



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 10:53 PM UTC

PDB ID : 5FJ9 / pdb_00005fj9
EMDB ID : EMD-3179
Title : Cryo-EM structure of yeast apo RNA polymerase III at 4.6 Å
Authors : Hoffmann, N.A.; Jakobi, A.J.; Moreno-Morcillo, M.; Glatt, S.; Kosinski, J.;
Hagen, W.J.; Sachse, C.; Muller, C.W.
Deposited on : 2015-10-06
Resolution : 4.60 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:

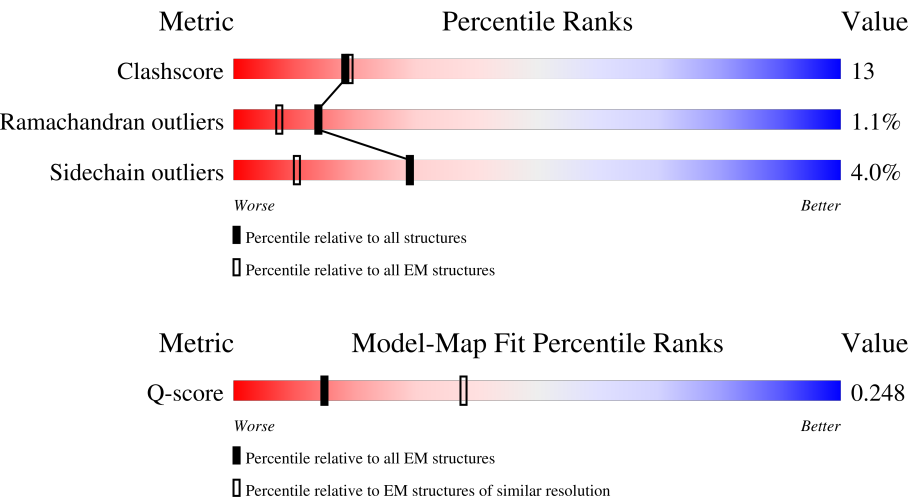
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 4.60 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	2407 (4.10 - 5.10)

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	1460	13% 69% 25% ••
2	B	1149	12% 66% 29% ••
3	C	335	7% 73% 24% •
4	D	161	16% 42% 31% • 26%

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Mol	Chain	Length	Quality of chain
5	E	215	
6	F	155	
7	G	212	
8	H	146	
9	I	110	
10	J	70	
11	K	142	
12	L	70	
13	M	282	
14	N	422	
15	O	654	
16	P	317	
17	Q	104	

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 38680 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-DIRECTED RNA POLYMERASE III SUBUNIT RPC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1412	Total	C	N	O	S	0	0
			11071	6979	1954	2079	59		

- Molecule 2 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	1114	Total	C	N	O	S	0	0
			8788	5558	1516	1654	60		

- Molecule 3 is a protein called DNA-DIRECTED RNA POLYMERASES I AND III SUBUNIT RPAC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	335	Total	C	N	O	S	0	0
			2655	1681	454	511	9		

- Molecule 4 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC9.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	119	Total	C	N	O	S	0	0
			977	628	156	187	6		

- Molecule 5 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	215	Total	C	N	O	S	0	0
			1759	1116	310	321	12		

- Molecule 6 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	83	Total	C	N	O	S	0	0
			671	429	114	125	3		

- Molecule 7 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC8.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	191	Total	C	N	O	S	0	0
			1544	1007	250	281	6		

- Molecule 8 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	140	Total	C	N	O	S	0	0
			1120	703	188	224	5		

- Molecule 9 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC10.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	92	Total	C	N	O	S	0	0
			728	455	117	145	11		

- Molecule 10 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	67	Total	C	N	O	S	0	0
			549	350	95	98	6		

- Molecule 11 is a protein called DNA-DIRECTED RNA POLYMERASES I AND III SUBUNIT RPAC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	101	Total	C	N	O	S	0	0
			792	496	130	161	5		

- Molecule 12 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	45	Total	C	N	O	S	0	0
			358	221	71	62	4		

- Molecule 13 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC5.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	165	Total	C	N	O	S	0	0
			1347	862	229	255	1		

- Molecule 14 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC4.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	105	Total	C	N	O	S	0	0
			802	508	144	147	3		

- Molecule 15 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	528	Total	C	N	O	S	0	0
			4247	2705	725	798	19		

- Molecule 16 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC6.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	98	Total	C	N	O	S	0	0
			820	533	126	157	4		

- Molecule 17 is a protein called DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC7.

Mol	Chain	Residues	Atoms				AltConf	Trace
17	Q	70	Total	C	N	O	0	0
			446	282	80	84		

- Molecule 18 is ZINC ION (CCD ID: ZN) (formula: Zn).

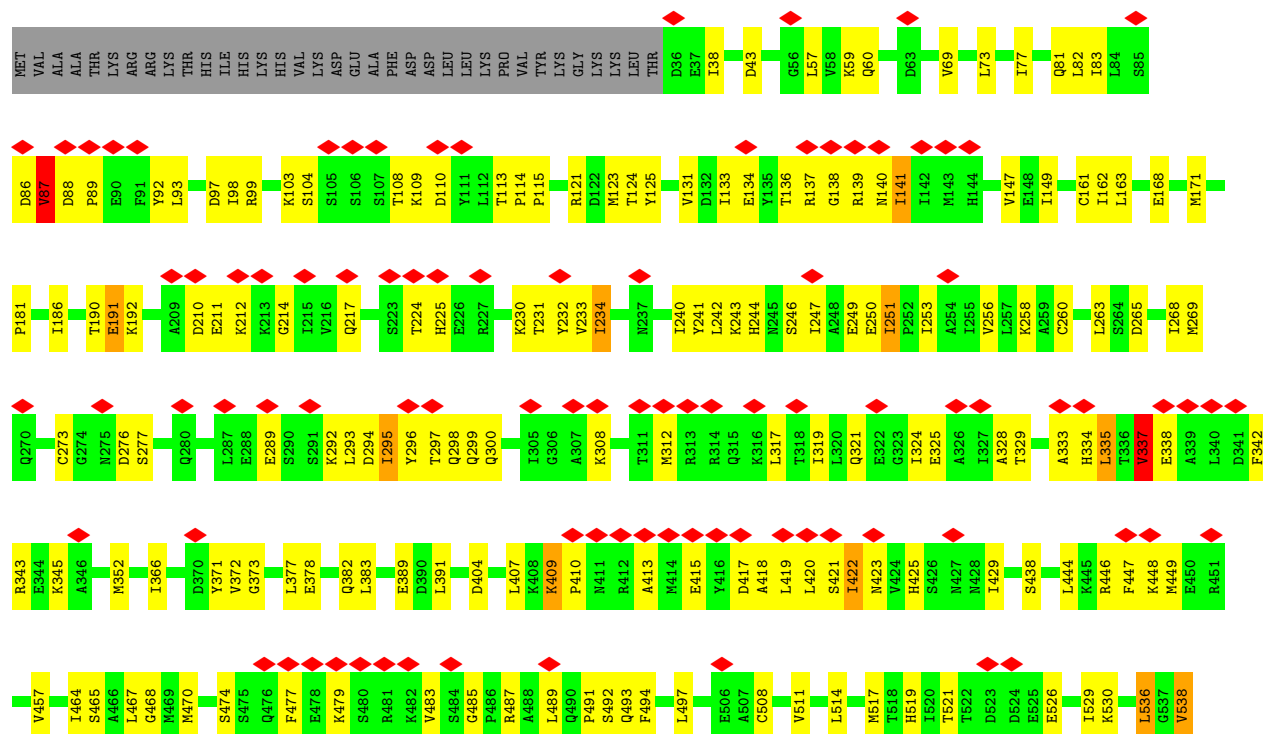
Mol	Chain	Residues	Atoms		AltConf
18	A	2	Total	Zn	0
			2	2	
18	B	1	Total	Zn	0
			1	1	
18	I	1	Total	Zn	0
			1	1	
18	J	1	Total	Zn	0
			1	1	
18	L	1	Total	Zn	0
			1	1	

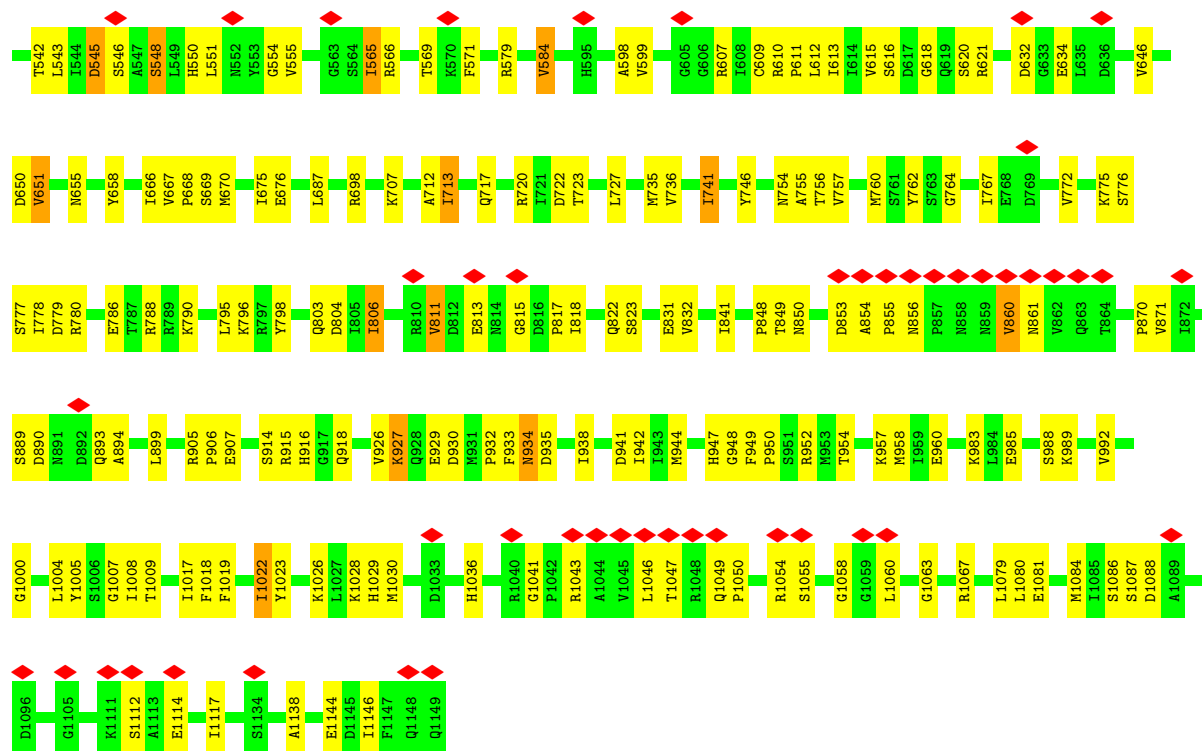
3 Residue-property plots

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.

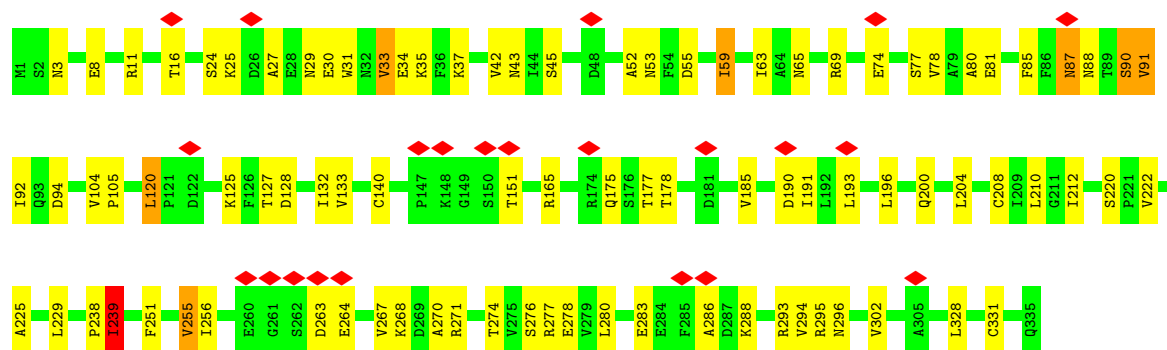
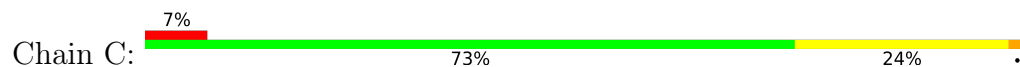
• Molecule 1: DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC1



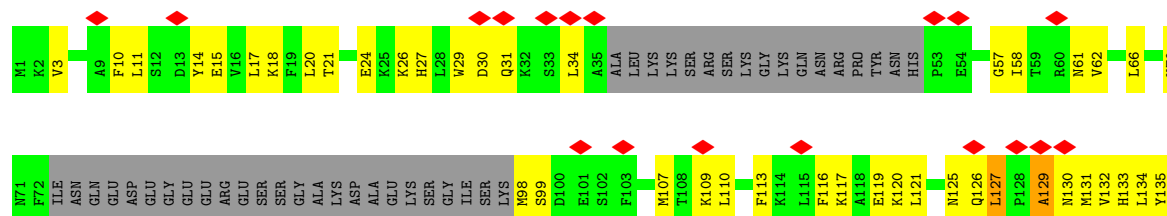
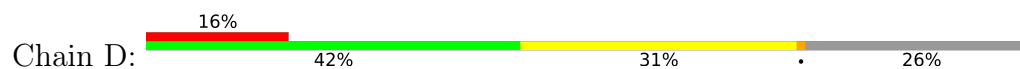


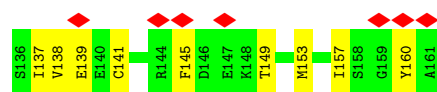


• Molecule 3: DNA-DIRECTED RNA POLYMERASES I AND III SUBUNIT RPAC1

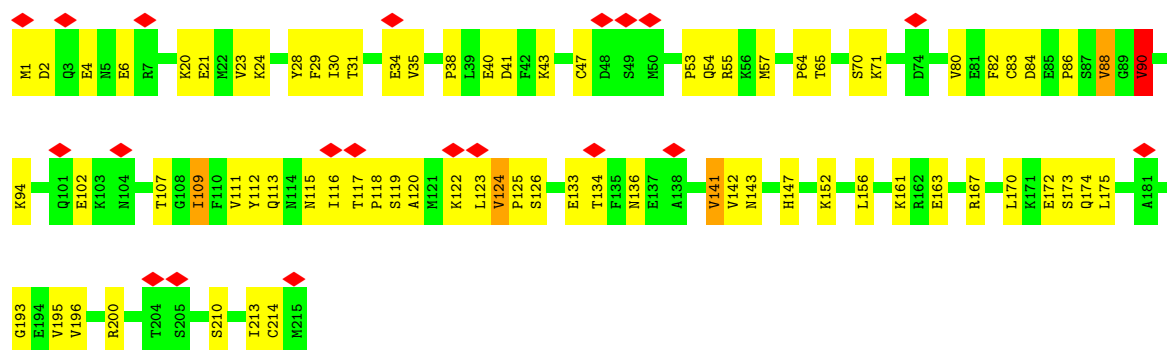


• Molecule 4: DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC9

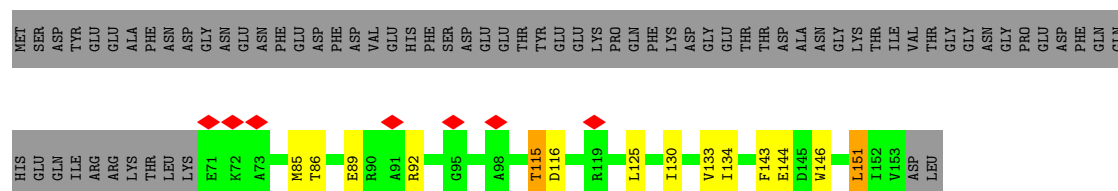
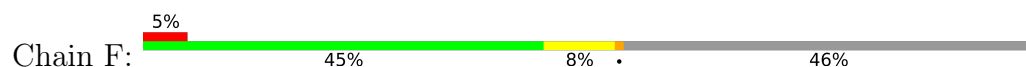




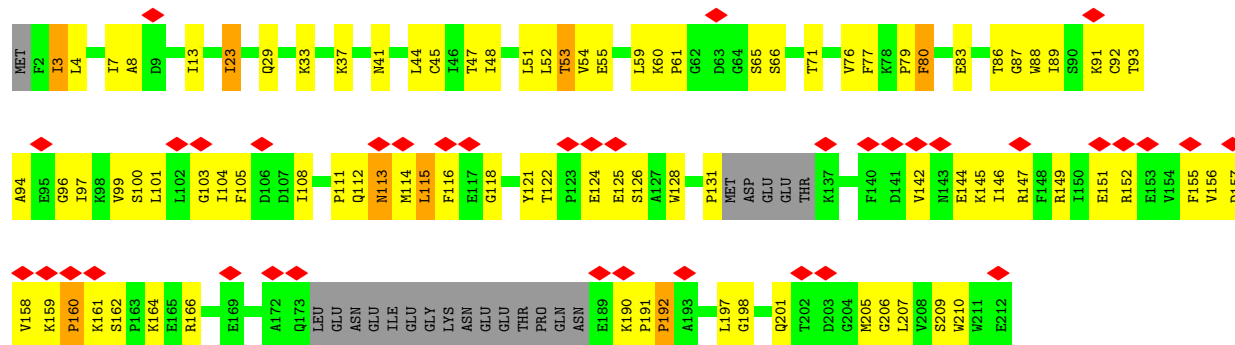
• Molecule 5: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 1



• Molecule 6: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 2

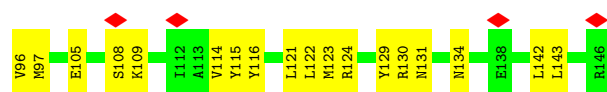


• Molecule 7: DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC8

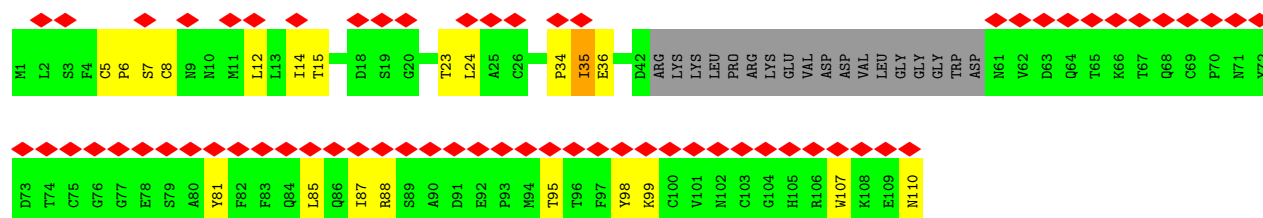


• Molecule 8: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 3

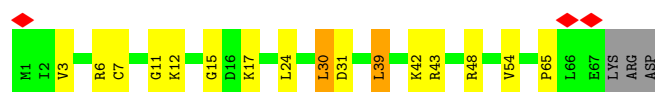




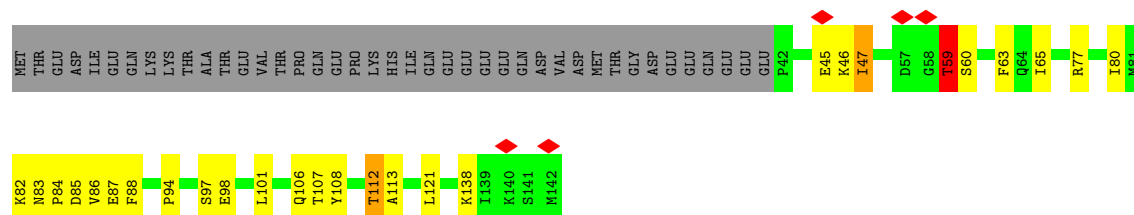
• Molecule 9: DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC10



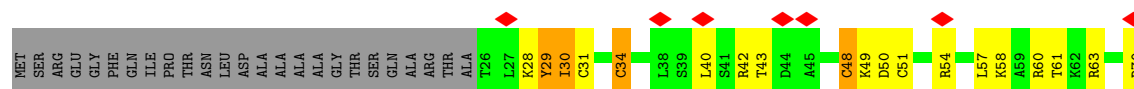
• Molecule 10: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 5



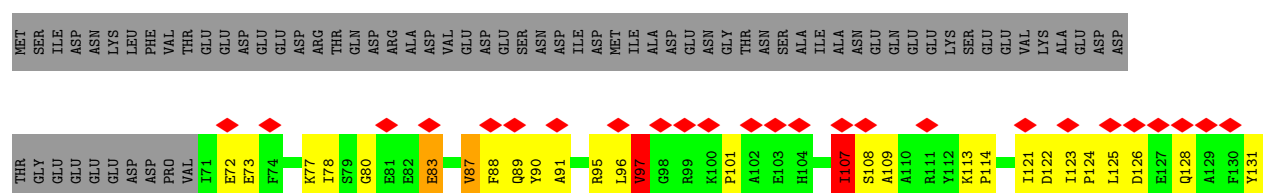
• Molecule 11: DNA-DIRECTED RNA POLYMERASES I AND III SUBUNIT RPAC2

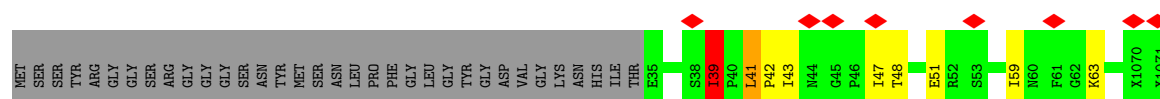


• Molecule 12: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 4



• Molecule 13: DNA-DIRECTED RNA POLYMERASE III SUBUNIT RPC5





WORLDWIDE
 **PDB**
PROTEIN DATA BANK

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	68818	Depositor
Resolution determination method	Not provided	
CTF correction method	EACH PARTICLE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	4200	Depositor
Magnification	75000	Depositor
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.322	Depositor
Minimum map value	-0.126	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.055	Depositor
Map size ($\text{\AA}$)	320.86398, 320.86398, 320.86398	wwPDB
Map dimensions	296, 296, 296	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.084, 1.084, 1.084	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.31	0/11267	0.85	31/15220 (0.2%)
2	B	0.29	0/8943	0.82	22/12068 (0.2%)
3	C	0.31	0/2711	0.83	4/3676 (0.1%)
4	D	0.31	0/991	0.87	1/1328 (0.1%)
5	E	0.30	0/1795	0.77	0/2416
6	F	0.30	0/683	0.77	0/923
7	G	0.32	0/1583	0.84	0/2146
8	H	0.28	0/1138	0.76	1/1540 (0.1%)
9	I	0.28	0/745	0.76	0/1007
10	J	0.26	0/558	0.72	0/750
11	K	0.29	0/803	0.79	1/1083 (0.1%)
12	L	0.27	0/360	0.78	1/478 (0.2%)
13	M	0.31	0/1378	0.83	2/1863 (0.1%)
14	N	0.35	0/810	0.88	0/1088
15	O	0.31	0/4310	0.86	9/5813 (0.2%)
16	P	0.43	0/837	0.91	3/1134 (0.3%)
17	Q	0.37	0/278	1.01	2/377 (0.5%)
All	All	0.31	0/39190	0.83	77/52910 (0.1%)

There are no bond length outliers.

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	858	PRO	CA-C-N	-8.16	109.64	119.84
1	A	858	PRO	C-N-CA	-8.16	109.64	119.84
17	Q	39	ILE	CA-C-N	7.58	127.74	120.31
17	Q	39	ILE	C-N-CA	7.58	127.74	120.31
2	B	479	LYS	CB-CA-C	-7.36	108.10	116.63
1	A	955	LEU	CA-C-N	7.36	126.87	118.85
1	A	955	LEU	C-N-CA	7.36	126.87	118.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	856	ASN	CA-C-N	6.82	126.85	119.90
2	B	856	ASN	C-N-CA	6.82	126.85	119.90
2	B	634	GLU	N-CA-C	6.63	116.61	107.73
1	A	1034	PRO	CA-C-N	6.60	126.62	119.89
1	A	1034	PRO	C-N-CA	6.60	126.62	119.89
15	O	328	ASP	CA-C-N	-6.54	111.67	119.84
15	O	328	ASP	C-N-CA	-6.54	111.67	119.84
2	B	815	GLY	N-CA-C	-6.26	104.98	113.37
1	A	1163	LYS	N-CA-C	6.21	117.20	108.00
1	A	1163	LYS	CB-CA-C	-6.21	107.64	116.34
2	B	477	PHE	CB-CA-C	-6.04	109.63	116.63
16	P	310	GLU	CB-CA-C	-6.00	109.67	116.63
2	B	319	ILE	N-CA-C	-5.99	106.48	112.17
16	P	306	LYS	N-CA-C	5.95	117.82	108.96
15	O	43	ASN	N-CA-C	5.94	119.29	109.79
1	A	1033	GLU	CA-C-N	-5.92	114.28	120.38
1	A	1033	GLU	C-N-CA	-5.92	114.28	120.38
2	B	1041	GLY	CA-C-N	5.90	125.90	119.89
2	B	1041	GLY	C-N-CA	5.90	125.90	119.89
2	B	479	LYS	N-CA-C	5.88	118.01	108.08
2	B	780	ARG	N-CA-C	-5.88	105.88	113.16
15	O	478	VAL	N-CA-C	5.86	113.67	108.63
13	M	174	GLU	CB-CA-C	-5.85	109.85	116.63
1	A	1050	ASP	CB-CA-C	-5.78	109.37	117.23
15	O	199	TYR	CB-CA-C	-5.75	109.42	117.23
2	B	276	ASP	N-CA-C	5.71	115.82	108.34
2	B	168	GLU	CB-CA-C	-5.59	109.62	117.23
1	A	219	MET	CB-CA-C	-5.58	110.16	116.63
2	B	409	LYS	CA-C-N	5.58	125.04	119.24
2	B	409	LYS	C-N-CA	5.58	125.04	119.24
1	A	1370	GLY	N-CA-C	5.55	118.11	110.56
15	O	559	ALA	N-CA-C	5.48	115.60	107.88
15	O	467	PHE	N-CA-C	5.47	117.88	109.07
1	A	448	ASP	CB-CA-C	-5.47	109.79	117.23
13	M	165	ASP	CB-CA-C	-5.46	110.29	116.63
16	P	314	PHE	CB-CA-C	-5.46	110.30	116.63
4	D	129	ALA	N-CA-C	5.44	119.86	113.23
2	B	214	GLY	N-CA-C	5.42	117.32	110.38
3	C	220	SER	CA-C-N	5.41	124.75	118.85
3	C	220	SER	C-N-CA	5.41	124.75	118.85
2	B	764	GLY	N-CA-C	-5.40	107.08	114.64
1	A	263	ALA	CA-C-N	5.39	123.54	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	ALA	C-N-CA	5.39	123.54	119.66
12	L	29	TYR	CB-CA-C	-5.39	109.38	115.79
15	O	42	LEU	N-CA-C	5.36	119.92	113.17
1	A	555	GLN	CB-CA-C	-5.36	109.95	117.23
1	A	247	THR	N-CA-C	-5.35	106.43	113.17
3	C	185	VAL	N-CA-C	5.34	113.57	109.19
3	C	90	SER	N-CA-C	5.33	117.43	107.99
1	A	1179	ASP	CB-CA-C	-5.32	109.99	117.23
2	B	934	ASN	N-CA-C	5.30	115.84	108.00
15	O	362	ASN	N-CA-C	5.29	117.70	110.24
1	A	1347	GLY	N-CA-C	5.28	118.46	110.75
1	A	1123	VAL	CA-C-N	-5.26	114.25	119.56
1	A	1123	VAL	C-N-CA	-5.26	114.25	119.56
1	A	686	GLY	CA-C-N	5.23	124.73	119.19
1	A	686	GLY	C-N-CA	5.23	124.73	119.19
1	A	979	ASN	CB-CA-C	-5.23	110.12	117.23
11	K	59	THR	CB-CA-C	-5.18	110.18	117.23
2	B	335	LEU	N-CA-C	5.17	114.96	108.45
2	B	277	SER	CB-CA-C	-5.16	109.11	116.34
8	H	32	THR	CB-CA-C	-5.15	110.22	117.23
1	A	757	ASN	N-CA-C	5.12	116.54	110.19
1	A	1084	GLU	CA-C-N	5.11	125.10	119.89
1	A	1084	GLU	C-N-CA	5.11	125.10	119.89
1	A	248	VAL	N-CA-C	5.11	119.92	108.88
2	B	676	GLU	CA-C-N	5.09	124.75	119.56
2	B	676	GLU	C-N-CA	5.09	124.75	119.56
1	A	1162	GLU	N-CA-C	5.06	117.15	108.90
1	A	756	CYS	CB-CA-C	-5.02	110.80	116.63

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	11071	0	11200	282	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	8788	0	8902	234	0
3	C	2655	0	2628	55	0
4	D	977	0	983	36	0
5	E	1759	0	1788	48	0
6	F	671	0	692	10	0
7	G	1544	0	1540	62	0
8	H	1120	0	1089	30	0
9	I	728	0	672	18	0
10	J	549	0	560	8	0
11	K	792	0	790	23	0
12	L	358	0	381	18	0
13	M	1347	0	1315	41	0
14	N	802	0	851	42	0
15	O	4247	0	4413	164	0
16	P	820	0	795	54	0
17	Q	446	0	324	10	0
18	A	2	0	0	0	0
18	B	1	0	0	0	0
18	I	1	0	0	0	0
18	J	1	0	0	0	0
18	L	1	0	0	0	0
All	All	38680	0	38923	1014	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 13.

All (1014) 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)
15:O:516:LEU:CD2	15:O:567:ARG:HE	1.08	1.58
15:O:516:LEU:HD23	15:O:567:ARG:NE	1.20	1.41
15:O:516:LEU:HD23	15:O:567:ARG:CD	1.60	1.30
15:O:516:LEU:CD2	15:O:567:ARG:NE	1.84	1.23
1:A:597:ILE:HB	1:A:600:PRO:HG3	1.15	1.15
15:O:516:LEU:HD22	15:O:567:ARG:HE	1.08	1.08
1:A:598:MET:C	1:A:600:PRO:HD3	1.78	1.08
1:A:597:ILE:CB	1:A:600:PRO:HG3	1.87	1.02
15:O:516:LEU:HB3	15:O:567:ARG:HG3	1.42	1.01
15:O:569:LYS:HE3	15:O:569:LYS:HA	1.43	1.00
3:C:251:PHE:HB3	3:C:255:VAL:HG11	1.51	0.92
2:B:548:SER:HB3	2:B:550:HIS:CE1	2.04	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ASN:HB3	15:O:544:ASN:HD21	1.37	0.90
1:A:43:GLU:HB2	1:A:48:PRO:HD2	1.55	0.86
1:A:597:ILE:HB	1:A:600:PRO:CG	2.02	0.86
4:D:126:GLN:HG3	4:D:127:LEU:HG	1.57	0.86
1:A:226:LYS:HE2	15:O:548:ILE:HG13	1.56	0.85
15:O:549:GLN:HG3	15:O:565:LEU:HB2	1.58	0.85
15:O:547:GLU:HG3	15:O:569:LYS:HD3	1.59	0.84
15:O:512:ARG:HG2	15:O:568:CYS:SG	2.17	0.84
15:O:569:LYS:HA	15:O:569:LYS:CE	2.07	0.83
15:O:516:LEU:HD23	15:O:567:ARG:CG	2.08	0.83
1:A:598:MET:C	1:A:600:PRO:CD	2.53	0.81
15:O:547:GLU:CG	15:O:569:LYS:HD3	2.10	0.80
13:M:113:LYS:HD3	13:M:241:ALA:HB2	1.63	0.79
2:B:294:ASP:HB3	2:B:300:GLN:HG3	1.64	0.79
5:E:124:VAL:HG13	5:E:125:PRO:HD3	1.64	0.79
15:O:493:GLU:OE2	15:O:512:ARG:NH1	2.16	0.78
15:O:516:LEU:CB	15:O:567:ARG:HG3	2.13	0.77
1:A:599:LYS:N	1:A:600:PRO:CD	2.46	0.77
1:A:229:ASN:CB	15:O:544:ASN:HD21	1.97	0.76
1:A:829:ASP:OD2	1:A:838:ASN:ND2	2.20	0.75
1:A:550:ILE:HG23	1:A:551:ILE:HG23	1.69	0.74
1:A:756:CYS:SG	1:A:757:ASN:N	2.59	0.73
15:O:570:GLU:HG3	15:O:571:THR:HG23	1.70	0.73
1:A:49:LYS:HD2	1:A:54:LEU:HB3	1.69	0.73
3:C:120:LEU:HD23	3:C:125:LYS:HG2	1.72	0.72
11:K:65:ILE:HB	11:K:101:LEU:HB3	1.72	0.72
15:O:538:ALA:HA	15:O:541:ILE:HG22	1.72	0.72
16:P:236:THR:HG22	16:P:237:PRO:HD2	1.69	0.72
1:A:597:ILE:CG2	1:A:600:PRO:HG3	2.19	0.72
7:G:146:ILE:HG23	7:G:206:GLY:HA2	1.70	0.72
14:N:363:ILE:HG12	14:N:373:VAL:HG13	1.72	0.72
15:O:552:PRO:HB3	15:O:557:ARG:H	1.55	0.71
1:A:215:VAL:HG21	15:O:554:THR:H	1.53	0.71
16:P:216:SER:HB3	16:P:222:LEU:HD11	1.71	0.71
7:G:207:LEU:HD23	7:G:209:SER:H	1.55	0.71
1:A:69:THR:HA	7:G:164:LYS:HZ3	1.55	0.70
2:B:822:GLN:HG3	2:B:823:SER:H	1.55	0.70
15:O:190:LEU:HG	15:O:193:GLN:HB3	1.71	0.70
13:M:87:VAL:HA	14:N:396:ALA:HA	1.73	0.70
15:O:124:GLU:HG3	15:O:125:GLU:N	2.06	0.70
12:L:48:CYS:SG	12:L:51:CYS:HB2	2.17	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:499:THR:HA	17:Q:41:LEU:HD21	1.72	0.70
15:O:555:ALA:HB1	15:O:558:SER:HB2	1.73	0.70
2:B:234:ILE:HB	2:B:240:ILE:HG22	1.74	0.70
4:D:127:LEU:HB2	4:D:133:HIS:HB3	1.73	0.69
1:A:436:ARG:HD3	1:A:459:GLY:HA3	1.75	0.69
13:M:148:LEU:HG	13:M:181:PRO:HB3	1.74	0.69
15:O:507:LEU:HD13	15:O:540:LEU:HD22	1.75	0.69
1:A:332:VAL:HG13	1:A:333:ASN:H	1.57	0.69
7:G:96:GLY:HA3	7:G:111:PRO:HA	1.73	0.69
9:I:81:TYR:HB2	9:I:99:LYS:HB3	1.74	0.69
1:A:1020:ALA:HB2	1:A:1032:LEU:HD11	1.75	0.68
2:B:265:ASP:O	2:B:268:ILE:N	2.25	0.68
2:B:548:SER:CB	2:B:550:HIS:CE1	2.76	0.68
3:C:16:THR:O	3:C:295:ARG:NH1	2.26	0.68
1:A:160:LEU:HD21	15:O:339:LEU:HG	1.76	0.68
1:A:153:ARG:HA	1:A:160:LEU:HD22	1.76	0.68
1:A:476:ARG:HG2	1:A:478:PRO:HD2	1.76	0.68
4:D:126:GLN:NE2	7:G:86:THR:OG1	2.24	0.68
8:H:24:CYS:SG	8:H:25:ARG:N	2.67	0.67
16:P:308:ASP:OD1	16:P:309:LYS:HG3	1.94	0.67
1:A:615:LYS:NZ	1:A:620:SER:O	2.26	0.67
16:P:248:VAL:HG23	16:P:255:LYS:HE2	1.76	0.67
1:A:600:PRO:O	1:A:601:TYR:HB2	1.92	0.67
2:B:760:MET:HE3	2:B:762:TYR:HB2	1.74	0.67
1:A:1178:LYS:HE2	9:I:36:GLU:HB3	1.76	0.67
1:A:221:ASP:HB3	15:O:550:GLU:HG3	1.76	0.67
11:K:88:PHE:HB3	11:K:106:GLN:HB3	1.74	0.67
11:K:47:ILE:HG12	11:K:65:ILE:HG12	1.76	0.66
15:O:123:ASN:O	15:O:124:GLU:HG2	1.95	0.66
16:P:223:GLU:O	16:P:227:ALA:N	2.22	0.66
2:B:778:ILE:HD11	2:B:906:PRO:HG2	1.78	0.66
8:H:58:THR:HB	8:H:143:LEU:HB3	1.79	0.65
10:J:17:LYS:HB3	10:J:39:LEU:HD21	1.78	0.65
15:O:541:ILE:HD11	15:O:548:ILE:HG12	1.78	0.65
1:A:226:LYS:HD3	15:O:547:GLU:OE1	1.96	0.65
2:B:247:ILE:HG21	2:B:251:ILE:HD11	1.79	0.65
2:B:419:LEU:O	2:B:423:ASN:ND2	2.23	0.65
15:O:341:GLU:HG2	15:O:344:SER:HB2	1.77	0.65
4:D:130:ASN:O	4:D:133:HIS:ND1	2.30	0.65
2:B:59:LYS:HE2	2:B:519:HIS:HA	1.78	0.65
2:B:1112:SER:HB2	2:B:1114:GLU:HG2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:VAL:HG11	15:O:554:THR:HA	1.78	0.64
7:G:161:LYS:HE3	7:G:166:ARG:HA	1.79	0.64
4:D:119:GLU:HG3	4:D:138:VAL:HB	1.80	0.64
9:I:12:LEU:HB3	9:I:24:LEU:HD13	1.79	0.64
16:P:308:ASP:O	16:P:309:LYS:HB2	1.98	0.64
3:C:222:VAL:HG11	3:C:225:ALA:HB2	1.78	0.64
16:P:233:VAL:HG12	16:P:234:GLU:H	1.62	0.64
1:A:599:LYS:H	1:A:602:TYR:HE1	1.43	0.64
2:B:258:LYS:HB3	2:B:297:THR:HA	1.80	0.64
12:L:31:CYS:HB3	12:L:34:CYS:HB2	1.80	0.64
15:O:150:GLU:HG3	15:O:151:GLU:HG2	1.80	0.64
16:P:221:ILE:HA	16:P:224:PHE:HB2	1.80	0.63
7:G:115:LEU:HD23	7:G:116:PHE:H	1.62	0.63
1:A:956:PRO:HG2	1:A:957:TYR:HD1	1.64	0.63
2:B:543:LEU:HD21	13:M:176:VAL:HG21	1.81	0.63
2:B:230:LYS:HE3	2:B:352:MET:HE1	1.80	0.63
7:G:104:ILE:HG23	7:G:105:PHE:H	1.63	0.63
15:O:185:TYR:HA	15:O:188:SER:HB2	1.78	0.63
16:P:307:HIS:O	16:P:308:ASP:HB3	1.98	0.63
2:B:81:GLN:HG2	2:B:82:LEU:HG	1.80	0.63
2:B:192:LYS:NZ	2:B:438:SER:O	2.31	0.63
1:A:573:ARG:HH21	11:K:87:GLU:HB3	1.63	0.62
2:B:137:ARG:HH11	2:B:139:ARG:HB2	1.63	0.62
15:O:516:LEU:HD23	15:O:567:ARG:HG3	1.81	0.62
1:A:1161:VAL:HG11	1:A:1303:VAL:HG23	1.82	0.62
2:B:1026:LYS:NZ	2:B:1030:MET:SD	2.69	0.62
2:B:775:LYS:HA	2:B:778:ILE:HG22	1.81	0.62
6:F:92:ARG:HH12	7:G:61:PRO:HB3	1.64	0.62
2:B:253:ILE:HD12	2:B:256:VAL:HB	1.82	0.62
1:A:67:CYS:SG	1:A:69:THR:OG1	2.57	0.62
15:O:488:LYS:NZ	15:O:650:VAL:O	2.33	0.62
1:A:1420:THR:HG22	2:B:1080:LEU:HD21	1.82	0.62
3:C:80:ALA:HA	3:C:208:CYS:HA	1.82	0.62
12:L:29:TYR:HE2	12:L:40:LEU:HB2	1.65	0.62
1:A:1169:VAL:HG13	1:A:1192:THR:HB	1.82	0.61
1:A:413:ARG:NH1	1:A:456:LEU:O	2.34	0.61
1:A:112:ALA:HA	1:A:234:ILE:HD11	1.81	0.61
1:A:213:ARG:O	1:A:213:ARG:NH1	2.32	0.61
1:A:393:ALA:HA	1:A:491:LYS:HB2	1.82	0.61
2:B:139:ARG:HE	2:B:141:ILE:HD11	1.64	0.61
7:G:87:GLY:O	7:G:146:ILE:N	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:712:ILE:HD12	2:B:949:PHE:HD2	1.65	0.61
1:A:1372:THR:O	1:A:1376:LEU:N	2.34	0.61
2:B:757:VAL:HG23	2:B:942:ILE:HB	1.82	0.61
12:L:42:ARG:HG3	12:L:43:THR:HG23	1.83	0.61
1:A:1257:PHE:HB2	9:I:14:ILE:HB	1.82	0.61
5:E:112:TYR:HB2	5:E:136:ASN:HA	1.82	0.61
16:P:314:PHE:C	16:P:316:GLU:H	2.08	0.61
1:A:252:ARG:O	1:A:254:GLU:N	2.33	0.61
15:O:106:TYR:HB3	15:O:208:TYR:HE2	1.65	0.61
15:O:198:GLY:HA3	15:O:286:ARG:HG2	1.83	0.60
2:B:1007:GLY:O	3:C:69:ARG:NH1	2.34	0.60
15:O:547:GLU:HG2	15:O:569:LYS:HD3	1.84	0.60
16:P:311:VAL:O	16:P:312:VAL:HG12	2.01	0.60
1:A:582:MET:HE3	1:A:703:ARG:HB2	1.83	0.60
15:O:584:ASN:OD1	16:P:312:VAL:HG23	2.00	0.60
1:A:599:LYS:N	1:A:600:PRO:HD2	2.17	0.60
7:G:89:ILE:HA	7:G:99:VAL:HA	1.82	0.60
15:O:286:ARG:NH1	15:O:320:GLU:O	2.35	0.60
1:A:62:SER:HB3	1:A:65:LEU:HG	1.84	0.60
1:A:121:ARG:NH2	15:O:212:GLU:OE1	2.35	0.60
1:A:1048:VAL:HG11	1:A:1053:LYS:HB2	1.83	0.60
4:D:127:LEU:HD11	4:D:137:ILE:HD13	1.83	0.60
16:P:313:TYR:HB2	17:Q:39:ILE:HD11	1.84	0.60
5:E:88:VAL:HG23	5:E:117:THR:HG22	1.83	0.60
13:M:91:ALA:HA	14:N:391:LEU:HD21	1.83	0.60
1:A:1144:VAL:HG23	1:A:1310:ILE:HG23	1.84	0.60
1:A:1164:THR:HB	1:A:1271:VAL:HA	1.83	0.60
15:O:31:VAL:HG12	15:O:32:MET:HG3	1.82	0.60
1:A:15:GLY:HA2	1:A:1407:ALA:HA	1.83	0.59
2:B:775:LYS:HB3	2:B:927:LYS:HA	1.83	0.59
15:O:357:PRO:HA	15:O:361:PHE:CD2	2.36	0.59
16:P:179:LEU:O	16:P:183:TRP:NE1	2.34	0.59
3:C:31:TRP:H	11:K:82:LYS:HE2	1.67	0.59
8:H:97:MET:HB2	8:H:142:LEU:HB3	1.82	0.59
14:N:305:MET:HG2	14:N:414:GLY:H	1.66	0.59
1:A:486:LEU:HD23	1:A:537:VAL:HG22	1.83	0.59
1:A:1084:GLU:HB2	6:F:86:THR:HG23	1.83	0.59
4:D:129:ALA:HB2	4:D:157:ILE:HG23	1.84	0.59
1:A:15:GLY:H	1:A:1408:VAL:HG12	1.66	0.59
7:G:157:ASP:OD1	7:G:157:ASP:N	2.36	0.59
9:I:87:ILE:HG23	9:I:88:ARG:HG3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:363:ILE:HG23	14:N:373:VAL:HG22	1.84	0.59
7:G:101:LEU:HD22	7:G:104:ILE:HG22	1.84	0.59
13:M:108:SER:OG	13:M:109:ALA:N	2.36	0.58
1:A:402:LEU:HD23	1:A:466:LEU:HD22	1.85	0.58
2:B:137:ARG:HD2	2:B:415:GLU:HB2	1.85	0.58
1:A:360:LYS:HD2	1:A:1389:PHE:O	2.04	0.58
2:B:616:SER:HB2	2:B:621:ARG:HH21	1.68	0.58
15:O:286:ARG:NH2	15:O:321:GLN:O	2.36	0.58
1:A:828:GLN:HB3	2:B:655:ASN:HD21	1.68	0.58
3:C:33:VAL:O	3:C:34:GLU:HG3	2.03	0.58
15:O:288:MET:SD	15:O:291:ARG:NH1	2.77	0.58
11:K:45:GLU:O	11:K:46:LYS:HD2	2.04	0.58
2:B:926:VAL:HG11	2:B:930:ASP:HB2	1.86	0.58
3:C:165:ARG:NH2	3:C:190:ASP:OD1	2.35	0.58
15:O:547:GLU:HG2	15:O:569:LYS:CD	2.33	0.58
1:A:358:LYS:HG2	1:A:1392:THR:HG21	1.86	0.58
2:B:832:VAL:HG12	12:L:60:ARG:HG3	1.86	0.58
15:O:357:PRO:HA	15:O:361:PHE:HD2	1.68	0.58
1:A:538:LYS:HB3	1:A:687:PRO:HB2	1.86	0.58
7:G:190:LYS:C	7:G:192:PRO:HD3	2.28	0.58
15:O:105:LYS:N	15:O:123:ASN:HB2	2.18	0.57
15:O:163:VAL:HG22	15:O:169:LEU:HD22	1.85	0.57
16:P:174:PHE:O	16:P:175:ILE:HG12	2.04	0.57
1:A:225:LEU:O	1:A:229:ASN:ND2	2.37	0.57
2:B:554:GLY:H	2:B:598:ALA:HA	1.68	0.57
2:B:1004:LEU:HD12	2:B:1017:ILE:HD12	1.86	0.57
13:M:247:TRP:HE1	14:N:408:LEU:HB2	1.69	0.57
15:O:191:PHE:HD2	15:O:274:VAL:HG21	1.69	0.57
1:A:1059:LEU:HD21	8:H:105:GLU:HA	1.86	0.57
2:B:87:VAL:HG11	2:B:407:LEU:HB3	1.86	0.57
1:A:412:ASN:HB3	1:A:416:LEU:HD13	1.86	0.57
2:B:343:ARG:HH22	2:B:546:SER:HB3	1.69	0.57
15:O:105:LYS:HB3	15:O:121:TYR:HB2	1.84	0.57
5:E:21:GLU:HB3	5:E:35:VAL:HG21	1.85	0.57
13:M:77:LYS:HE3	13:M:262:GLU:HB3	1.85	0.57
13:M:183:PHE:HB2	13:M:186:ILE:HG13	1.87	0.57
15:O:124:GLU:HG3	15:O:125:GLU:H	1.67	0.57
15:O:338:ASP:OD1	15:O:339:LEU:N	2.38	0.57
1:A:270:PRO:HD3	2:B:1046:LEU:HD13	1.87	0.57
15:O:652:GLN:HE21	15:O:653:MET:HE3	1.69	0.57
1:A:232:LYS:NZ	16:P:316:GLU:OE1	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:389:GLU:OE2	2:B:446:ARG:NH2	2.37	0.57
9:I:5:CYS:HB2	9:I:12:LEU:HD21	1.85	0.57
11:K:107:THR:OG1	11:K:108:TYR:N	2.36	0.57
12:L:48:CYS:CB	12:L:51:CYS:HB2	2.30	0.57
5:E:90:VAL:HG12	5:E:94:LYS:HG2	1.86	0.57
1:A:317:ASP:OD1	1:A:320:GLN:NE2	2.38	0.57
2:B:373:GLY:HA2	2:B:607:ARG:HH22	1.70	0.57
12:L:49:LYS:N	12:L:51:CYS:SG	2.78	0.57
15:O:158:GLU:OE1	15:O:161:GLN:NE2	2.38	0.57
2:B:536:LEU:HD22	2:B:571:PHE:HD1	1.70	0.56
11:K:80:ILE:HG22	11:K:86:VAL:HG11	1.87	0.56
7:G:118:GLY:HA3	7:G:131:PRO:HD3	1.86	0.56
1:A:574:ALA:O	11:K:77:ARG:NH2	2.38	0.56
2:B:493:GLN:HG3	2:B:497:LEU:HB2	1.88	0.56
2:B:698:ARG:HE	2:B:952:ARG:HB3	1.69	0.56
4:D:17:LEU:HB2	4:D:66:LEU:HD23	1.87	0.56
7:G:91:LYS:HG2	7:G:92:CYS:H	1.70	0.56
7:G:112:GLN:HG2	7:G:115:LEU:HD13	1.86	0.56
11:K:46:LYS:O	11:K:47:ILE:C	2.48	0.56
15:O:540:LEU:HD21	15:O:546:VAL:HG11	1.86	0.56
16:P:186:ILE:HG23	16:P:253:LEU:HD21	1.87	0.56
2:B:57:LEU:HD22	2:B:467:LEU:HD11	1.86	0.56
2:B:325:GLU:O	2:B:329:THR:N	2.34	0.56
12:L:28:LYS:HG3	12:L:29:TYR:HD2	1.70	0.56
15:O:221:LYS:O	15:O:225:ASN:ND2	2.39	0.56
1:A:276:ASP:N	1:A:276:ASP:OD1	2.39	0.56
15:O:347:ASP:HA	15:O:350:GLU:HB2	1.88	0.56
5:E:83:CYS:SG	5:E:84:ASP:N	2.77	0.56
1:A:272:VAL:HG23	1:A:281:ASN:HB3	1.86	0.56
1:A:712:ILE:HG23	1:A:790:ALA:HB3	1.87	0.56
1:A:799:SER:OG	1:A:800:LYS:N	2.39	0.56
1:A:1332:ARG:HB2	1:A:1363:THR:HG21	1.88	0.56
1:A:1364:TYR:OH	1:A:1379:MET:SD	2.62	0.56
3:C:140:CYS:HB2	3:C:196:LEU:HD13	1.87	0.56
1:A:252:ARG:HB3	1:A:253:PRO:HD2	1.87	0.56
2:B:161:CYS:SG	2:B:162:ILE:N	2.79	0.56
7:G:45:CYS:HA	7:G:76:VAL:HA	1.87	0.56
11:K:65:ILE:HD12	11:K:101:LEU:HD23	1.86	0.56
2:B:244:HIS:HD1	2:B:246:SER:HG	1.53	0.56
7:G:3:ILE:HG13	7:G:76:VAL:HG23	1.87	0.56
1:A:645:MET:HE1	8:H:124:ARG:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:50:ASP:OD1	12:L:51:CYS:N	2.39	0.55
15:O:316:LEU:HA	15:O:319:THR:HG22	1.89	0.55
15:O:505:MET:HB3	16:P:250:ASP:OD2	2.07	0.55
2:B:618:GLY:HA2	2:B:668:PRO:HB3	1.88	0.55
15:O:228:ARG:HH12	15:O:231:PRO:HD3	1.72	0.55
16:P:314:PHE:O	16:P:316:GLU:N	2.39	0.55
1:A:476:ARG:NH1	1:A:509:ASN:O	2.39	0.55
1:A:1084:GLU:O	1:A:1087:THR:OG1	2.25	0.55
2:B:233:VAL:HG11	2:B:243:LYS:HE3	1.89	0.55
9:I:14:ILE:HD13	9:I:24:LEU:HB3	1.87	0.55
16:P:312:VAL:HG13	16:P:312:VAL:O	2.05	0.55
1:A:1343:MET:HE3	1:A:1350:VAL:HG21	1.89	0.55
6:F:125:LEU:HA	6:F:130:ILE:HD11	1.88	0.55
7:G:207:LEU:HD22	7:G:210:TRP:CD1	2.41	0.55
2:B:1036:HIS:CE1	2:B:1058:GLY:HA3	2.41	0.55
7:G:13:ILE:HG13	7:G:66:SER:HB3	1.88	0.55
13:M:123:ILE:O	13:M:146:GLN:N	2.36	0.55
2:B:519:HIS:O	2:B:609:CYS:N	2.38	0.55
5:E:38:PRO:HB2	5:E:40:GLU:HG2	1.87	0.55
14:N:366:HIS:HB3	14:N:370:LYS:HB2	1.89	0.55
15:O:185:TYR:O	15:O:189:SER:N	2.39	0.55
16:P:223:GLU:HG2	16:P:240:ILE:HA	1.87	0.55
2:B:483:VAL:HG12	2:B:485:GLY:H	1.71	0.55
5:E:117:THR:HG23	5:E:119:SER:H	1.72	0.55
15:O:328:ASP:OD1	15:O:328:ASP:N	2.40	0.55
16:P:235:LEU:HD12	16:P:236:THR:H	1.72	0.55
1:A:155:LEU:HG	15:O:336:LEU:HD23	1.89	0.54
1:A:541:LEU:HG	1:A:551:ILE:HD11	1.89	0.54
5:E:156:LEU:HD11	5:E:195:VAL:HB	1.89	0.54
15:O:495:VAL:HG11	15:O:577:MET:HE3	1.89	0.54
16:P:223:GLU:HB3	16:P:240:ILE:HD13	1.89	0.54
2:B:115:PRO:HG3	2:B:163:LEU:HD11	1.89	0.54
2:B:242:LEU:HD11	2:B:253:ILE:HD13	1.88	0.54
13:M:88:PHE:HD2	14:N:397:LEU:HD13	1.72	0.54
16:P:251:ASP:OD1	16:P:252:LYS:N	2.36	0.54
1:A:1141:ILE:N	1:A:1295:VAL:O	2.34	0.54
2:B:860:VAL:HG13	2:B:861:ASN:H	1.71	0.54
4:D:11:LEU:HD22	7:G:4:LEU:HD13	1.88	0.54
5:E:31:THR:HG22	5:E:34:GLU:HG2	1.89	0.54
11:K:85:ASP:O	11:K:107:THR:OG1	2.20	0.54
13:M:83:GLU:HB2	14:N:400:ALA:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:343:ARG:NH2	2:B:546:SER:HB3	2.23	0.54
14:N:397:LEU:HD11	14:N:408:LEU:HD12	1.90	0.54
15:O:578:ARG:HG2	15:O:648:TRP:HZ3	1.73	0.54
2:B:521:THR:OG1	2:B:609:CYS:SG	2.66	0.54
5:E:109:ILE:HG22	5:E:133:GLU:HB3	1.90	0.54
5:E:141:VAL:HG23	5:E:142:VAL:H	1.73	0.54
15:O:652:GLN:HG3	15:O:653:MET:HG3	1.90	0.54
2:B:295:ILE:O	2:B:296:TYR:CG	2.60	0.54
12:L:30:ILE:N	12:L:57:LEU:O	2.41	0.54
1:A:22:SER:O	1:A:24:ALA:N	2.37	0.54
2:B:137:ARG:HG2	2:B:138:GLY:H	1.73	0.54
5:E:20:LYS:HB3	5:E:35:VAL:HG23	1.90	0.54
15:O:516:LEU:CD2	15:O:567:ARG:CD	2.55	0.54
16:P:252:LYS:HD3	16:P:253:LEU:H	1.72	0.54
2:B:914:SER:HB3	2:B:918:GLN:HB2	1.89	0.53
3:C:270:ALA:HB3	3:C:271:ARG:HH11	1.73	0.53
4:D:125:ASN:OD1	4:D:126:GLN:N	2.40	0.53
14:N:398:SER:HB3	14:N:407:GLU:HB3	1.90	0.53
1:A:247:THR:HG23	1:A:249:PRO:HD3	1.90	0.53
2:B:163:LEU:HD22	2:B:171:MET:HE1	1.90	0.53
2:B:712:ALA:O	2:B:717:GLN:NE2	2.42	0.53
3:C:8:GLU:HB3	3:C:11:ARG:HG2	1.89	0.53
3:C:229:LEU:O	3:C:293:ARG:NH2	2.39	0.53
5:E:28:TYR:HA	5:E:64:PRO:HA	1.90	0.53
15:O:337:GLN:O	15:O:338:ASP:HB2	2.09	0.53
1:A:100:ILE:HG13	1:A:166:LYS:HD3	1.89	0.53
1:A:473:LEU:HA	1:A:487:SER:HA	1.90	0.53
1:A:1223:ASN:HB3	1:A:1231:ALA:HB3	1.91	0.53
1:A:1311:GLY:O	5:E:147:HIS:NE2	2.41	0.53
2:B:727:LEU:HD11	2:B:788:ARG:HE	1.73	0.53
7:G:89:ILE:HG23	7:G:142:VAL:HA	1.91	0.53
7:G:114:MET:O	7:G:201:GLN:HB3	2.08	0.53
1:A:1325:VAL:HG23	1:A:1326:LEU:H	1.74	0.53
14:N:290:ILE:HG13	14:N:293:LYS:HE2	1.90	0.53
15:O:50:LYS:HA	15:O:53:VAL:HG12	1.91	0.53
16:P:254:GLU:HB2	16:P:262:ARG:HB2	1.91	0.53
1:A:372:ARG:HG2	2:B:1049:GLN:HG2	1.90	0.53
15:O:292:ARG:HE	15:O:326:ILE:HD11	1.73	0.53
3:C:81:GLU:OE2	12:L:70:ARG:NH2	2.42	0.53
7:G:44:LEU:O	7:G:77:PHE:N	2.42	0.53
2:B:934:ASN:OD1	2:B:935:ASP:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:123:LEU:HD11	5:E:126:SER:HB3	1.91	0.53
13:M:122:ASP:HB3	13:M:145:VAL:HG21	1.90	0.53
1:A:1443:PRO:HG3	7:G:52:LEU:HD11	1.90	0.53
2:B:832:VAL:HB	12:L:60:ARG:HA	1.89	0.53
3:C:45:SER:HB2	3:C:53:ASN:HB3	1.91	0.53
1:A:200:GLU:HA	15:O:515:LYS:HE3	1.90	0.53
2:B:526:GLU:O	2:B:529:ILE:HG22	2.09	0.53
10:J:30:LEU:HD22	10:J:31:ASP:H	1.73	0.53
1:A:200:GLU:HG2	15:O:515:LYS:HD2	1.90	0.52
7:G:113:ASN:OD1	7:G:113:ASN:N	2.40	0.52
8:H:114:VAL:HG21	8:H:130:ARG:HH11	1.74	0.52
15:O:150:GLU:HG3	15:O:151:GLU:H	1.72	0.52
16:P:255:LYS:HE3	16:P:261:TYR:CE1	2.43	0.52
3:C:94:ASP:N	3:C:94:ASP:OD1	2.41	0.52
1:A:535:MET:O	1:A:540:ASN:ND2	2.42	0.52
2:B:289:GLU:HA	2:B:292:LYS:HG2	1.91	0.52
9:I:34:PRO:O	9:I:35:ILE:HB	2.09	0.52
1:A:378:ARG:HE	1:A:516:GLU:HB3	1.75	0.52
4:D:15:GLU:HA	4:D:18:LYS:HB3	1.90	0.52
5:E:161:LYS:NZ	5:E:193:GLY:O	2.37	0.52
15:O:338:ASP:HB3	15:O:342:ALA:N	2.25	0.52
1:A:1207:VAL:O	1:A:1211:ARG:HG2	2.09	0.52
2:B:417:ASP:HA	2:B:420:LEU:HG	1.90	0.52
15:O:343:LYS:HA	15:O:346:GLN:HB3	1.91	0.52
15:O:620:LEU:HD12	15:O:621:PRO:HD2	1.91	0.52
1:A:5:VAL:HG22	7:G:37:LYS:HB3	1.92	0.52
1:A:362:GLY:O	1:A:363:ARG:HB2	2.10	0.52
1:A:652:VAL:HG21	1:A:668:VAL:HG11	1.91	0.52
2:B:249:GLU:O	2:B:308:LYS:NZ	2.42	0.52
8:H:39:THR:N	8:H:124:ARG:O	2.41	0.52
2:B:299:GLN:HG3	13:M:186:ILE:HG22	1.90	0.52
4:D:57:GLY:O	4:D:61:ASN:ND2	2.43	0.52
16:P:228:ALA:HA	16:P:233:VAL:HG13	1.92	0.52
1:A:665:ASP:OD1	1:A:797:CYS:HB3	2.10	0.52
1:A:1221:ASP:OD1	1:A:1222:VAL:N	2.43	0.52
1:A:1408:VAL:HG23	1:A:1413:GLU:HG3	1.92	0.52
9:I:15:THR:HG23	9:I:23:THR:HG23	1.92	0.52
1:A:182:THR:HG22	1:A:185:TRP:HE1	1.74	0.52
2:B:667:VAL:O	2:B:669:SER:N	2.41	0.52
11:K:47:ILE:HG23	11:K:63:PHE:HB3	1.92	0.52
15:O:174:TYR:O	15:O:177:SER:OG	2.26	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:LEU:HD11	1:A:140:ILE:HG21	1.91	0.52
1:A:1140:ILE:HG22	1:A:1296:GLU:HG3	1.92	0.52
3:C:239:ILE:HG22	3:C:288:LYS:HE2	1.92	0.52
7:G:121:TYR:HB2	7:G:128:TRP:CZ3	2.45	0.52
1:A:1420:THR:HA	2:B:1080:LEU:HD11	1.92	0.51
3:C:132:ILE:H	3:C:208:CYS:HB3	1.75	0.51
1:A:988:ASP:HA	1:A:991:LYS:HB3	1.91	0.51
12:L:29:TYR:CE2	12:L:40:LEU:HB2	2.45	0.51
13:M:80:GLY:HA3	13:M:261:LYS:HE2	1.90	0.51
1:A:305:LYS:HG3	1:A:306:GLY:H	1.75	0.51
2:B:372:VAL:O	2:B:607:ARG:NH1	2.41	0.51
4:D:20:LEU:HD12	4:D:62:VAL:HG11	1.91	0.51
1:A:614:ILE:HD13	1:A:624:ILE:HD12	1.93	0.51
2:B:796:LYS:NZ	2:B:798:TYR:OH	2.44	0.51
5:E:88:VAL:HB	5:E:120:ALA:HB2	1.91	0.51
9:I:23:THR:OG1	9:I:24:LEU:N	2.44	0.51
13:M:89:GLN:HB3	14:N:394:VAL:HG22	1.93	0.51
1:A:714:ILE:HD13	2:B:958:MET:HE3	1.91	0.51
2:B:723:THR:HA	2:B:790:LYS:HG2	1.92	0.51
7:G:104:ILE:HG13	7:G:105:PHE:CD2	2.45	0.51
1:A:11:LYS:HG3	2:B:1117:ILE:HD13	1.92	0.51
1:A:244:ILE:HG21	1:A:253:PRO:HD3	1.91	0.51
1:A:1323:PHE:HA	1:A:1327:GLY:HA2	1.93	0.51
5:E:112:TYR:HE2	5:E:134:THR:HB	1.75	0.51
10:J:7:CYS:SG	10:J:11:GLY:N	2.77	0.51
1:A:252:ARG:HH12	15:O:46:LEU:HD22	1.74	0.51
1:A:896:LEU:HB3	1:A:1090:GLY:HA3	1.92	0.51
8:H:8:ASP:OD1	8:H:9:ILE:N	2.43	0.51
1:A:578:GLN:O	1:A:582:MET:N	2.43	0.51
2:B:333:ALA:HB1	2:B:345:LYS:HD3	1.91	0.51
8:H:116:TYR:HB2	8:H:123:MET:HB3	1.91	0.51
15:O:488:LYS:HZ1	15:O:651:PHE:HA	1.75	0.51
1:A:376:SER:HB2	2:B:1060:LEU:HD12	1.93	0.51
1:A:1143:ALA:HA	1:A:1314:THR:HG22	1.92	0.51
2:B:338:GLU:H	2:B:345:LYS:HZ3	1.58	0.51
16:P:309:LYS:HD3	17:Q:43:ILE:HD11	1.93	0.51
1:A:12:ARG:HG3	2:B:1146:ILE:HG12	1.92	0.51
1:A:1347:GLY:O	1:A:1348:MET:HG2	2.12	0.51
2:B:140:ASN:O	2:B:140:ASN:ND2	2.44	0.51
2:B:337:VAL:HG23	2:B:338:GLU:H	1.75	0.51
5:E:88:VAL:HG21	5:E:117:THR:H	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:200:ARG:NH1	5:E:210:SER:OG	2.44	0.51
15:O:192:VAL:HG12	15:O:274:VAL:H	1.76	0.51
1:A:482:ARG:HH12	1:A:924:LEU:HD11	1.75	0.50
4:D:3:VAL:HG13	7:G:7:ILE:HG22	1.92	0.50
7:G:124:GLU:O	7:G:125:GLU:HB3	2.12	0.50
1:A:905:ARG:HH22	5:E:170:LEU:HD21	1.76	0.50
2:B:698:ARG:HH21	2:B:952:ARG:HG2	1.76	0.50
15:O:190:LEU:HD23	15:O:194:LEU:HG	1.92	0.50
1:A:984:VAL:HG21	1:A:988:ASP:HB3	1.94	0.50
6:F:115:THR:HG22	6:F:116:ASP:H	1.76	0.50
15:O:195:CYS:SG	15:O:196:GLU:N	2.84	0.50
1:A:235:LYS:HD2	1:A:252:ARG:HH11	1.76	0.50
2:B:98:ILE:HG13	2:B:131:VAL:HG12	1.94	0.50
2:B:211:GLU:HG3	2:B:212:LYS:HD2	1.94	0.50
2:B:756:THR:HG23	2:B:941:ASP:H	1.76	0.50
7:G:156:VAL:HG11	7:G:190:LYS:HZ1	1.76	0.50
15:O:343:LYS:O	15:O:347:ASP:N	2.33	0.50
15:O:547:GLU:HG2	15:O:569:LYS:HG2	1.94	0.50
17:Q:48:THR:O	17:Q:51:GLU:N	2.44	0.50
1:A:524:THR:HG23	1:A:527:ALA:H	1.76	0.50
2:B:404:ASP:HA	2:B:407:LEU:HB2	1.92	0.50
1:A:1372:THR:HG23	1:A:1375:GLY:H	1.77	0.50
2:B:210:ASP:OD1	2:B:211:GLU:N	2.45	0.50
1:A:997:GLN:O	1:A:998:TYR:HB2	2.12	0.50
14:N:305:MET:HE3	14:N:414:GLY:HA3	1.93	0.50
15:O:265:VAL:HG12	15:O:267:PRO:HD3	1.94	0.50
1:A:372:ARG:HD3	2:B:1087:SER:HB3	1.93	0.50
1:A:723:LEU:HD21	1:A:811:ALA:HA	1.94	0.50
15:O:54:LYS:HA	15:O:58:GLY:HA2	1.94	0.50
16:P:314:PHE:C	16:P:316:GLU:N	2.69	0.50
2:B:487:ARG:HH21	2:B:508:CYS:HB3	1.76	0.49
2:B:615:VAL:HG12	2:B:620:SER:HA	1.94	0.49
5:E:115:ASN:OD1	5:E:116:ILE:N	2.45	0.49
15:O:500:LEU:HD23	15:O:504:ALA:HB2	1.94	0.49
2:B:121:ARG:HH12	12:L:54:ARG:HH21	1.58	0.49
14:N:389:THR:OG1	14:N:390:PHE:N	2.40	0.49
15:O:303:ARG:HD2	15:O:467:PHE:HE2	1.77	0.49
1:A:653:ILE:HG23	1:A:660:LEU:HB2	1.92	0.49
13:M:95:ARG:NH1	14:N:393:ASP:OD1	2.45	0.49
1:A:1273:LYS:HG2	1:A:1274:GLY:H	1.78	0.49
2:B:297:THR:HG23	2:B:298:GLN:HG2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:422:ILE:HA	2:B:425:HIS:HB3	1.94	0.49
7:G:23:ILE:HG22	7:G:51:LEU:HD23	1.95	0.49
8:H:105:GLU:HG2	8:H:115:TYR:HE1	1.77	0.49
12:L:29:TYR:HA	12:L:58:LYS:HA	1.93	0.49
15:O:592:LYS:HG3	15:O:637:VAL:HG11	1.93	0.49
2:B:296:TYR:HB2	2:B:300:GLN:OE1	2.13	0.49
2:B:312:MET:HB3	2:B:317:LEU:HG	1.93	0.49
2:B:391:LEU:HD23	2:B:429:ILE:HA	1.93	0.49
7:G:89:ILE:HD12	7:G:97:ILE:HB	1.94	0.49
2:B:989:LYS:HA	2:B:992:VAL:HG12	1.93	0.49
4:D:61:ASN:HB3	7:G:103:GLY:O	2.13	0.49
13:M:90:TYR:HB3	13:M:179:LEU:HB3	1.93	0.49
1:A:46:ARG:HG3	1:A:48:PRO:HG3	1.94	0.49
1:A:387:LEU:HD21	1:A:393:ALA:HB2	1.94	0.49
1:A:393:ALA:HB3	1:A:499:ARG:HG3	1.95	0.49
1:A:629:LYS:HB3	1:A:633:PHE:HE2	1.76	0.49
2:B:804:ASP:HB3	2:B:848:PRO:HD3	1.93	0.49
2:B:831:GLU:OE1	2:B:831:GLU:N	2.43	0.49
2:B:934:ASN:HA	2:B:1005:TYR:CD2	2.48	0.49
3:C:24:SER:HB2	3:C:27:ALA:HB2	1.93	0.49
5:E:213:ILE:HG12	5:E:214:CYS:H	1.77	0.49
13:M:142:GLY:O	13:M:143:VAL:HG12	2.12	0.49
15:O:145:ARG:O	15:O:146:VAL:HB	2.13	0.49
1:A:862:LEU:HD21	2:B:491:PRO:HA	1.95	0.49
1:A:1205:ILE:O	1:A:1209:ILE:N	2.42	0.49
3:C:191:ILE:HG23	10:J:15:GLY:HA3	1.93	0.49
13:M:96:LEU:O	13:M:97:VAL:HG12	2.13	0.49
1:A:408:VAL:HG23	1:A:412:ASN:HB2	1.95	0.49
14:N:304:PHE:O	14:N:413:ASP:N	2.31	0.49
14:N:392:GLN:HB3	14:N:394:VAL:HG23	1.93	0.49
2:B:97:ASP:OD2	2:B:99:ARG:NH1	2.41	0.48
2:B:371:TYR:HB2	2:B:492:SER:OG	2.13	0.48
2:B:543:LEU:HG	13:M:176:VAL:HG11	1.95	0.48
14:N:292:ARG:O	14:N:296:LYS:N	2.44	0.48
15:O:188:SER:HA	15:O:191:PHE:HB2	1.94	0.48
15:O:620:LEU:HD23	15:O:623:GLU:HB2	1.95	0.48
1:A:1064:GLU:HB3	1:A:1065:LYS:H	1.42	0.48
2:B:536:LEU:HD22	2:B:571:PHE:CD1	2.48	0.48
2:B:853:ASP:HB2	2:B:855:PRO:HD3	1.95	0.48
2:B:1022:ILE:HG22	2:B:1023:TYR:H	1.76	0.48
6:F:134:ILE:N	6:F:146:TRP:O	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:556:ASP:O	15:O:557:ARG:HG2	2.13	0.48
1:A:1284:ASN:OD1	1:A:1285:ILE:N	2.45	0.48
2:B:464:ILE:O	2:B:468:GLY:N	2.46	0.48
14:N:389:THR:O	14:N:390:PHE:HB3	2.12	0.48
15:O:326:ILE:HG23	15:O:653:MET:HE1	1.94	0.48
2:B:569:THR:OG1	14:N:275:GLU:OE2	2.29	0.48
3:C:331:CYS:SG	11:K:46:LYS:HB2	2.53	0.48
15:O:488:LYS:NZ	15:O:651:PHE:HA	2.28	0.48
16:P:248:VAL:O	16:P:255:LYS:NZ	2.33	0.48
1:A:60:VAL:HG23	1:A:74:LEU:HD23	1.96	0.48
1:A:372:ARG:HB3	2:B:1050:PRO:HD2	1.93	0.48
1:A:444:LEU:HD11	1:A:449:ARG:H	1.78	0.48
1:A:1451:LEU:HG	4:D:107:MET:HB3	1.96	0.48
5:E:170:LEU:HB3	5:E:174:GLN:NE2	2.29	0.48
14:N:308:GLN:HB2	14:N:417:VAL:HG12	1.94	0.48
1:A:62:SER:H	1:A:65:LEU:HD12	1.79	0.48
2:B:776:SER:HA	2:B:779:ASP:HB3	1.94	0.48
1:A:311:ASN:OD1	1:A:312:MET:N	2.41	0.48
1:A:1225:ILE:HG23	1:A:1226:GLY:H	1.78	0.48
5:E:86:PRO:HA	5:E:113:GLN:HB3	1.96	0.48
7:G:147:ARG:NH2	7:G:205:MET:SD	2.87	0.48
15:O:641:LEU:O	15:O:645:LEU:HB2	2.13	0.48
16:P:313:TYR:CB	17:Q:39:ILE:HD11	2.43	0.48
1:A:90:VAL:HG13	1:A:258:TRP:HB2	1.94	0.48
1:A:113:ILE:HG12	1:A:115:LEU:H	1.79	0.48
1:A:220:ASP:HA	15:O:550:GLU:HB3	1.95	0.48
1:A:980:SER:HB3	5:E:163:GLU:HG2	1.96	0.48
3:C:256:ILE:HG22	3:C:267:VAL:HA	1.96	0.48
15:O:553:ARG:HB2	15:O:562:ALA:HA	1.96	0.48
16:P:215:TYR:CZ	16:P:262:ARG:HD3	2.49	0.48
17:Q:41:LEU:HB3	17:Q:42:PRO:HD2	1.96	0.48
1:A:132:VAL:HG21	1:A:137:ARG:HH22	1.79	0.47
1:A:325:MET:HE3	1:A:350:ILE:HG22	1.96	0.47
1:A:953:GLY:HA2	1:A:1063:SER:HB2	1.95	0.47
1:A:1187:ARG:HG2	1:A:1229:ARG:HG3	1.95	0.47
4:D:110:LEU:HB3	4:D:120:LYS:NZ	2.28	0.47
2:B:823:SER:OG	12:L:63:ARG:NH1	2.42	0.47
2:B:1008:ILE:HG13	2:B:1009:THR:H	1.78	0.47
4:D:116:PHE:O	4:D:120:LYS:HB2	2.14	0.47
13:M:72:GLU:O	14:N:364:ARG:HG2	2.14	0.47
14:N:293:LYS:HA	14:N:296:LYS:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1145:LEU:HD21	1:A:1153:ALA:HB1	1.96	0.47
2:B:811:VAL:HG12	2:B:813:GLU:H	1.78	0.47
15:O:291:ARG:HA	15:O:294:LYS:HD3	1.95	0.47
15:O:493:GLU:OE1	16:P:249:TYR:OH	2.28	0.47
15:O:570:GLU:O	15:O:571:THR:OG1	2.29	0.47
1:A:113:ILE:HD13	1:A:120:LYS:HE2	1.97	0.47
1:A:221:ASP:OD1	15:O:550:GLU:HB2	2.13	0.47
2:B:217:GLN:HG2	2:B:232:TYR:CD1	2.49	0.47
2:B:550:HIS:NE2	2:B:551:LEU:HD23	2.29	0.47
3:C:74:GLU:HA	3:C:212:ILE:HD11	1.95	0.47
5:E:173:SER:C	5:E:175:LEU:H	2.23	0.47
13:M:107:ILE:HD12	13:M:124:PRO:HG3	1.95	0.47
1:A:395:PRO:HG2	1:A:398:VAL:HG22	1.96	0.47
1:A:1020:ALA:HA	1:A:1032:LEU:HD21	1.96	0.47
1:A:34:VAL:HG23	1:A:35:SER:H	1.79	0.47
1:A:706:GLY:HA2	2:B:762:TYR:HA	1.97	0.47
2:B:121:ARG:NH1	2:B:123:MET:SD	2.88	0.47
2:B:775:LYS:HE3	2:B:927:LYS:HA	1.96	0.47
2:B:854:ALA:N	2:B:855:PRO:HD3	2.29	0.47
2:B:1000:GLY:HA3	2:B:1018:PHE:HD1	1.80	0.47
4:D:98:MET:SD	4:D:160:TYR:OH	2.71	0.47
15:O:322:LYS:HB3	15:O:361:PHE:CE1	2.50	0.47
16:P:314:PHE:CG	16:P:315:ASP:N	2.82	0.47
16:P:315:ASP:O	16:P:317:TRP:HD1	1.97	0.47
1:A:666:LYS:O	1:A:667:SER:HB3	2.15	0.47
1:A:1202:ILE:HD12	1:A:1228:ASP:HA	1.97	0.47
2:B:613:ILE:HG12	2:B:646:VAL:HG12	1.96	0.47
2:B:727:LEU:HB2	2:B:786:GLU:HB2	1.97	0.47
3:C:43:ASN:HB2	3:C:55:ASP:HB3	1.95	0.47
8:H:130:ARG:O	8:H:134:ASN:ND2	2.39	0.47
13:M:78:ILE:HG22	13:M:170:LEU:HD22	1.96	0.47
1:A:598:MET:CA	1:A:600:PRO:HD3	2.43	0.47
12:L:28:LYS:HG3	12:L:29:TYR:CD2	2.48	0.47
15:O:259:LEU:HD21	15:O:261:GLN:HB2	1.97	0.47
15:O:541:ILE:HG13	15:O:546:VAL:HG23	1.97	0.47
1:A:1190:LEU:HA	1:A:1193:ILE:HD12	1.97	0.47
2:B:741:ILE:HB	2:B:746:TYR:HB3	1.97	0.47
2:B:849:THR:HG23	2:B:850:ASN:H	1.79	0.47
5:E:143:ASN:OD1	5:E:143:ASN:N	2.47	0.47
15:O:40:ARG:O	16:P:317:TRP:NE1	2.47	0.47
1:A:133:ASP:OD1	1:A:134:ASN:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ASP:OD1	1:A:220:ASP:N	2.48	0.47
1:A:890:MET:O	1:A:894:GLU:N	2.48	0.47
1:A:1134:LYS:HA	1:A:1320:LEU:HD23	1.96	0.47
2:B:108:THR:OG1	2:B:109:LYS:N	2.47	0.47
1:A:364:PHE:CD1	1:A:364:PHE:C	2.94	0.46
1:A:383:PRO:HB3	1:A:502:GLU:HG2	1.97	0.46
2:B:83:ILE:HG13	2:B:93:LEU:HB3	1.96	0.46
5:E:70:SER:OG	5:E:71:LYS:N	2.47	0.46
7:G:151:GLU:HB3	7:G:198:GLY:HA2	1.97	0.46
15:O:255:LYS:HB3	15:O:256:PRO:HD3	1.96	0.46
16:P:253:LEU:HB3	16:P:261:TYR:HB3	1.96	0.46
2:B:465:SER:OG	2:B:707:LYS:O	2.34	0.46
2:B:545:ASP:CG	14:N:391:LEU:HD22	2.40	0.46
7:G:88:TRP:HB3	7:G:145:LYS:HA	1.97	0.46
14:N:290:ILE:HD13	14:N:384:LYS:HE3	1.97	0.46
15:O:203:ILE:HG13	15:O:203:ILE:O	2.13	0.46
15:O:282:ILE:HG22	15:O:283:ASN:H	1.80	0.46
15:O:516:LEU:O	15:O:565:LEU:HD23	2.16	0.46
16:P:221:ILE:O	16:P:225:ILE:N	2.39	0.46
16:P:255:LYS:HD2	16:P:255:LYS:N	2.30	0.46
1:A:664:MET:HB2	1:A:669:LEU:HD13	1.97	0.46
1:A:1145:LEU:HD23	1:A:1146:VAL:HG23	1.97	0.46
2:B:269:MET:O	2:B:273:CYS:HB3	2.15	0.46
2:B:1047:THR:HG22	2:B:1049:GLN:HG3	1.97	0.46
14:N:375:ILE:HG22	14:N:376:GLY:H	1.80	0.46
1:A:913:GLN:NE2	1:A:915:THR:O	2.49	0.46
2:B:1080:LEU:O	2:B:1084:MET:HB3	2.16	0.46
7:G:80:PHE:CE2	7:G:83:GLU:HB2	2.50	0.46
15:O:516:LEU:CD2	15:O:567:ARG:HG3	2.44	0.46
2:B:687:LEU:O	2:B:915:ARG:NH2	2.48	0.46
5:E:54:GLN:HG3	5:E:57:MET:HE2	1.97	0.46
11:K:112:THR:HG22	11:K:113:ALA:H	1.81	0.46
15:O:87:ASP:O	15:O:88:VAL:HG12	2.16	0.46
2:B:806:ILE:H	2:B:806:ILE:HG13	1.60	0.46
1:A:1391:LYS:HD2	1:A:1391:LYS:HA	1.70	0.46
1:A:1441:LEU:HD21	7:G:54:VAL:HG12	1.97	0.46
7:G:88:TRP:O	7:G:100:SER:N	2.30	0.46
7:G:144:GLU:OE2	7:G:207:LEU:HG	2.16	0.46
13:M:255:PHE:O	13:M:259:ILE:HG12	2.16	0.46
2:B:77:ILE:HD13	2:B:98:ILE:HB	1.98	0.46
2:B:103:LYS:HG2	2:B:104:SER:H	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:736:VAL:HG21	2:B:960:GLU:HG3	1.97	0.46
2:B:992:VAL:HG23	3:C:278:GLU:HG2	1.98	0.46
3:C:3:ASN:HD21	3:C:296:ASN:HB2	1.80	0.46
5:E:55:ARG:HD2	5:E:82:PHE:HE2	1.81	0.46
5:E:117:THR:OG1	5:E:118:PRO:HD2	2.15	0.46
13:M:131:TYR:OH	13:M:143:VAL:HG23	2.16	0.46
15:O:467:PHE:C	15:O:468:LEU:HG	2.40	0.46
15:O:512:ARG:CG	15:O:568:CYS:SG	2.97	0.46
1:A:47:ALA:N	1:A:48:PRO:HD3	2.31	0.46
1:A:308:SER:OG	1:A:309:ILE:N	2.49	0.46
1:A:415:LYS:HD3	6:F:115:THR:HG21	1.97	0.46
1:A:1145:LEU:HG	1:A:1309:VAL:HG12	1.97	0.46
3:C:175:GLN:O	3:C:178:THR:HG23	2.16	0.46
4:D:130:ASN:HB3	4:D:133:HIS:CE1	2.51	0.46
13:M:95:ARG:HD2	13:M:101:PRO:HB2	1.97	0.46
13:M:123:ILE:C	13:M:145:VAL:HG23	2.40	0.46
15:O:516:LEU:CG	15:O:567:ARG:HG3	2.45	0.46
1:A:597:ILE:HG22	1:A:598:MET:H	1.81	0.46
1:A:869:ARG:HH21	2:B:489:LEU:HD12	1.80	0.46
2:B:69:VAL:HA	2:B:73:LEU:HD23	1.97	0.46
2:B:181:PRO:HB2	2:B:470:MET:HE1	1.97	0.46
2:B:244:HIS:ND1	2:B:246:SER:OG	2.43	0.46
7:G:162:SER:O	7:G:166:ARG:N	2.42	0.46
9:I:95:THR:HA	9:I:110:ASN:HA	1.97	0.46
16:P:306:LYS:HB3	16:P:307:HIS:CD2	2.50	0.46
3:C:193:LEU:HD12	3:C:193:LEU:H	1.80	0.45
1:A:22:SER:C	1:A:24:ALA:H	2.21	0.45
1:A:645:MET:HE1	8:H:124:ARG:HD2	1.98	0.45
1:A:1033:GLU:HG2	1:A:1034:PRO:HD2	1.98	0.45
2:B:916:HIS:CD2	2:B:957:LYS:HB2	2.52	0.45
3:C:29:ASN:ND2	3:C:35:LYS:HE2	2.32	0.45
3:C:34:GLU:HA	3:C:37:LYS:HB3	1.98	0.45
4:D:10:PHE:HD2	4:D:117:LYS:HE3	1.80	0.45
5:E:29:PHE:HB2	5:E:65:THR:HB	1.98	0.45
7:G:97:ILE:HG22	7:G:128:TRP:NE1	2.31	0.45
14:N:395:ILE:HG12	14:N:408:LEU:HG	1.99	0.45
15:O:466:PRO:HG2	15:O:467:PHE:HD2	1.81	0.45
2:B:1081:GLU:OE2	2:B:1086:SER:OG	2.34	0.45
16:P:264:THR:HG23	16:P:265:LEU:H	1.81	0.45
1:A:542:LEU:HD21	1:A:679:TYR:HE1	1.81	0.45
2:B:92:TYR:CZ	2:B:136:THR:HG21	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:804:ASP:OD1	2:B:804:ASP:N	2.50	0.45
9:I:98:TYR:HB2	9:I:107:TRP:HE3	1.81	0.45
11:K:97:SER:OG	11:K:98:GLU:N	2.50	0.45
1:A:308:SER:N	1:A:311:ASN:HD21	2.15	0.45
1:A:444:LEU:HD21	1:A:449:ARG:HG2	1.97	0.45
1:A:712:ILE:HG22	1:A:713:GLY:H	1.81	0.45
2:B:241:TYR:HB3	2:B:250:GLU:HB3	1.97	0.45
2:B:517:MET:HE2	2:B:611:PRO:HG2	1.99	0.45
2:B:933:PHE:O	2:B:934:ASN:ND2	2.50	0.45
8:H:21:ASN:OD1	8:H:22:LYS:N	2.40	0.45
15:O:201:ILE:HG22	15:O:202:GLN:H	1.81	0.45
16:P:300:LYS:O	16:P:301:MET:HB2	2.16	0.45
1:A:1365:LYS:HE3	1:A:1379:MET:HB2	1.98	0.45
16:P:253:LEU:O	16:P:255:LYS:NZ	2.43	0.45
1:A:386:ASN:ND2	11:K:94:PRO:O	2.50	0.45
2:B:325:GLU:HA	2:B:328:ALA:HB3	1.99	0.45
3:C:229:LEU:HB3	3:C:293:ARG:CZ	2.46	0.45
16:P:312:VAL:HG11	17:Q:41:LEU:HB2	1.98	0.45
1:A:541:LEU:HD22	1:A:682:LEU:HD22	1.99	0.45
1:A:545:LYS:HG3	1:A:546:SER:H	1.82	0.45
2:B:139:ARG:HB3	2:B:141:ILE:HD11	1.99	0.45
2:B:343:ARG:HE	2:B:542:THR:HA	1.82	0.45
2:B:409:LYS:HA	2:B:410:PRO:HD3	1.85	0.45
3:C:280:LEU:HA	3:C:286:ALA:HB2	1.98	0.45
3:C:283:GLU:O	3:C:286:ALA:N	2.50	0.45
8:H:7:ASP:OD1	8:H:58:THR:OG1	2.34	0.45
1:A:163:VAL:HG22	1:A:164:VAL:H	1.81	0.45
5:E:47:CYS:HA	5:E:53:PRO:HA	1.97	0.45
14:N:300:LYS:N	14:N:301:PRO:HD2	2.32	0.45
1:A:360:LYS:HE3	1:A:361:GLN:HG3	1.99	0.45
2:B:240:ILE:HG13	2:B:253:ILE:HG21	1.98	0.45
4:D:31:GLN:HA	4:D:34:LEU:HD12	1.98	0.45
6:F:143:PHE:O	6:F:144:GLU:HB2	2.17	0.45
13:M:88:PHE:HE2	14:N:397:LEU:HD22	1.81	0.45
15:O:191:PHE:O	15:O:194:LEU:HB2	2.17	0.45
15:O:509:ARG:HA	15:O:512:ARG:HB2	1.99	0.45
1:A:235:LYS:HA	1:A:235:LYS:HD3	1.77	0.44
2:B:230:LYS:HD2	2:B:334:HIS:HB3	1.99	0.44
2:B:260:CYS:O	2:B:342:PHE:HB3	2.17	0.44
2:B:889:SER:OG	2:B:893:GLN:HG2	2.17	0.44
3:C:328:LEU:HG	11:K:121:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:20:LEU:HD21	7:G:47:THR:HG21	1.98	0.44
9:I:35:ILE:HD12	9:I:35:ILE:HA	1.87	0.44
13:M:122:ASP:HB3	13:M:145:VAL:CG2	2.47	0.44
1:A:183:PHE:CD2	1:A:184:ARG:HG3	2.52	0.44
1:A:475:ASN:HB3	1:A:518:ASN:HB2	1.99	0.44
1:A:979:ASN:O	5:E:167:ARG:NH2	2.50	0.44
1:A:1026:ARG:NH1	8:H:129:TYR:OH	2.50	0.44
1:A:1151:GLU:HG3	1:A:1152:ARG:H	1.82	0.44
2:B:88:ASP:CG	2:B:89:PRO:HD2	2.41	0.44
2:B:795:LEU:HB2	2:B:894:ALA:HB3	1.99	0.44
2:B:1043:ARG:HG2	2:B:1050:PRO:HD3	1.98	0.44
3:C:25:LYS:HE2	8:H:85:GLY:HA2	2.00	0.44
5:E:40:GLU:HG3	5:E:41:ASP:H	1.82	0.44
15:O:489:SER:O	15:O:493:GLU:HG2	2.17	0.44
15:O:581:LEU:HD13	15:O:648:TRP:CE2	2.52	0.44
1:A:1123:VAL:HG13	1:A:1124:PRO:HD3	1.98	0.44
2:B:803:GLN:HG2	2:B:804:ASP:H	1.82	0.44
2:B:932:PRO:HB3	2:B:1004:LEU:HB3	2.00	0.44
5:E:172:GLU:HG2	5:E:173:SER:H	1.82	0.44
1:A:200:GLU:HB3	15:O:516:LEU:HD21	2.00	0.44
2:B:983:LYS:HG3	2:B:985:GLU:H	1.83	0.44
3:C:92:ILE:H	3:C:92:ILE:HG13	1.56	0.44
7:G:122:THR:OG1	7:G:126:SER:O	2.34	0.44
10:J:12:LYS:HE3	10:J:43:ARG:NH2	2.32	0.44
13:M:108:SER:O	13:M:122:ASP:HB2	2.17	0.44
1:A:93:ILE:HG23	1:A:94:GLY:H	1.83	0.44
1:A:817:ILE:HG13	1:A:821:ASN:O	2.18	0.44
2:B:38:ILE:HD11	2:B:43:ASP:HB3	2.00	0.44
2:B:1067:ARG:NH1	2:B:1079:LEU:HD22	2.33	0.44
5:E:195:VAL:HG22	5:E:213:ILE:HG13	1.99	0.44
7:G:159:LYS:N	7:G:160:PRO:HD3	2.32	0.44
13:M:73:GLU:C	14:N:364:ARG:HE	2.26	0.44
15:O:289:LYS:HD3	15:O:325:LYS:HB3	1.99	0.44
1:A:627:ASP:OD1	1:A:627:ASP:N	2.51	0.44
1:A:862:LEU:HD22	2:B:494:PHE:HB2	1.98	0.44
7:G:52:LEU:O	7:G:53:THR:HG22	2.17	0.44
15:O:150:GLU:HG3	15:O:151:GLU:N	2.32	0.44
15:O:163:VAL:HA	15:O:169:LEU:HD13	1.98	0.44
15:O:516:LEU:HD22	15:O:567:ARG:NE	1.91	0.44
15:O:528:MET:HG3	16:P:175:ILE:HG13	2.00	0.44
1:A:1323:PHE:HD1	1:A:1328:ILE:H	1.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:506:ARG:HH12	15:O:526:ALA:HB1	1.81	0.44
1:A:297:SER:HA	1:A:300:LYS:HB3	2.00	0.44
1:A:626:LEU:HD21	1:A:675:HIS:HB2	2.00	0.44
1:A:653:ILE:HG12	1:A:660:LEU:HD22	1.99	0.44
1:A:664:MET:HE2	1:A:664:MET:HB3	1.93	0.44
1:A:809:MET:HE1	2:B:950:PRO:HA	1.99	0.44
1:A:1455:ALA:HB2	4:D:120:LYS:HZ3	1.81	0.44
2:B:1067:ARG:HH22	2:B:1079:LEU:HB3	1.83	0.44
5:E:24:LYS:HB2	5:E:30:ILE:HD11	1.98	0.44
15:O:553:ARG:C	15:O:555:ALA:H	2.25	0.44
1:A:367:ASN:O	1:A:371:LYS:HB2	2.18	0.44
2:B:224:THR:OG1	2:B:225:HIS:N	2.50	0.44
2:B:717:GLN:OE1	2:B:727:LEU:HD22	2.16	0.44
2:B:1008:ILE:O	3:C:65:ASN:ND2	2.51	0.44
3:C:85:PHE:CE1	3:C:204:LEU:HD22	2.53	0.44
3:C:270:ALA:HB3	3:C:271:ARG:NH1	2.32	0.44
14:N:296:LYS:HD3	14:N:296:LYS:HA	1.78	0.44
14:N:366:HIS:N	14:N:370:LYS:O	2.51	0.44
1:A:179:ILE:HG12	1:A:221:ASP:HB3	2.00	0.43
1:A:1411:VAL:HG13	1:A:1412:SER:H	1.83	0.43
2:B:325:GLU:O	2:B:329:THR:OG1	2.26	0.43
2:B:929:GLU:HB3	3:C:69:ARG:HG2	2.00	0.43
7:G:149:ARG:O	7:G:198:GLY:HA3	2.18	0.43
12:L:60:ARG:HG2	12:L:61:THR:H	1.83	0.43
15:O:242:LYS:NZ	15:O:341:GLU:OE2	2.45	0.43
1:A:481:HIS:CD2	1:A:483:LEU:HB2	2.52	0.43
2:B:87:VAL:HG12	2:B:407:LEU:HD22	2.00	0.43
4:D:70:LYS:HG3	4:D:98:MET:SD	2.59	0.43
6:F:85:MET:HE2	6:F:151:LEU:HD22	1.99	0.43
7:G:47:THR:OG1	7:G:48:ILE:N	2.50	0.43
1:A:126:GLU:OE2	1:A:136:ARG:NH2	2.38	0.43
1:A:431:ASN:OD1	1:A:465:HIS:NE2	2.51	0.43
1:A:661:SER:HB2	8:H:122:LEU:HD11	2.01	0.43
1:A:985:LYS:O	1:A:987:GLU:N	2.51	0.43
1:A:1300:LEU:O	1:A:1303:VAL:HG12	2.17	0.43
1:A:19:SER:O	2:B:1138:ALA:N	2.52	0.43
1:A:634:VAL:HA	1:A:635:PRO:HD3	1.82	0.43
1:A:774:ARG:HH22	1:A:804:LEU:HD11	1.84	0.43
1:A:862:LEU:O	1:A:866:ILE:HG12	2.19	0.43
1:A:1059:LEU:HD22	1:A:1060:TYR:CE1	2.53	0.43
2:B:190:THR:OG1	2:B:191:GLU:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:29:ASN:HB2	3:C:35:LYS:NZ	2.33	0.43
2:B:538:VAL:HG12	2:B:565:ILE:HD12	1.99	0.43
5:E:2:ASP:O	5:E:6:GLU:HG2	2.18	0.43
8:H:12:VAL:HG12	8:H:13:SER:H	1.82	0.43
15:O:329:PRO:C	15:O:331:THR:H	2.23	0.43
15:O:533:ILE:HD13	15:O:533:ILE:HA	1.93	0.43
1:A:42:LEU:HD12	1:A:42:LEU:H	1.84	0.43
1:A:424:PRO:HG3	1:A:444:LEU:O	2.18	0.43
1:A:988:ASP:O	1:A:992:ALA:N	2.42	0.43
2:B:86:ASP:O	2:B:87:VAL:HG23	2.18	0.43
2:B:373:GLY:HA2	2:B:607:ARG:NH2	2.34	0.43
2:B:722:ASP:OD1	2:B:723:THR:N	2.47	0.43
2:B:757:VAL:O	2:B:1019:PHE:HB2	2.19	0.43
1:A:272:VAL:HB	1:A:274:MET:HG2	2.01	0.43
2:B:110:ASP:OD1	2:B:110:ASP:N	2.52	0.43
2:B:225:HIS:HA	2:B:446:ARG:HH11	1.83	0.43
3:C:59:ILE:HB	3:C:63:ILE:HD11	2.00	0.43
9:I:85:LEU:HD23	9:I:85:LEU:HA	1.87	0.43
15:O:327:ARG:NH1	15:O:330:LEU:HD22	2.33	0.43
15:O:581:LEU:HD13	15:O:648:TRP:CD2	2.53	0.43
1:A:49:LYS:HG3	1:A:56:PRO:HD3	2.01	0.43
1:A:353:PHE:O	1:A:356:ARG:HG2	2.19	0.43
2:B:124:THR:OG1	2:B:125:TYR:N	2.51	0.43
2:B:905:ARG:O	2:B:907:GLU:N	2.51	0.43
7:G:152:ARG:HB3	7:G:197:LEU:HG	2.00	0.43
8:H:44:VAL:HG13	8:H:45:GLU:H	1.84	0.43
14:N:373:VAL:HG12	14:N:374:LYS:H	1.83	0.43
15:O:74:LEU:HD21	15:O:79:LEU:HG	2.01	0.43
17:Q:59:ILE:HG22	17:Q:63:LYS:HE3	2.00	0.43
1:A:1097:ILE:HD13	1:A:1358:LEU:HD22	2.01	0.43
2:B:444:LEU:O	2:B:448:LYS:HA	2.19	0.43
2:B:650:ASP:OD1	2:B:651:VAL:N	2.47	0.43
3:C:255:VAL:O	3:C:268:LYS:HB2	2.19	0.43
5:E:20:LYS:O	5:E:23:VAL:HG12	2.19	0.43
13:M:135:LYS:HE3	13:M:140:TRP:CH2	2.54	0.43
13:M:247:TRP:HA	14:N:406:ALA:HB3	2.01	0.43
15:O:53:VAL:HG21	15:O:65:ILE:HD12	2.00	0.43
1:A:182:THR:HG22	1:A:185:TRP:NE1	2.35	0.42
1:A:493:ARG:HB3	1:A:494:PRO:HD2	2.01	0.42
1:A:955:LEU:HA	1:A:956:PRO:HD2	1.89	0.42
4:D:113:PHE:CE2	4:D:153:MET:HG2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:7:ILE:O	7:G:71:THR:OG1	2.27	0.42
15:O:492:TYR:HE1	15:O:576:PHE:HE1	1.66	0.42
1:A:387:LEU:O	1:A:501:ASN:ND2	2.53	0.42
1:A:1302:ASP:OD1	1:A:1303:VAL:N	2.52	0.42
2:B:417:ASP:O	2:B:421:SER:HB2	2.19	0.42
2:B:720:ARG:NH2	2:B:722:ASP:OD2	2.41	0.42
2:B:1028:LYS:HG2	2:B:1029:HIS:H	1.83	0.42
4:D:139:GLU:C	4:D:141:CYS:H	2.28	0.42
5:E:43:LYS:O	5:E:47:CYS:HB2	2.18	0.42
15:O:500:LEU:HD12	15:O:500:LEU:HA	1.87	0.42
15:O:597:GLN:HA	15:O:600:SER:HB2	2.00	0.42
1:A:350:ILE:HD12	1:A:350:ILE:HA	1.91	0.42
1:A:1145:LEU:O	1:A:1310:ILE:HG22	2.20	0.42
1:A:1308:GLY:HA2	9:I:88:ARG:HA	2.00	0.42
2:B:113:THR:HA	2:B:114:PRO:HD3	1.88	0.42
2:B:137:ARG:NH2	2:B:413:ALA:O	2.38	0.42
2:B:418:ALA:O	2:B:422:ILE:HG12	2.19	0.42
2:B:775:LYS:C	2:B:777:SER:H	2.27	0.42
2:B:914:SER:C	2:B:916:HIS:H	2.27	0.42
7:G:94:ALA:HB2	7:G:121:TYR:CD2	2.54	0.42
15:O:159:ILE:O	15:O:163:VAL:HG23	2.19	0.42
15:O:197:MET:HG3	15:O:198:GLY:H	1.84	0.42
15:O:208:TYR:O	15:O:209:THR:OG1	2.37	0.42
1:A:409:THR:C	1:A:411:TYR:H	2.28	0.42
1:A:598:MET:O	1:A:600:PRO:HD3	2.11	0.42
2:B:378:GLU:OE2	2:B:382:GLN:HB2	2.20	0.42
2:B:551:LEU:HD22	14:N:388:THR:H	1.85	0.42
3:C:263:ASP:OD1	3:C:264:GLU:N	2.51	0.42
8:H:96:VAL:HG12	8:H:143:LEU:HA	2.02	0.42
1:A:238:ASP:OD1	1:A:238:ASP:N	2.53	0.42
1:A:468:ASP:OD1	1:A:469:GLY:N	2.53	0.42
7:G:104:ILE:HG23	7:G:105:PHE:N	2.32	0.42
15:O:275:LYS:HA	15:O:276:PRO:HD3	1.90	0.42
16:P:237:PRO:O	16:P:240:ILE:HB	2.19	0.42
1:A:1148:ASP:OD1	1:A:1291:ARG:NH1	2.50	0.42
2:B:263:LEU:HD12	2:B:297:THR:HG22	2.00	0.42
2:B:292:LYS:HG3	2:B:293:LEU:H	1.85	0.42
7:G:115:LEU:HD23	7:G:116:PHE:N	2.33	0.42
8:H:3:ASN:ND2	8:H:62:SER:OG	2.38	0.42
15:O:201:ILE:O	15:O:281:THR:N	2.48	0.42
15:O:338:ASP:O	15:O:342:ALA:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:576:PHE:O	15:O:579:GLN:HG2	2.19	0.42
1:A:1080:LYS:HD3	1:A:1080:LYS:HA	1.94	0.42
1:A:1166:LEU:O	1:A:1170:ALA:HB2	2.20	0.42
2:B:474:SER:HA	2:B:511:VAL:HG12	2.02	0.42
2:B:529:ILE:HG23	2:B:530:LYS:N	2.35	0.42
2:B:735:MET:HB2	2:B:754:ASN:HD21	1.85	0.42
7:G:55:GLU:OE1	7:G:55:GLU:N	2.53	0.42
13:M:95:ARG:NH2	14:N:392:GLN:O	2.38	0.42
16:P:310:GLU:C	16:P:311:VAL:HG23	2.44	0.42
1:A:978:ASP:N	1:A:982:CYS:O	2.48	0.42
1:A:1173:VAL:HG22	1:A:1186:VAL:HG12	2.01	0.42
4:D:17:LEU:O	4:D:21:THR:OG1	2.20	0.42
7:G:29:GLN:O	7:G:33:LYS:HG2	2.20	0.42
7:G:59:LEU:HD13	7:G:65:SER:O	2.19	0.42
1:A:959:ILE:HG23	1:A:1067:VAL:HG21	2.01	0.42
1:A:1200:LEU:HB2	1:A:1204:ASP:HB2	2.02	0.42
1:A:1428:PHE:CZ	6:F:89:GLU:HA	2.55	0.42
2:B:131:VAL:HG21	2:B:149:ILE:HG23	2.01	0.42
2:B:755:ALA:HA	10:J:48:ARG:NH1	2.35	0.42
4:D:26:LYS:HE2	4:D:26:LYS:HB2	1.95	0.42
4:D:145:PHE:HB2	4:D:149:THR:OG1	2.19	0.42
7:G:41:ASN:HA	7:G:155:PHE:CE2	2.54	0.42
15:O:262:ILE:HG12	15:O:278:VAL:HG21	2.02	0.42
16:P:297:ASN:O	16:P:298:TYR:HB3	2.19	0.42
1:A:409:THR:O	1:A:411:TYR:N	2.53	0.42
2:B:447:PHE:HB3	2:B:449:MET:HG2	2.01	0.42
2:B:1088:ASP:OD1	2:B:1088:ASP:N	2.51	0.42
3:C:77:SER:O	3:C:210:LEU:HD12	2.20	0.42
4:D:24:GLU:OE2	4:D:30:ASP:HB2	2.20	0.42
8:H:50:ALA:N	8:H:53:ASP:OD2	2.53	0.42
15:O:325:LYS:HA	15:O:325:LYS:HD2	1.77	0.42
2:B:915:ARG:HD2	2:B:1023:TYR:HD2	1.85	0.41
3:C:105:PRO:HG3	10:J:6:ARG:HH21	1.85	0.41
13:M:77:LYS:HB3	13:M:77:LYS:HE2	1.79	0.41
16:P:309:LYS:HD3	17:Q:43:ILE:CD1	2.49	0.41
3:C:87:ASN:O	3:C:88:ASN:HB2	2.20	0.41
4:D:27:HIS:C	4:D:29:TRP:H	2.27	0.41
11:K:46:LYS:O	11:K:47:ILE:O	2.39	0.41
14:N:395:ILE:HD12	14:N:396:ALA:H	1.84	0.41
15:O:123:ASN:C	15:O:124:GLU:HG2	2.45	0.41
15:O:193:GLN:NE2	15:O:196:GLU:OE1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:617:ASN:O	1:A:620:SER:OG	2.35	0.41
2:B:321:GLN:O	2:B:324:ILE:HG22	2.20	0.41
3:C:90:SER:HB2	3:C:200:GLN:HB3	2.01	0.41
4:D:132:VAL:O	4:D:135:TYR:HB3	2.20	0.41
1:A:247:THR:C	1:A:249:PRO:HD3	2.45	0.41
1:A:984:VAL:HG22	1:A:985:LYS:H	1.86	0.41
1:A:1442:VAL:HA	1:A:1443:PRO:HD3	1.77	0.41
2:B:217:GLN:NE2	2:B:352:MET:HG2	2.34	0.41
2:B:1036:HIS:NE2	2:B:1058:GLY:HA3	2.35	0.41
2:B:1054:ARG:O	2:B:1055:SER:OG	2.31	0.41
3:C:104:VAL:HA	3:C:105:PRO:HD3	1.89	0.41
8:H:93:TYR:CD2	8:H:143:LEU:HG	2.56	0.41
8:H:131:ASN:OD1	8:H:131:ASN:N	2.52	0.41
9:I:6:PRO:O	9:I:7:SER:OG	2.26	0.41
11:K:59:THR:HB	11:K:60:SER:H	1.67	0.41
15:O:640:ARG:NH2	17:Q:43:ILE:HG13	2.35	0.41
16:P:229:GLN:C	16:P:231:ALA:H	2.27	0.41
1:A:610:PHE:HZ	1:A:697:MET:HE1	1.86	0.41
1:A:967:LEU:HD11	1:A:1009:ARG:HD3	2.03	0.41
5:E:82:PHE:CE1	5:E:111:VAL:HB	2.55	0.41
5:E:152:LYS:HE2	5:E:152:LYS:HB2	1.96	0.41
13:M:88:PHE:CD2	14:N:397:LEU:HD13	2.53	0.41
15:O:349:ALA:O	15:O:352:VAL:HG12	2.20	0.41
1:A:1301:ARG:O	1:A:1305:CYS:N	2.45	0.41
3:C:42:VAL:HG13	11:K:138:LYS:HD3	2.03	0.41
4:D:58:ILE:H	4:D:58:ILE:HG13	1.74	0.41
8:H:108:SER:OG	8:H:109:LYS:N	2.53	0.41
13:M:121:ILE:HD13	13:M:121:ILE:HA	1.91	0.41
1:A:1153:ALA:HA	1:A:1156:VAL:HB	2.02	0.41
1:A:1184:ILE:O	1:A:1231:ALA:HA	2.21	0.41
2:B:944:MET:HE2	2:B:944:MET:HB3	1.94	0.41
4:D:14:TYR:OH	4:D:99:SER:O	2.35	0.41
15:O:583:TRP:HE1	16:P:315:ASP:HB3	1.84	0.41
1:A:18:PHE:HA	2:B:1138:ALA:O	2.20	0.41
1:A:372:ARG:CD	2:B:1087:SER:HB3	2.51	0.41
1:A:1365:LYS:HE2	1:A:1365:LYS:HB2	1.97	0.41
3:C:276:SER:O	3:C:277:ARG:HG2	2.20	0.41
5:E:112:TYR:CE2	5:E:134:THR:HB	2.55	0.41
14:N:387:GLU:CD	14:N:387:GLU:H	2.29	0.41
15:O:105:LYS:HD2	15:O:209:THR:N	2.36	0.41
1:A:1083:LEU:HD23	1:A:1083:LEU:HA	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:60:GLN:H	2:B:60:GLN:CD	2.28	0.41
2:B:565:ILE:HG12	2:B:566:ARG:N	2.36	0.41
2:B:1063:GLY:O	2:B:1067:ARG:N	2.54	0.41
3:C:52:ALA:HB3	3:C:302:VAL:HG12	2.02	0.41
4:D:130:ASN:CG	4:D:131:MET:HG2	2.45	0.41
5:E:1:MET:HB3	5:E:4:GLU:HG2	2.02	0.41
8:H:41:ASP:HB2	8:H:121:LEU:HB3	2.03	0.41
9:I:7:SER:OG	9:I:8:CYS:N	2.53	0.41
9:I:35:ILE:HG13	9:I:36:GLU:H	1.86	0.41
10:J:42:LYS:HG3	10:J:43:ARG:H	1.86	0.41
11:K:83:ASN:HA	11:K:84:PRO:HD3	1.89	0.41
14:N:297:MET:SD	14:N:303:ARG:NH1	2.94	0.41
15:O:286:ARG:HA	15:O:289:LYS:HB2	2.02	0.41
1:A:285:LEU:HD23	1:A:285:LEU:HA	1.91	0.41
1:A:946:THR:HG21	1:A:1066:SER:HA	2.03	0.41
2:B:841:ILE:HG21	2:B:870:PRO:HB2	2.03	0.41
2:B:948:GLY:HA2	2:B:952:ARG:NH1	2.36	0.41
3:C:256:ILE:HA	3:C:268:LYS:H	1.86	0.41
7:G:60:LYS:HA	7:G:61:PRO:HD3	1.89	0.41
8:H:17:PRO:HB3	8:H:24:CYS:HB2	2.03	0.41
8:H:76:THR:C	8:H:78:SER:H	2.29	0.41
13:M:126:ASP:HB3	13:M:128:GLN:HG2	2.03	0.41
15:O:195:CYS:HB2	15:O:263:LEU:HG	2.03	0.41
15:O:226:ILE:HG23	15:O:229:ASN:H	1.86	0.41
1:A:630:ASN:ND2	1:A:650:GLY:O	2.53	0.40
1:A:1147:ASN:ND2	1:A:1149:ASN:OD1	2.54	0.40
2:B:247:ILE:HD11	2:B:308:LYS:HZ1	1.86	0.40
2:B:548:SER:HA	2:B:550:HIS:ND1	2.36	0.40
2:B:610:ARG:HA	2:B:611:PRO:HD3	1.77	0.40
2:B:811:VAL:HA	2:B:817:PRO:HB3	2.03	0.40
2:B:934:ASN:HB3	2:B:1004:LEU:HA	2.01	0.40
3:C:127:THR:HG22	3:C:128:ASP:H	1.85	0.40
11:K:65:ILE:O	11:K:101:LEU:N	2.54	0.40
15:O:329:PRO:O	15:O:331:THR:N	2.44	0.40
1:A:597:ILE:HG22	1:A:598:MET:N	2.36	0.40
1:A:1439:LYS:HA	1:A:1439:LYS:HE2	2.03	0.40
2:B:579:ARG:HG3	2:B:584:VAL:O	2.21	0.40
3:C:238:PRO:O	3:C:239:ILE:HG13	2.21	0.40
7:G:8:ALA:HA	7:G:71:THR:HA	2.02	0.40
8:H:58:THR:O	8:H:143:LEU:N	2.48	0.40
13:M:160:ALA:HA	14:N:306:VAL:HA	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:148:ASP:OD1	15:O:148:ASP:N	2.46	0.40
15:O:596:LYS:O	15:O:600:SER:N	2.54	0.40
16:P:311:VAL:C	16:P:312:VAL:HG12	2.47	0.40
1:A:12:ARG:NH2	2:B:1144:GLU:OE2	2.55	0.40
1:A:620:SER:HA	1:A:621:PRO:HD3	1.87	0.40
1:A:1428:PHE:CE2	6:F:92:ARG:HD2	2.57	0.40
2:B:985:GLU:O	2:B:988:SER:OG	2.29	0.40
2:B:658:TYR:CE2	2:B:670:MET:HG2	2.57	0.40
2:B:890:ASP:N	2:B:890:ASP:OD1	2.54	0.40
5:E:170:LEU:HB3	5:E:174:GLN:HE21	1.86	0.40
16:P:221:ILE:HA	16:P:221:ILE:HD13	1.98	0.40
1:A:506:THR:OG1	1:A:507:PRO:HD3	2.22	0.40
1:A:521:VAL:HA	1:A:522:PRO:HD3	1.91	0.40
1:A:568:ASP:OD1	8:H:22:LYS:HG2	2.21	0.40
1:A:800:LYS:HG3	2:B:947:HIS:O	2.21	0.40
2:B:93:LEU:HD12	2:B:134:GLU:O	2.20	0.40
2:B:497:LEU:HD23	2:B:514:LEU:HD23	2.03	0.40
13:M:253:GLU:H	13:M:253:GLU:CD	2.28	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	1402/1460 (96%)	1157 (82%)	233 (17%)	12 (1%)	14	49
2	B	1112/1149 (97%)	948 (85%)	155 (14%)	9 (1%)	16	53
3	C	333/335 (99%)	282 (85%)	47 (14%)	4 (1%)	10	42
4	D	113/161 (70%)	80 (71%)	33 (29%)	0	100	100
5	E	213/215 (99%)	172 (81%)	38 (18%)	3 (1%)	9	39
6	F	81/155 (52%)	73 (90%)	8 (10%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	G	185/212 (87%)	154 (83%)	25 (14%)	6 (3%)	3	21
8	H	136/146 (93%)	116 (85%)	20 (15%)	0	100	100
9	I	88/110 (80%)	77 (88%)	10 (11%)	1 (1%)	11	45
10	J	65/70 (93%)	53 (82%)	11 (17%)	1 (2%)	8	38
11	K	99/142 (70%)	83 (84%)	15 (15%)	1 (1%)	12	47
12	L	43/70 (61%)	33 (77%)	10 (23%)	0	100	100
13	M	161/282 (57%)	135 (84%)	22 (14%)	4 (2%)	4	26
14	N	101/422 (24%)	78 (77%)	22 (22%)	1 (1%)	12	47
15	O	522/654 (80%)	428 (82%)	88 (17%)	6 (1%)	11	45
16	P	92/317 (29%)	60 (65%)	26 (28%)	6 (6%)	1	13
17	Q	33/104 (32%)	29 (88%)	4 (12%)	0	100	100
All	All	4779/6004 (80%)	3958 (83%)	767 (16%)	54 (1%)	14	45

All (54) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	587	ILE
1	A	632	VAL
1	A	1371	ILE
9	I	35	ILE
11	K	47	ILE
13	M	97	VAL
13	M	107	ILE
13	M	143	VAL
15	O	88	VAL
15	O	146	VAL
16	P	175	ILE
16	P	230	VAL
1	A	362	GLY
2	B	87	VAL
3	C	239	ILE
16	P	233	VAL
1	A	307	ILE
1	A	975	VAL
2	B	811	VAL
2	B	927	LYS
3	C	255	VAL
14	N	411	ARG

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Mol	Chain	Res	Type
16	P	308	ASP
16	P	315	ASP
1	A	23	ALA
1	A	1328	ILE
2	B	632	ASP
2	B	651	VAL
2	B	713	ILE
2	B	767	ILE
3	C	87	ASN
3	C	91	VAL
15	O	338	ASP
16	P	309	LYS
1	A	1389	PHE
7	G	79	PRO
7	G	80	PHE
7	G	191	PRO
10	J	65	PRO
15	O	330	LEU
15	O	554	THR
1	A	666	LYS
5	E	122	LYS
15	O	285	ASP
5	E	88	VAL
1	A	278	PRO
7	G	158	VAL
13	M	114	PRO
1	A	600	PRO
2	B	584	VAL
5	E	90	VAL
7	G	192	PRO
2	B	337	VAL
7	G	160	PRO

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	1225/1257 (98%)	1183 (97%)	42 (3%)	32	54
2	B	975/1006 (97%)	937 (96%)	38 (4%)	28	49
3	C	296/296 (100%)	284 (96%)	12 (4%)	27	49
4	D	110/145 (76%)	106 (96%)	4 (4%)	31	52
5	E	197/197 (100%)	189 (96%)	8 (4%)	27	49
6	F	73/137 (53%)	70 (96%)	3 (4%)	27	49
7	G	170/190 (90%)	163 (96%)	7 (4%)	27	49
8	H	123/128 (96%)	118 (96%)	5 (4%)	27	49
9	I	83/98 (85%)	83 (100%)	0	100	100
10	J	62/65 (95%)	57 (92%)	5 (8%)	11	31
11	K	91/130 (70%)	89 (98%)	2 (2%)	45	64
12	L	40/57 (70%)	37 (92%)	3 (8%)	12	33
13	M	143/249 (57%)	137 (96%)	6 (4%)	26	48
14	N	88/360 (24%)	83 (94%)	5 (6%)	18	40
15	O	486/593 (82%)	463 (95%)	23 (5%)	23	45
16	P	94/285 (33%)	87 (93%)	7 (7%)	13	33
17	Q	31/56 (55%)	28 (90%)	3 (10%)	8	25
All	All	4287/5249 (82%)	4114 (96%)	173 (4%)	29	49

All (173) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	34	VAL
1	A	104	GLN
1	A	124	LEU
1	A	132	VAL
1	A	160	LEU
1	A	164	VAL
1	A	222	LEU
1	A	247	THR
1	A	289	LEU
1	A	327	ILE
1	A	332	VAL
1	A	381	ILE
1	A	390	ASP
1	A	408	VAL

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Mol	Chain	Res	Type
1	A	409	THR
1	A	542	LEU
1	A	712	ILE
1	A	747	LYS
1	A	813	VAL
1	A	817	ILE
1	A	848	VAL
1	A	904	VAL
1	A	955	LEU
1	A	1022	LEU
1	A	1064	GLU
1	A	1123	VAL
1	A	1141	ILE
1	A	1161	VAL
1	A	1162	GLU
1	A	1186	VAL
1	A	1193	ILE
1	A	1224	ILE
1	A	1271	VAL
1	A	1272	VAL
1	A	1294	LEU
1	A	1303	VAL
1	A	1310	ILE
1	A	1350	VAL
1	A	1408	VAL
1	A	1411	VAL
1	A	1423	ILE
2	B	87	VAL
2	B	133	ILE
2	B	141	ILE
2	B	147	VAL
2	B	186	ILE
2	B	191	GLU
2	B	231	THR
2	B	234	ILE
2	B	251	ILE
2	B	295	ILE
2	B	335	LEU
2	B	337	VAL
2	B	366	ILE
2	B	377	LEU
2	B	383	LEU

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Mol	Chain	Res	Type
2	B	422	ILE
2	B	457	VAL
2	B	536	LEU
2	B	538	VAL
2	B	545	ASP
2	B	548	SER
2	B	555	VAL
2	B	565	ILE
2	B	599	VAL
2	B	612	LEU
2	B	666	ILE
2	B	675	ILE
2	B	713	ILE
2	B	741	ILE
2	B	772	VAL
2	B	806	ILE
2	B	818	ILE
2	B	860	VAL
2	B	871	VAL
2	B	899	LEU
2	B	938	ILE
2	B	954	THR
2	B	1022	ILE
3	C	30	GLU
3	C	33	VAL
3	C	59	ILE
3	C	78	VAL
3	C	91	VAL
3	C	120	LEU
3	C	133	VAL
3	C	151	THR
3	C	177	THR
3	C	239	ILE
3	C	274	THR
3	C	294	VAL
4	D	109	LYS
4	D	121	LEU
4	D	127	LEU
4	D	134	LEU
5	E	80	VAL
5	E	90	VAL
5	E	102	GLU

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Mol	Chain	Res	Type
5	E	107	THR
5	E	109	ILE
5	E	124	VAL
5	E	141	VAL
5	E	196	VAL
6	F	115	THR
6	F	133	VAL
6	F	151	LEU
7	G	3	ILE
7	G	23	ILE
7	G	53	THR
7	G	93	THR
7	G	108	ILE
7	G	113	ASN
7	G	115	LEU
8	H	12	VAL
8	H	26	ILE
8	H	32	THR
8	H	44	VAL
8	H	59	ILE
10	J	3	VAL
10	J	24	LEU
10	J	30	LEU
10	J	39	LEU
10	J	54	VAL
11	K	59	THR
11	K	112	THR
12	L	30	ILE
12	L	34	CYS
12	L	48	CYS
13	M	83	GLU
13	M	87	VAL
13	M	97	VAL
13	M	107	ILE
13	M	125	LEU
13	M	143	VAL
14	N	313	LEU
14	N	363	ILE
14	N	373	VAL
14	N	412	VAL
14	N	416	ILE
15	O	33	THR

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Mol	Chain	Res	Type
15	O	42	LEU
15	O	64	VAL
15	O	88	VAL
15	O	112	ILE
15	O	131	LEU
15	O	169	LEU
15	O	186	THR
15	O	201	ILE
15	O	223	TYR
15	O	226	ILE
15	O	264	THR
15	O	271	LEU
15	O	282	ILE
15	O	336	LEU
15	O	352	VAL
15	O	468	LEU
15	O	516	LEU
15	O	565	LEU
15	O	567	ARG
15	O	569	LYS
15	O	644	LEU
15	O	645	LEU
16	P	175	ILE
16	P	180	THR
16	P	233	VAL
16	P	235	LEU
16	P	236	THR
16	P	240	ILE
16	P	264	THR
17	Q	39	ILE
17	Q	41	LEU
17	Q	47	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	207	HIS
1	A	412	ASN
1	A	422	ASN
1	A	437	ASN
1	A	475	ASN
1	A	523	GLN

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Mol	Chain	Res	Type
1	A	843	GLN
1	A	934	ASN
1	A	1233	ASN
1	A	1354	HIS
2	B	299	GLN
2	B	552	ASN
2	B	626	HIS
2	B	655	ASN
2	B	702	GLN
2	B	717	GLN
2	B	865	GLN
3	C	29	ASN
3	C	175	GLN
4	D	27	HIS
4	D	122	GLN
5	E	3	GLN
5	E	5	ASN
5	E	54	GLN
5	E	174	GLN
6	F	104	ASN
9	I	71	ASN
9	I	105	HIS
9	I	110	ASN
13	M	254	GLN
14	N	287	HIS
15	O	147	ASN
15	O	244	ASN
15	O	337	GLN
15	O	346	GLN
15	O	544	ASN
15	O	575	ASN
15	O	580	ASN
15	O	652	GLN
16	P	176	ASN
16	P	307	HIS

5.3.3 RNA ⓘ

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 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
17	Q	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	Q	1088:UNK	C	1090:UNK	N	15.14
1	Q	69:PRO	C	1070:UNK	N	8.39

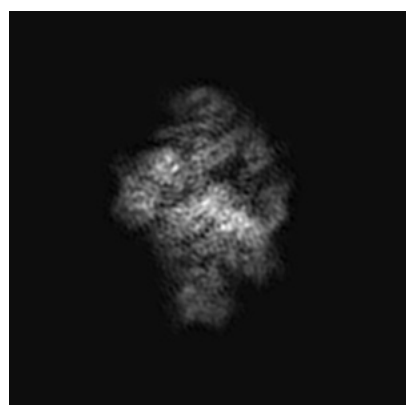
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-3179. These allow visual inspection of the internal detail of the map and identification of artifacts.

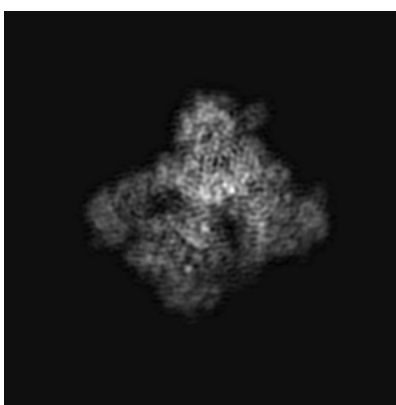
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

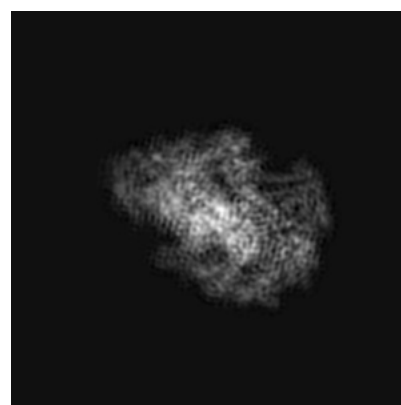
6.1.1 Primary map



X



Y

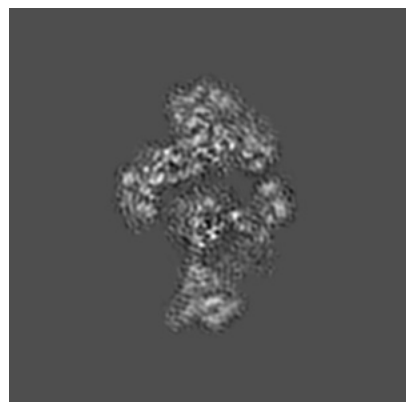


Z

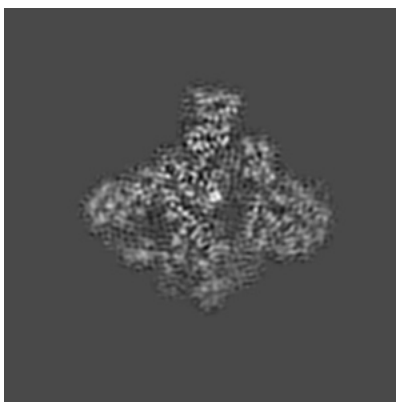
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

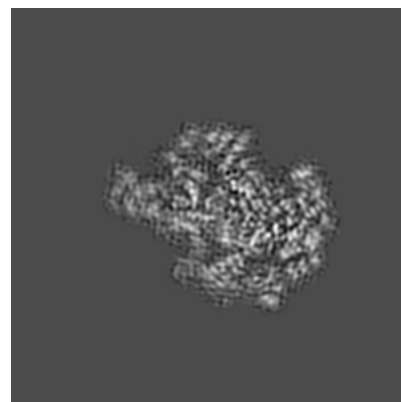
6.2.1 Primary map



X Index: 148



Y Index: 148

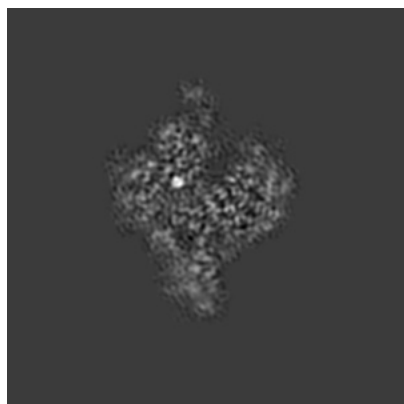


Z Index: 148

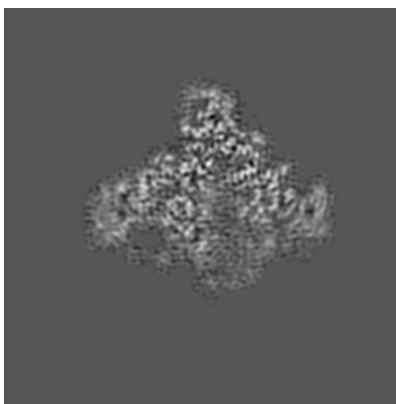
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

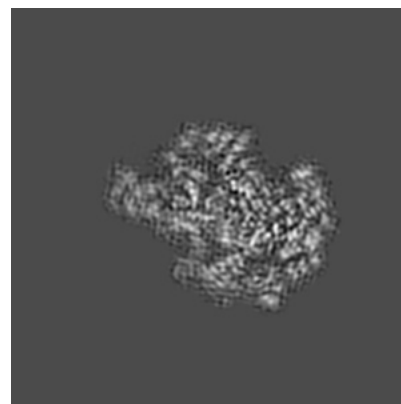
6.3.1 Primary map



X Index: 164



Y Index: 139

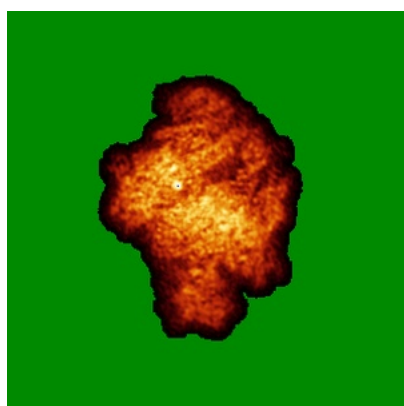


Z Index: 148

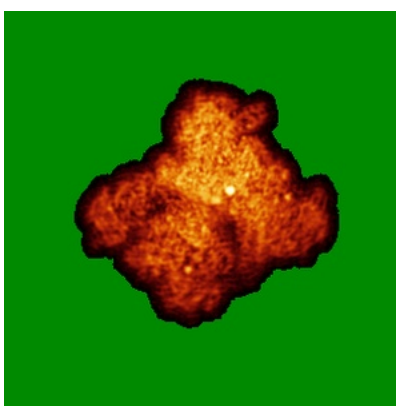
The images above show the largest variance slices of the map in three orthogonal directions.

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

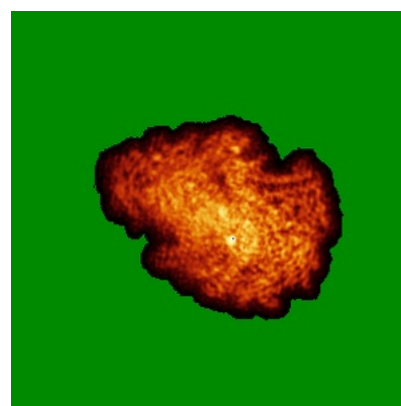
6.4.1 Primary map



X



Y

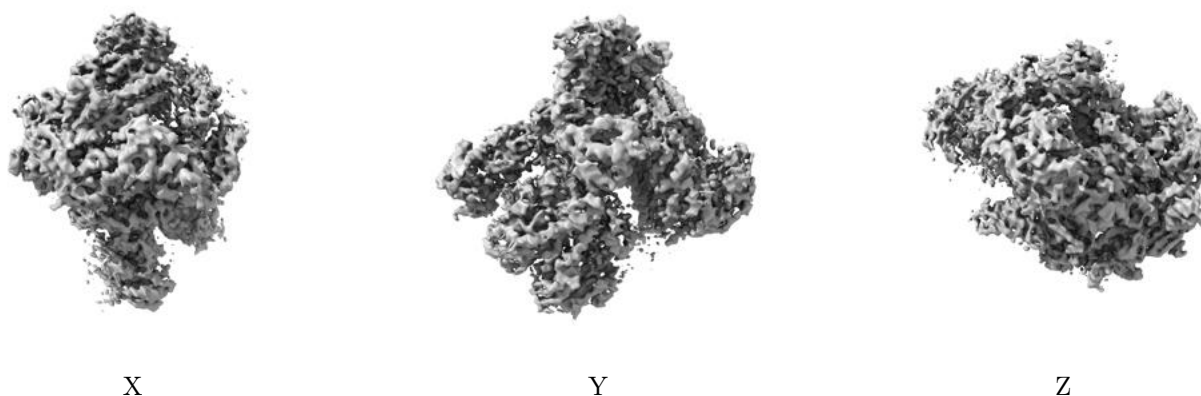


Z

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 0.055. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

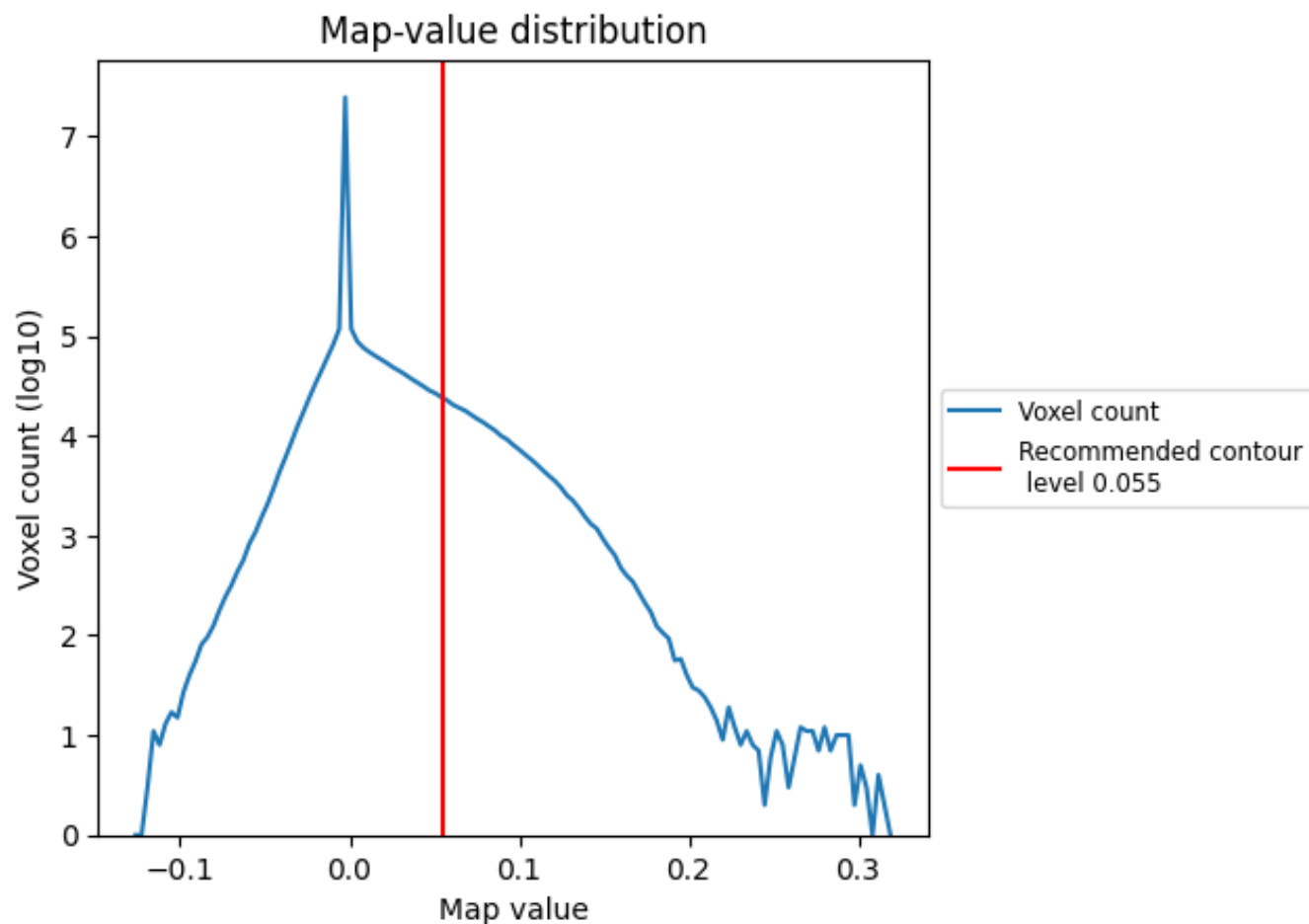
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

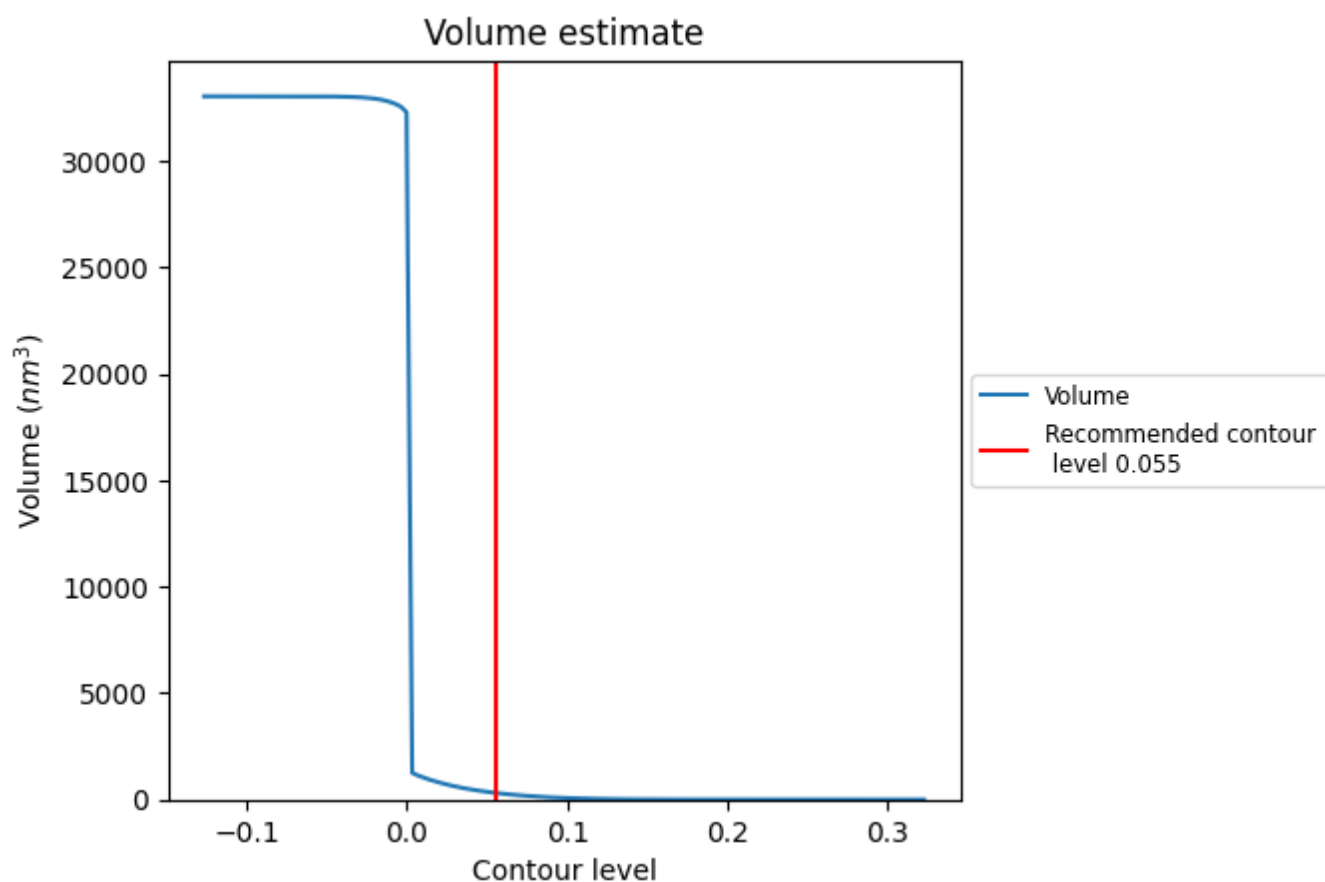
This section contains the results of statistical analysis of the map.

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.

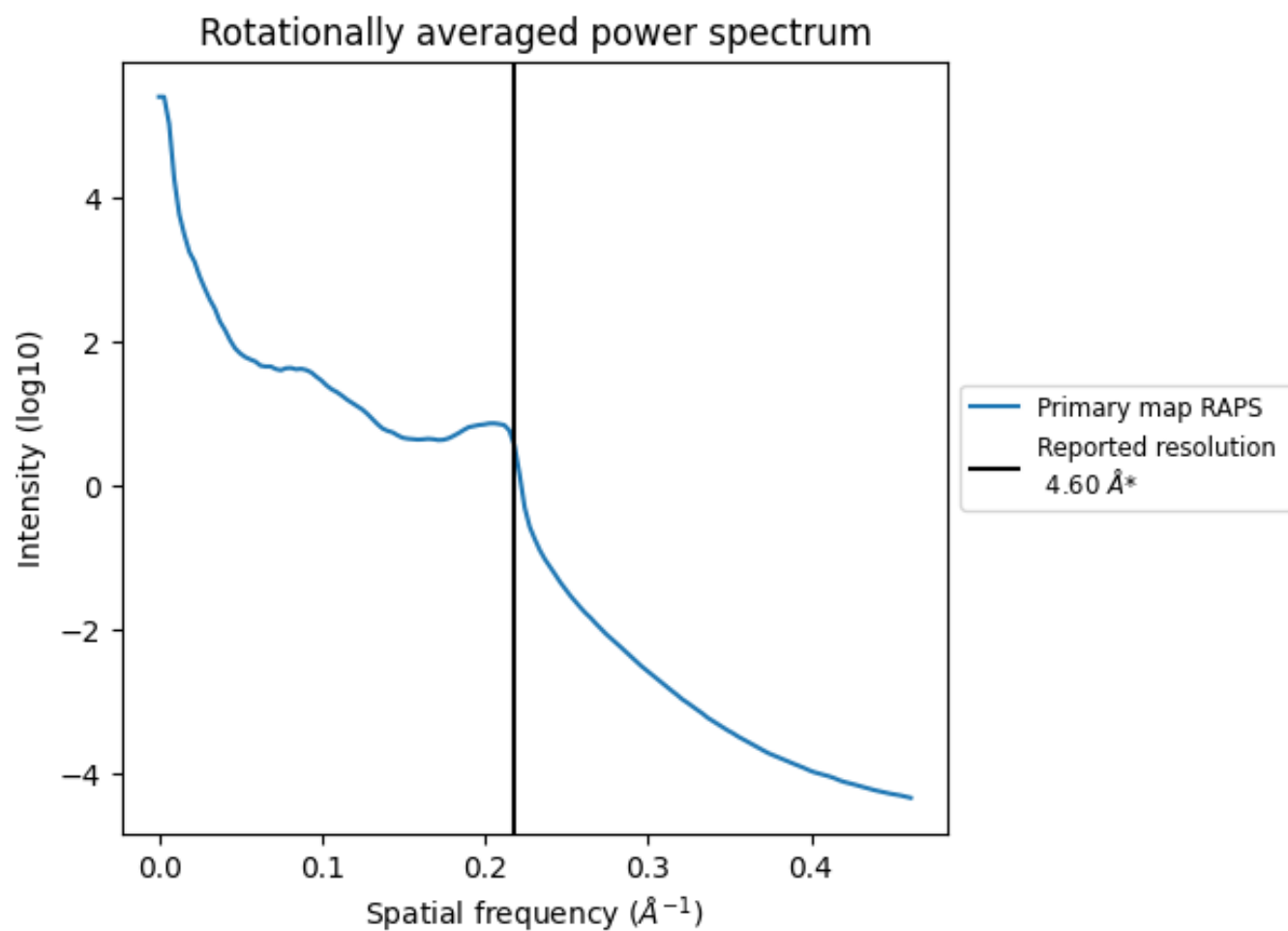
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 315 nm³; this corresponds to an approximate mass of 285 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 ⓘ

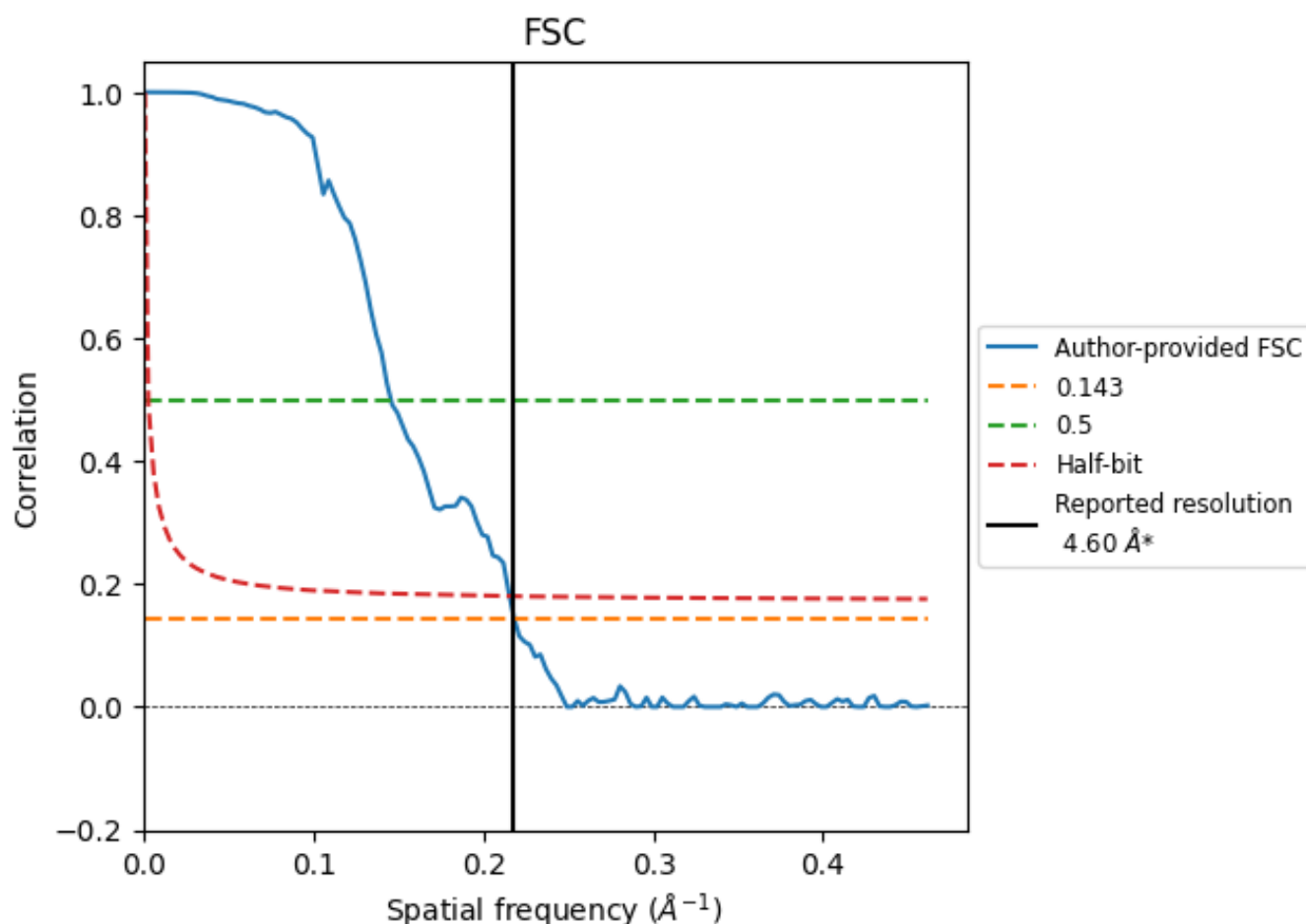


*Reported resolution corresponds to spatial frequency of 0.217 Å⁻¹

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 0.217 Å⁻¹

8.2 Resolution estimates [i](#)

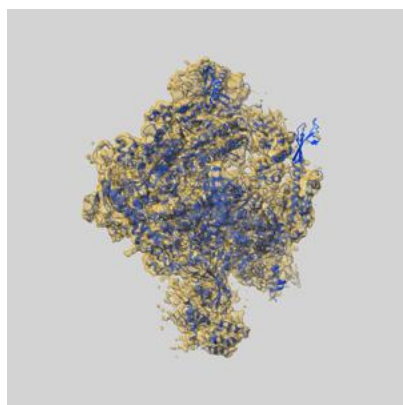
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	-	-	-
Author-provided FSC curve	4.58	6.86	4.64
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

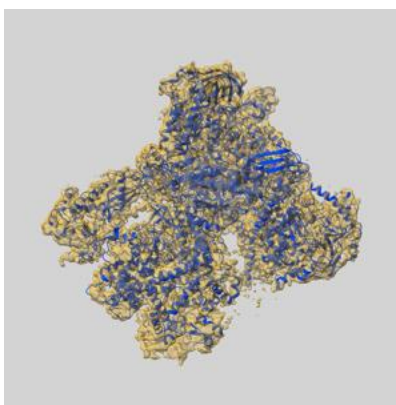
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-3179 and PDB model 5FJ9. Per-residue inclusion information can be found in section 3 on page 7.

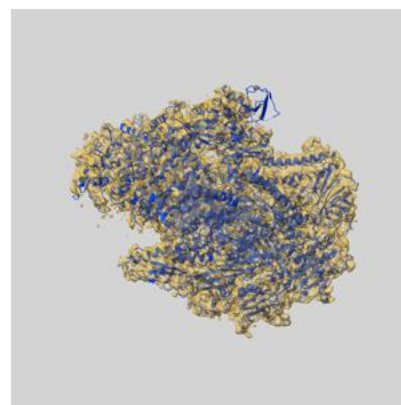
9.1 Map-model overlay [i](#)



X



Y



Z

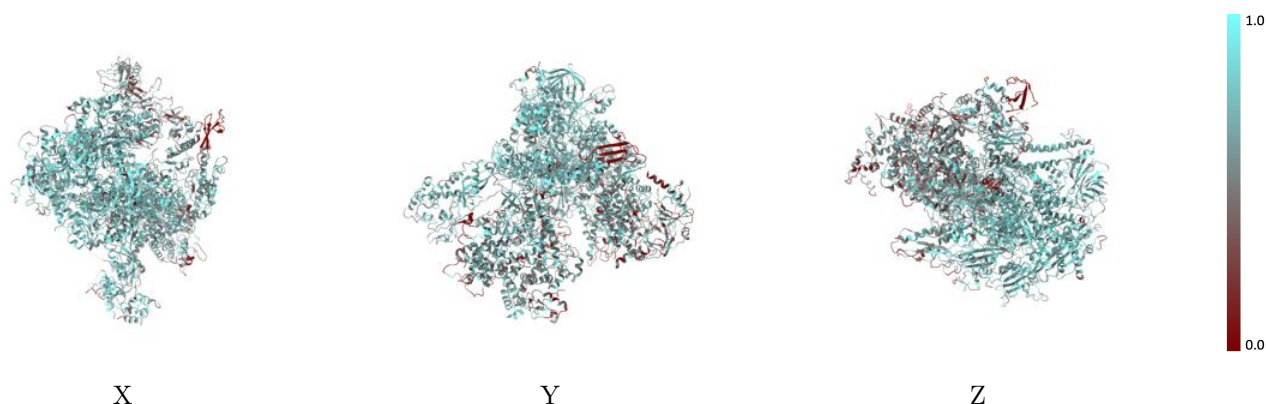
The images above show the 3D surface view of the map at the recommended contour level 0.055 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



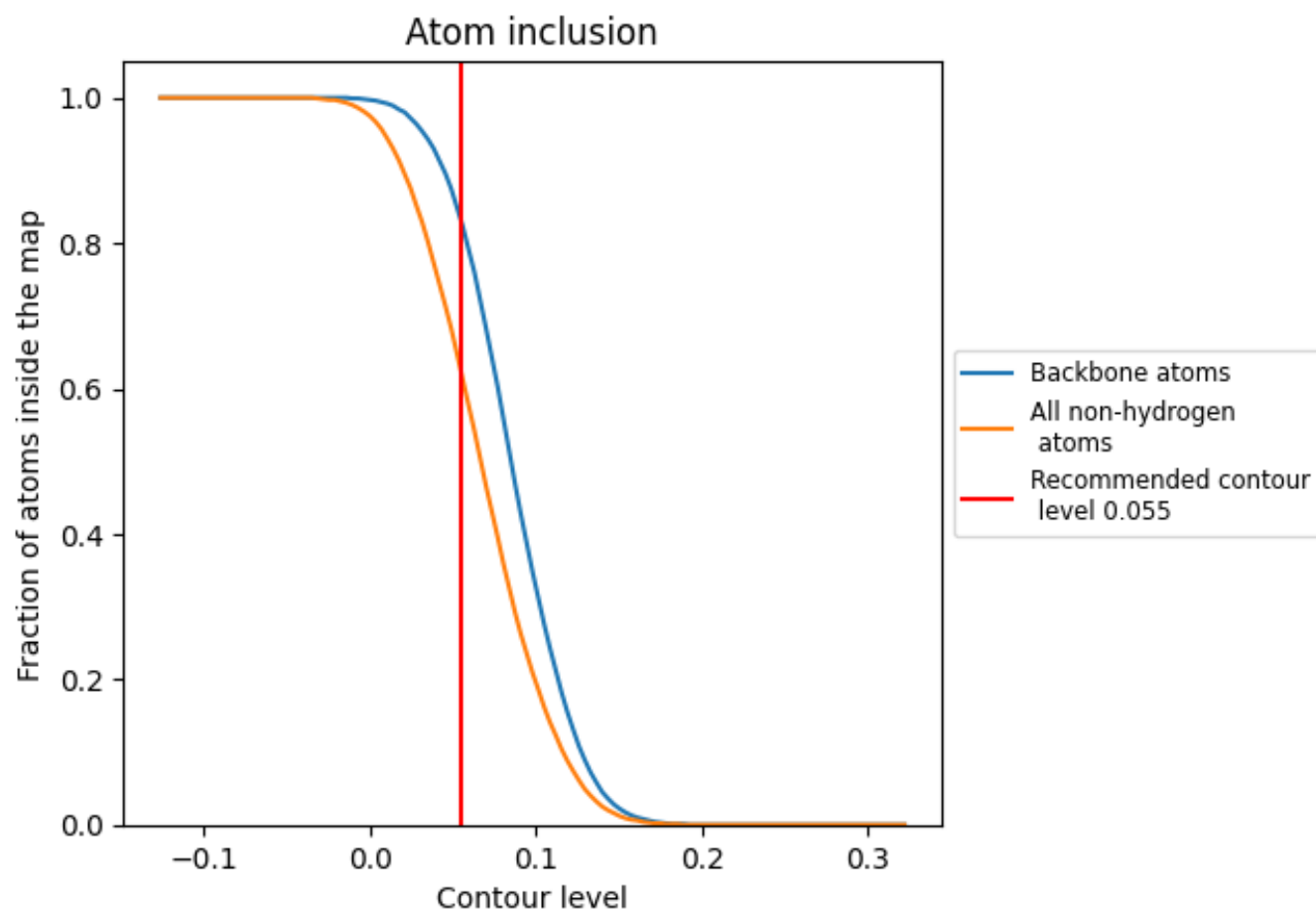
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.055).

9.4 Atom inclusion ⓘ



At the recommended contour level, 83% of all backbone atoms, 62% 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 (0.055) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6230	 0.2480
A	 0.6460	 0.2650
B	 0.6530	 0.2820
C	 0.7070	 0.2880
D	 0.6110	 0.1860
E	 0.6500	 0.2410
F	 0.7270	 0.3350
G	 0.6220	 0.2270
H	 0.6880	 0.2780
I	 0.2140	 0.1050
J	 0.7300	 0.2940
K	 0.6900	 0.2940
L	 0.6500	 0.2430
M	 0.4770	 0.1820
N	 0.4720	 0.1830
O	 0.5560	 0.1840
P	 0.4800	 0.1130
Q	 0.5440	 0.2160

