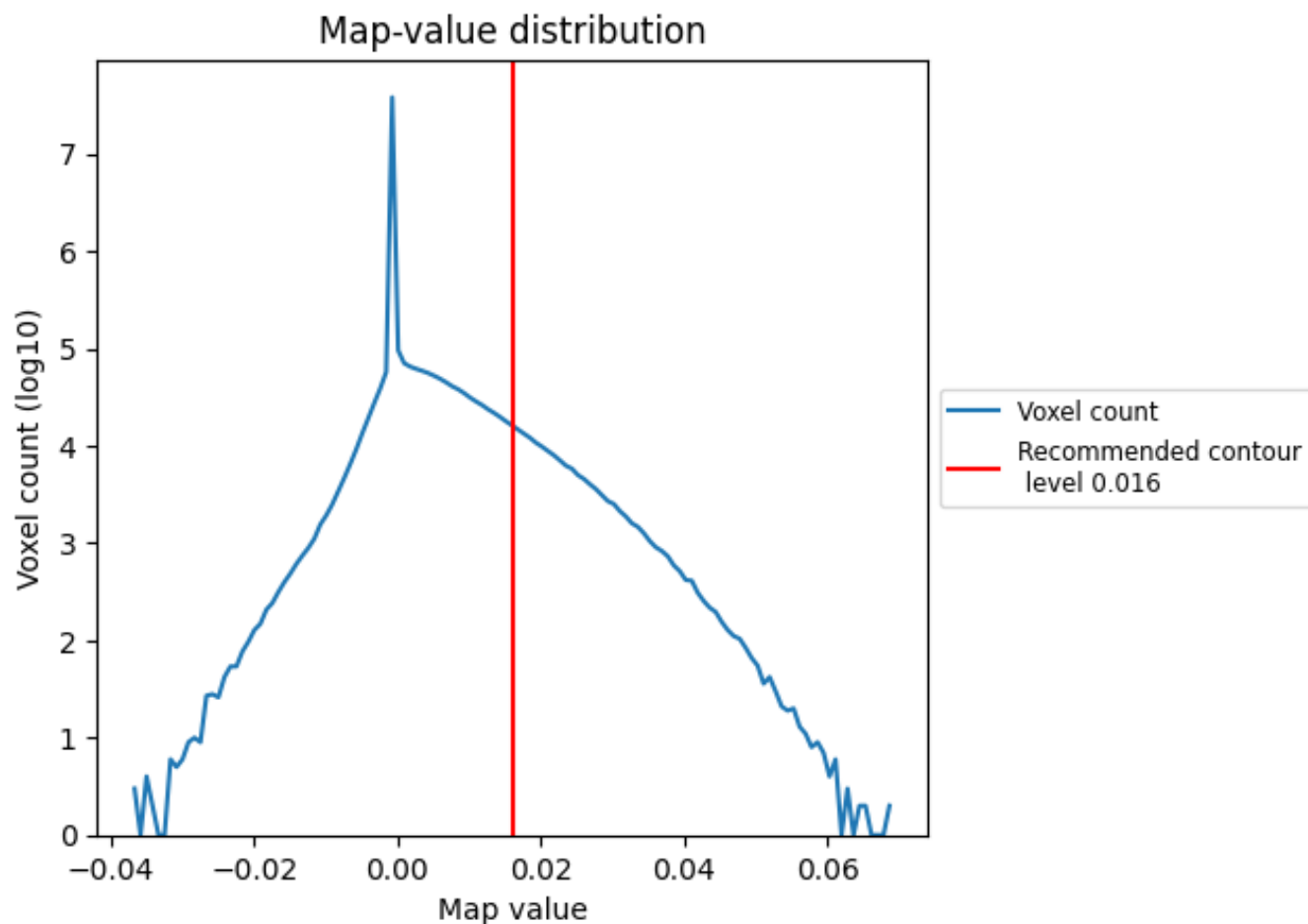


Z

7 Map analysis [i](#)

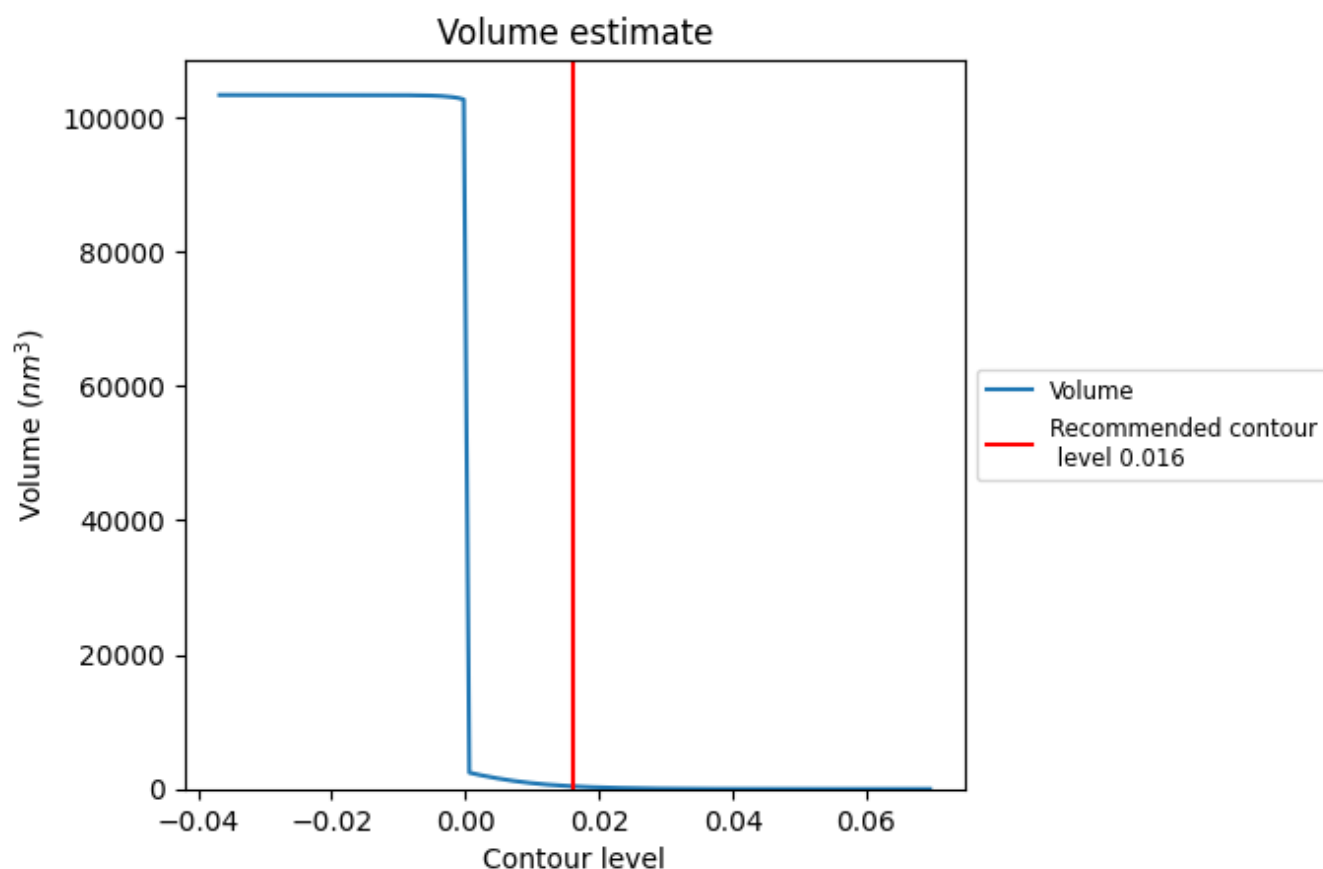
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

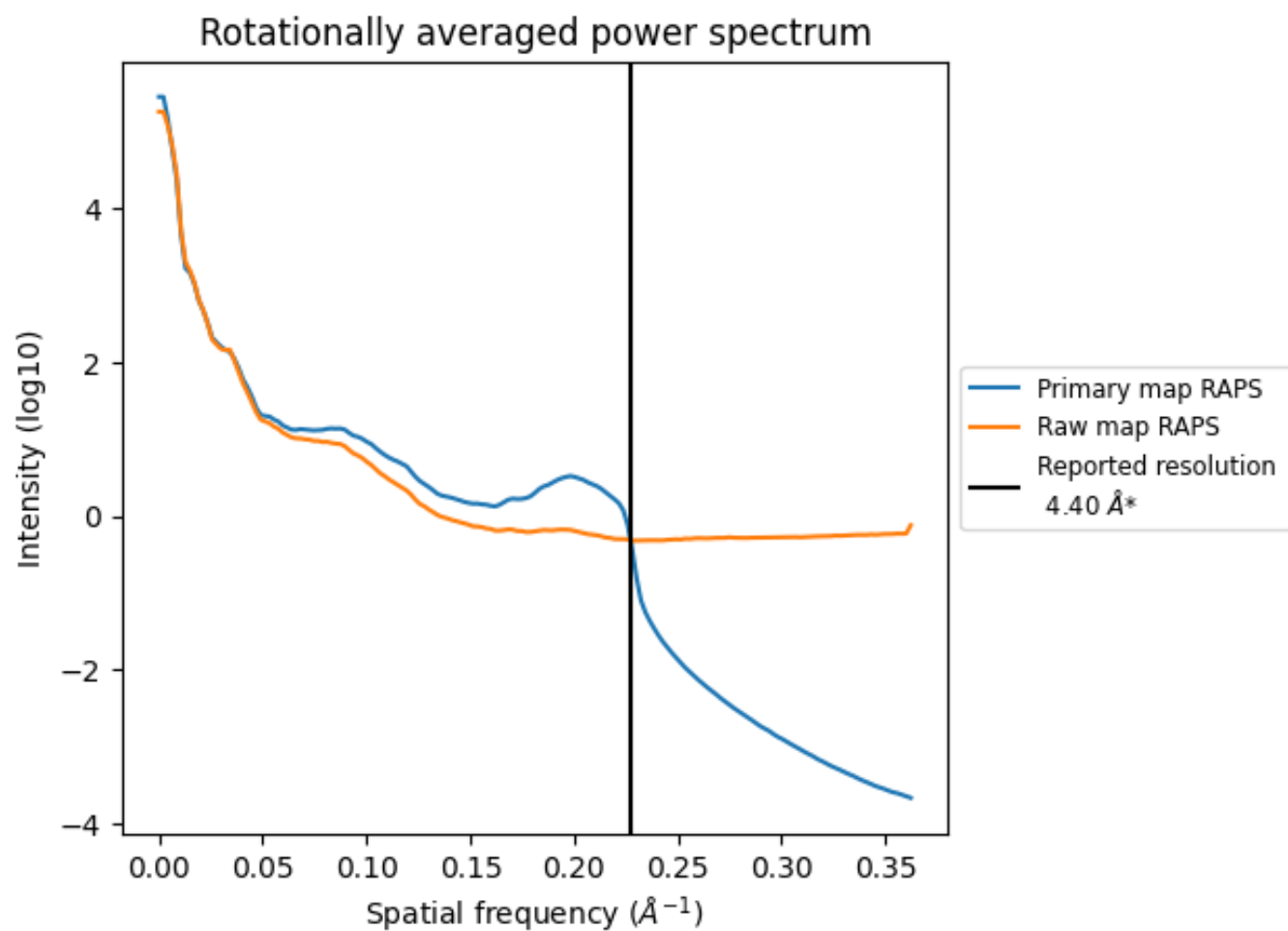
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 404 nm^3 ; this corresponds to an approximate mass of 365 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 ⓘ

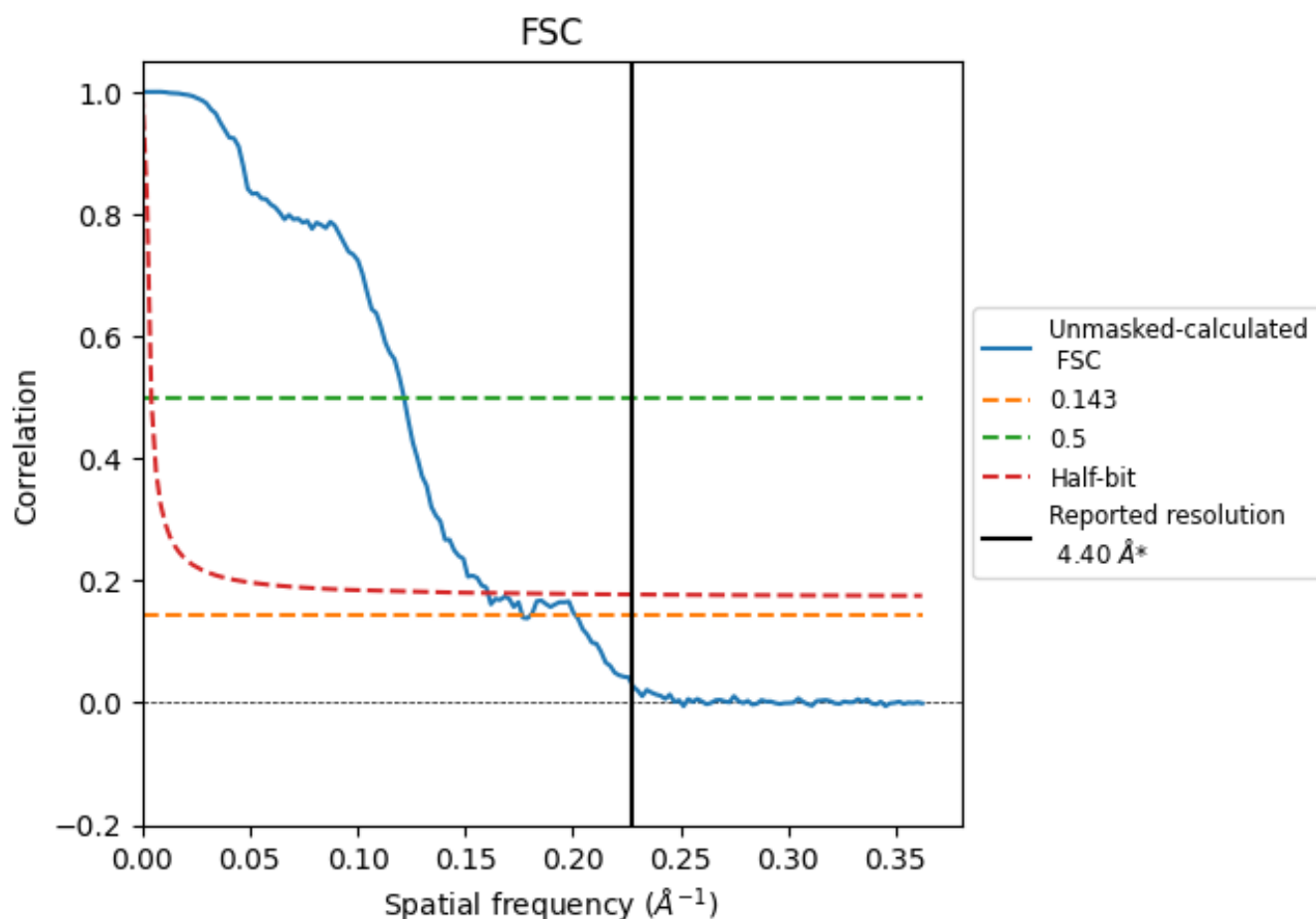


*Reported resolution corresponds to spatial frequency of 0.227 Å⁻¹

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 0.227 Å⁻¹

8.2 Resolution estimates [i](#)

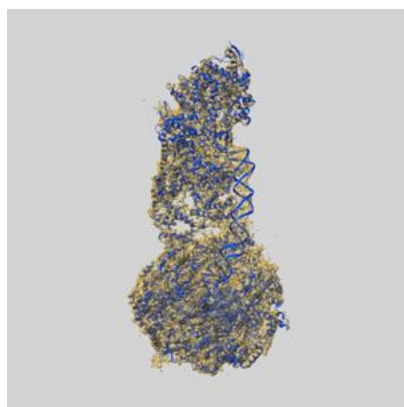
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.40	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	5.66	8.22	6.23

*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 5.66 differs from the reported value 4.4 by more than 10 %

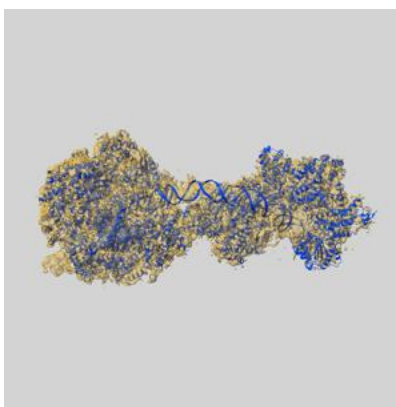
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-4980 and PDB model 9I3I. Per-residue inclusion information can be found in section [3](#) on page [10](#).

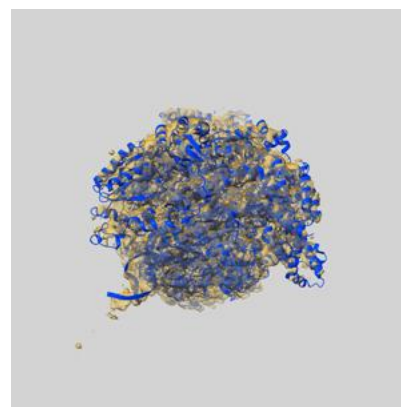
9.1 Map-model overlay [i](#)



X



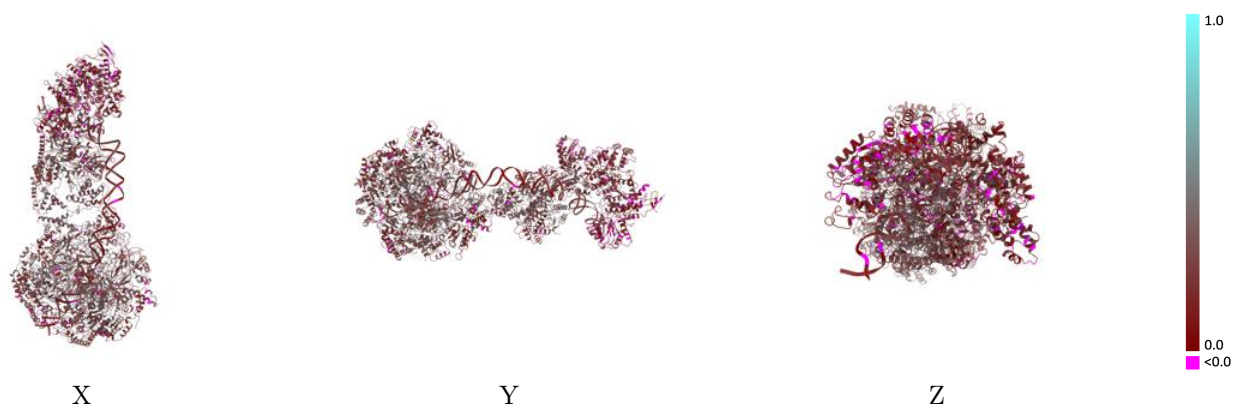
Y



Z

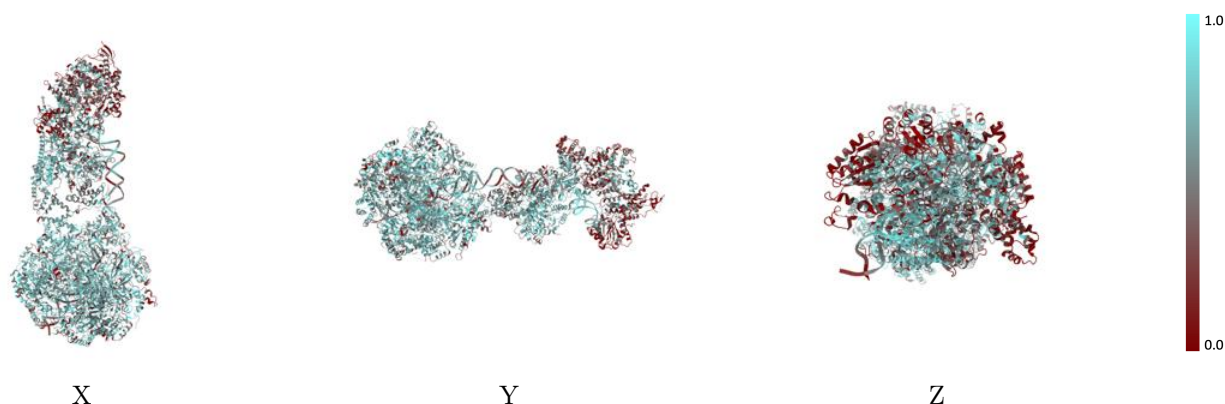
The images above show the 3D surface view of the map at the recommended contour level 0.016 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



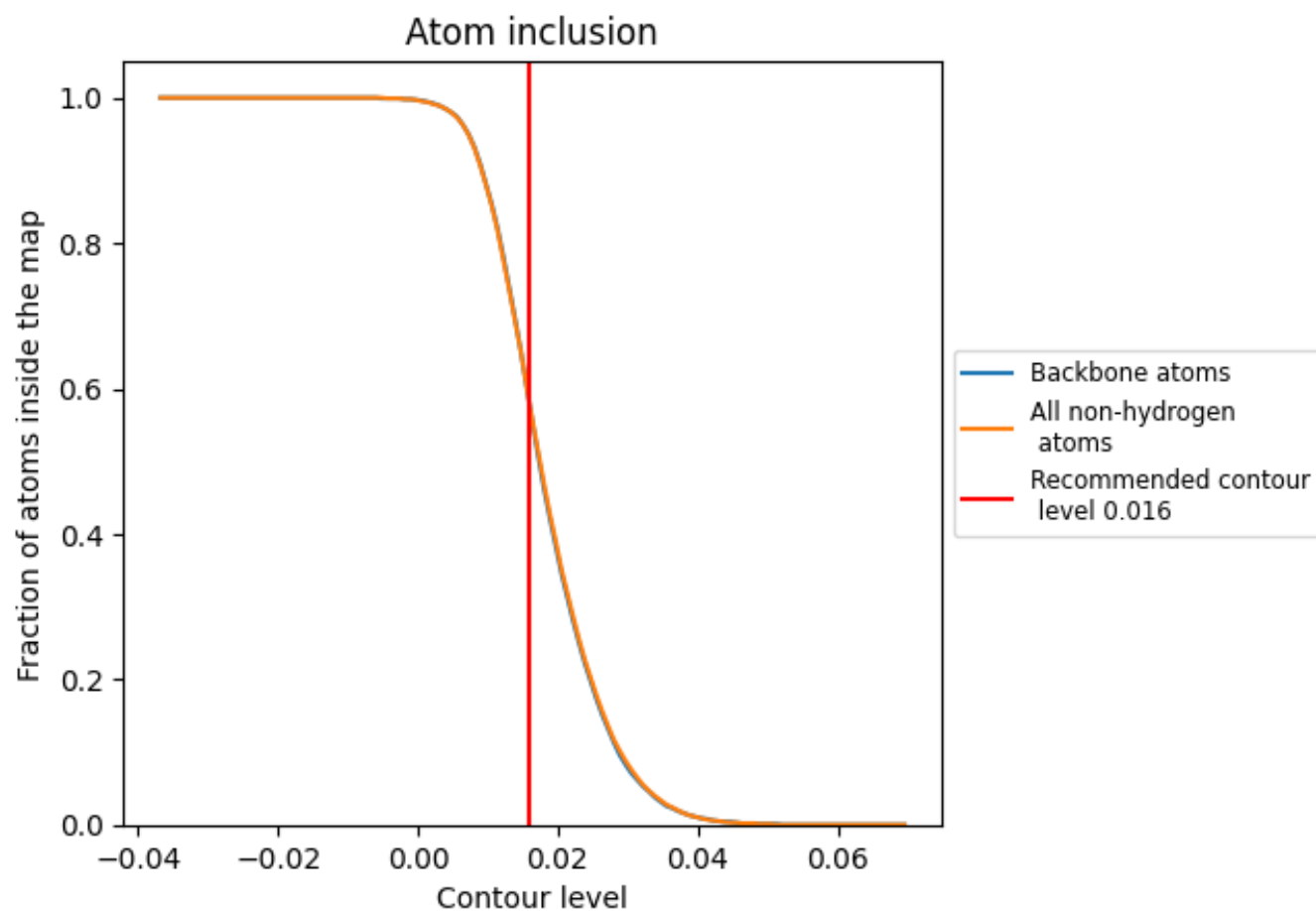
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.016).

9.4 Atom inclusion [i](#)



At the recommended contour level, 58% of all backbone atoms, 58% 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 (0.016) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5770	0.2680
2	0.6830	0.3130
3	0.6490	0.3030
4	0.6480	0.3030
5	0.7110	0.3240
6	0.6250	0.2840
7	0.6330	0.2950
A	0.2570	0.1500
B	0.6560	0.3160
C	0.6230	0.2820
D	0.4710	0.2110
E	0.3850	0.2090
F	0.5620	0.2810
X	0.5580	0.1650
Y	0.5760	0.1730

