Full wwPDB EM Validation Report ⓘ

May 7, 2024 – 02:59 pm BST

PDB ID	:	8S55
EMDB ID	:	EMD-19726
Title	:	RNA polymerase II early elongation complex bound to TFIIE and TFIIF - state a (composite structure)
Authors	:	Zhan, Y.; Grabbe, F.; Oberbeckmann, E.; Dienemann, C.; Cramer, P.
Deposited on	:	2024-02-22
Resolution	:	3.40 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:

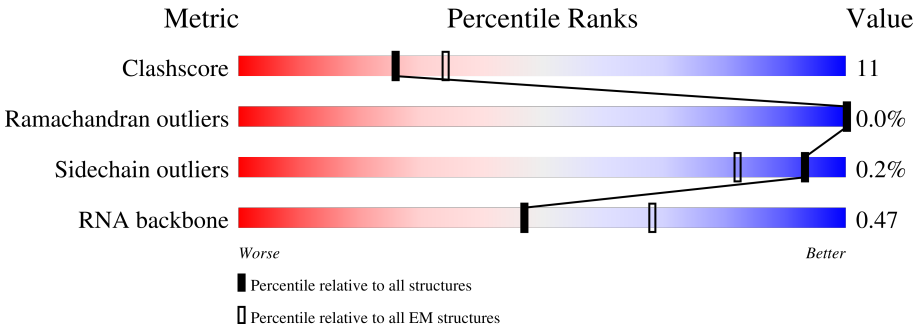
EMDB validation analysis	:	0.0.1.dev92
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	FAILED
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance

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

The reported resolution of this entry is 3.40 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	1984	60% 11% 29%
2	B	1300	73% 14% 13%
3	C	275	79% 15% 7%
4	D	184	35% 35% 30%
5	E	210	87% 12%
6	F	127	53% 12% 35%
7	G	172	48% 52% .

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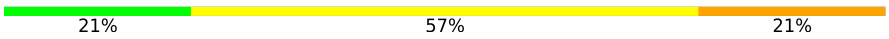
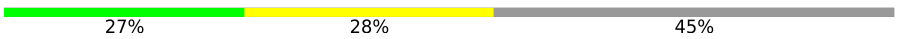
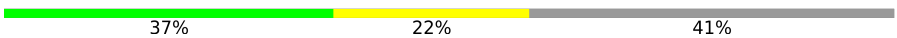
Mol	Chain	Length	Quality of chain
8	H	150	
9	I	125	
10	J	67	
11	K	117	
12	L	58	
13	N	139	
14	P	14	
15	Q	517	
16	R	249	
17	T	139	
18	W	439	
19	X	291	

2 Entry composition

There are 21 unique types of molecules in this entry. The entry contains 38148 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 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1417	11234	7070	2010	2082	72	0	0

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	1130	9030	5714	1587	1665	64	0	0

- Molecule 3 is a protein called DNA-directed RNA polymerase II subunit RPB3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	257	2059	1294	351	408	6	0	0

- Molecule 4 is a protein called RNA polymerase II subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	128	1050	656	178	212	4	0	0

- Molecule 5 is a protein called DNA-directed RNA polymerase II subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	209	1720	1089	300	323	8	0	0

- Molecule 6 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	F	82	657	418	113	121	5	0	0

- Molecule 7 is a protein called DNA-directed RNA polymerase II subunit RPB7.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	171	Total	C	N	O	S	0	0
			1351	875	219	249	8		

- Molecule 8 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	148	Total	C	N	O	S	0	0
			1186	750	194	237	5		

- Molecule 9 is a protein called DNA-directed RNA polymerase II subunit RPB9.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	114	Total	C	N	O	S	0	0
			927	571	166	179	11		

- Molecule 10 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC5.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	67	Total	C	N	O	S	0	0
			533	345	90	92	6		

- Molecule 11 is a protein called DNA-directed RNA polymerase II subunit RPB11-a.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	114	Total	C	N	O	S	0	0
			916	591	151	172	2		

- Molecule 12 is a protein called RNA polymerase II subunit K.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	44	Total	C	N	O	S	0	0
			372	231	72	63	6		

- Molecule 13 is a DNA chain called Non-template DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	N	35	Total	C	N	O	P	0	0
			712	339	132	207	34		

- Molecule 14 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	P	14	Total	C	N	O	P	0	0
			296	133	56	93	14		

- Molecule 15 is a protein called General transcription factor IIF subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	Q	138	Total	C	N	O	S	0	0
			1138	719	208	208	3		

- Molecule 16 is a protein called General transcription factor IIF subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	R	137	Total	C	N	O	S	0	0
			1070	666	193	209	2		

- Molecule 17 is a DNA chain called Template DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	T	46	Total	C	N	O	P	0	0
			949	449	175	279	46		

- Molecule 18 is a protein called General transcription factor IIE subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	W	187	Total	C	N	O	S	0	0
			1535	964	275	285	11		

- Molecule 19 is a protein called Transcription initiation factor IIE subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	X	171	Total	C	N	O	S	0	0
			1403	895	243	261	4		

- Molecule 20 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
20	A	2	Total	Zn	0
			2	2	
20	B	1	Total	Zn	0
			1	1	

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Mol	Chain	Residues	Atoms	AltConf
20	C	1	Total Zn 1 1	0
20	I	2	Total Zn 2 2	0
20	J	1	Total Zn 1 1	0
20	L	1	Total Zn 1 1	0
20	W	1	Total Zn 1 1	0

- Molecule 21 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

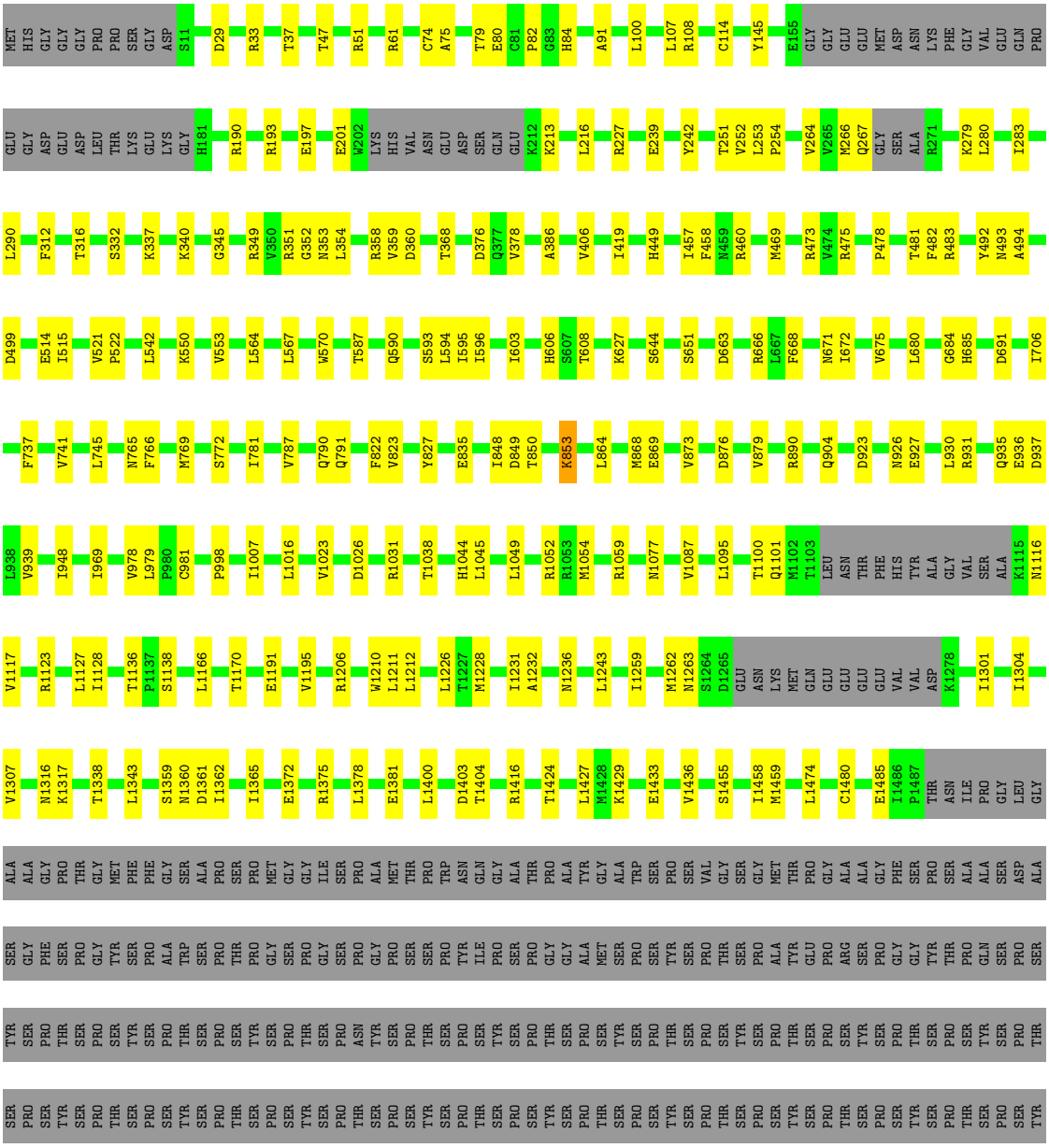
Mol	Chain	Residues	Atoms	AltConf
21	P	1	Total Mg 1 1	0

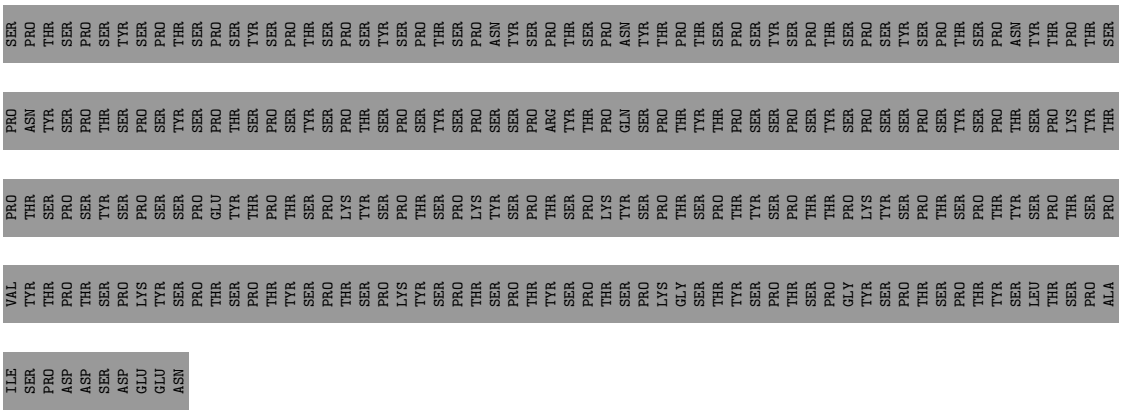
3 Residue-property plots (i)

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. 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. 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 subunit

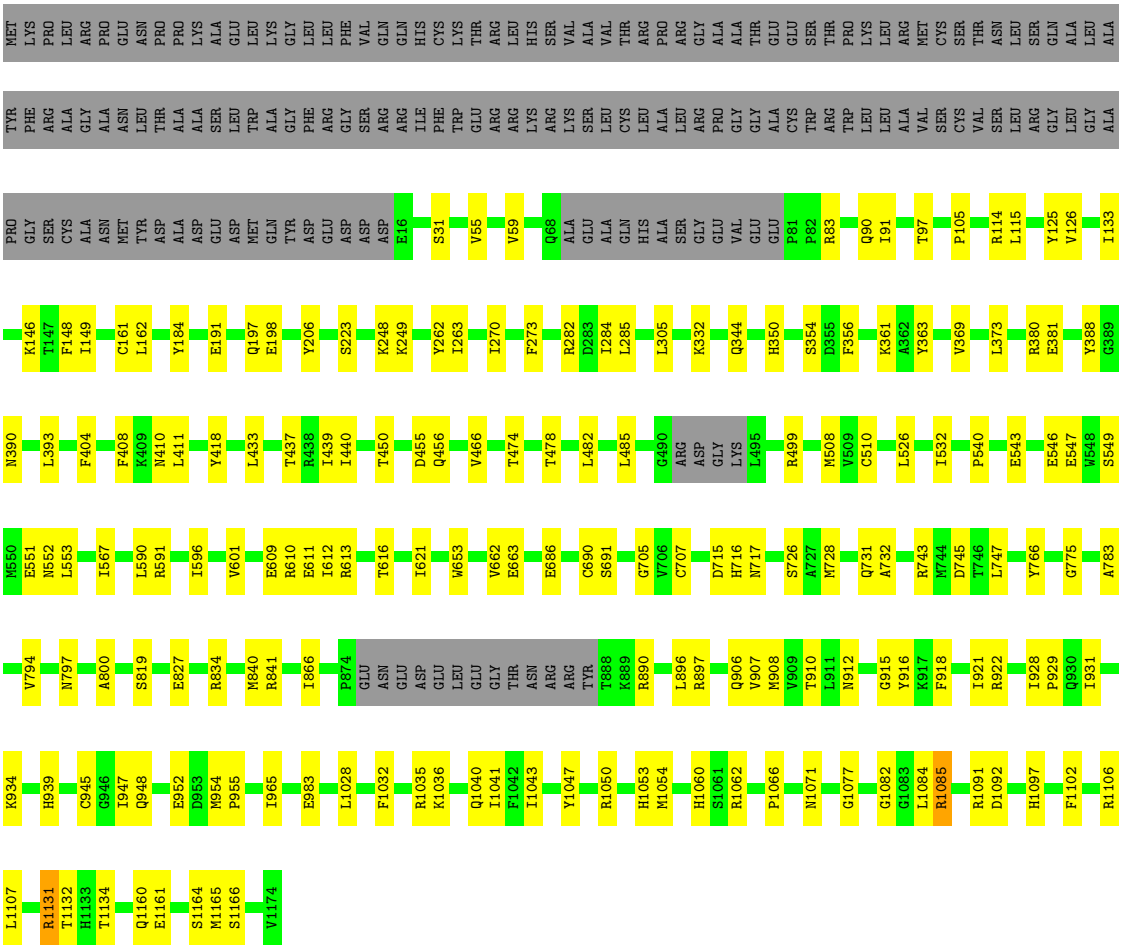
Chain A:





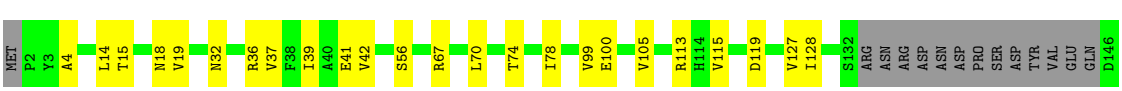
• Molecule 2: DNA-directed RNA polymerase subunit beta

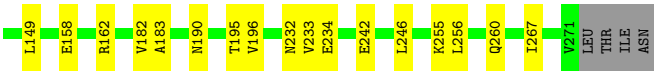
Chain B:



• Molecule 3: DNA-directed RNA polymerase II subunit RPB3

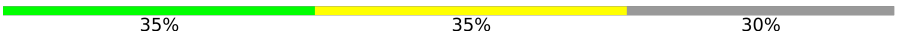
Chain C:

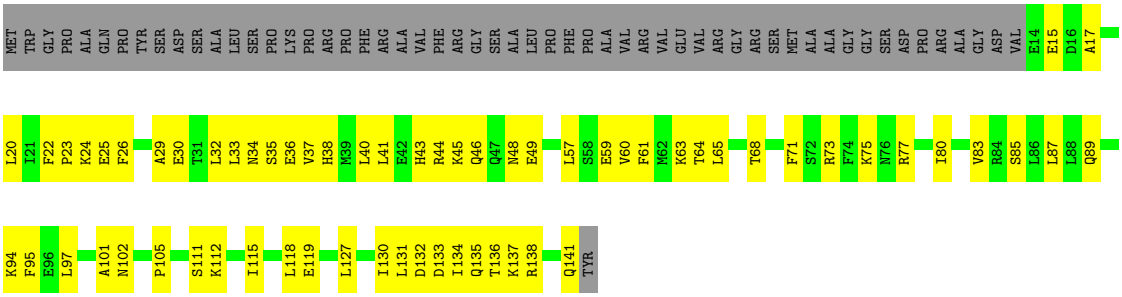




• Molecule 4: RNA polymerase II subunit D

Chain D:



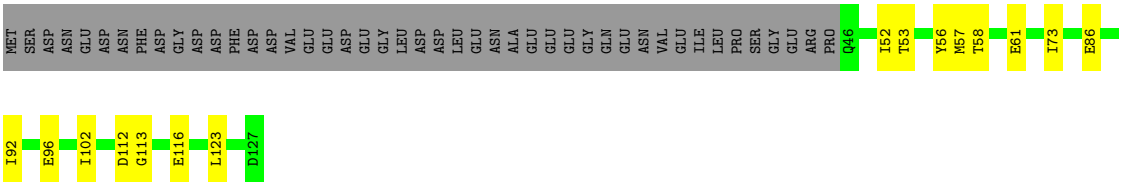


• Molecule 5: DNA-directed RNA polymerase II subunit E

Chain E:

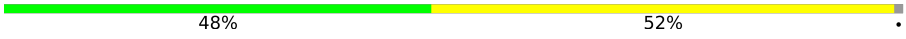
• Molecule 6: DNA-directed RNA polymerases I, II, and III subunit RPABC2

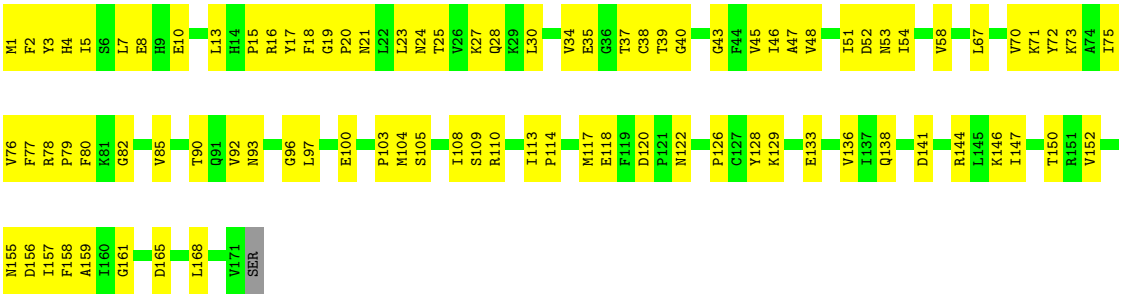
Chain F:



• Molecule 7: DNA-directed RNA polymerase II subunit RPB7

Chain G:



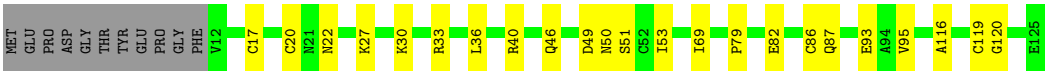


• Molecule 8: DNA-directed RNA polymerases I, II, and III subunit RPABC3

Chain H:

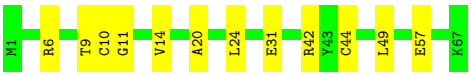
• Molecule 9: DNA-directed RNA polymerase II subunit RPB9

Chain I:



• Molecule 10: DNA-directed RNA polymerases I, II, and III subunit RPABC5

Chain J:

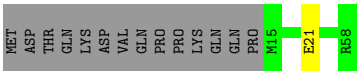


• Molecule 11: DNA-directed RNA polymerase II subunit RPB11-a

Chain K:

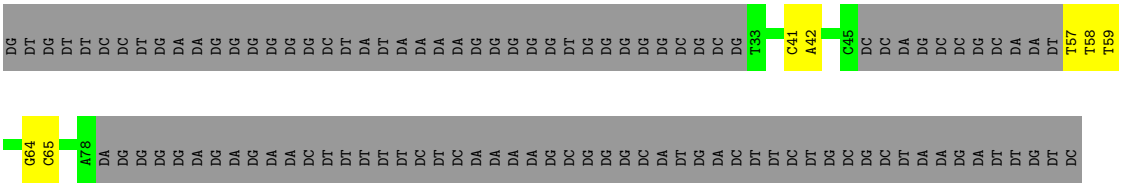
• Molecule 12: RNA polymerase II subunit K

Chain L:



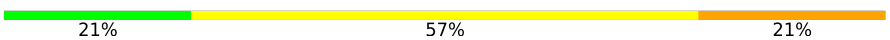
• Molecule 13: Non-template DNA

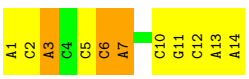
Chain N:



• Molecule 14: RNA

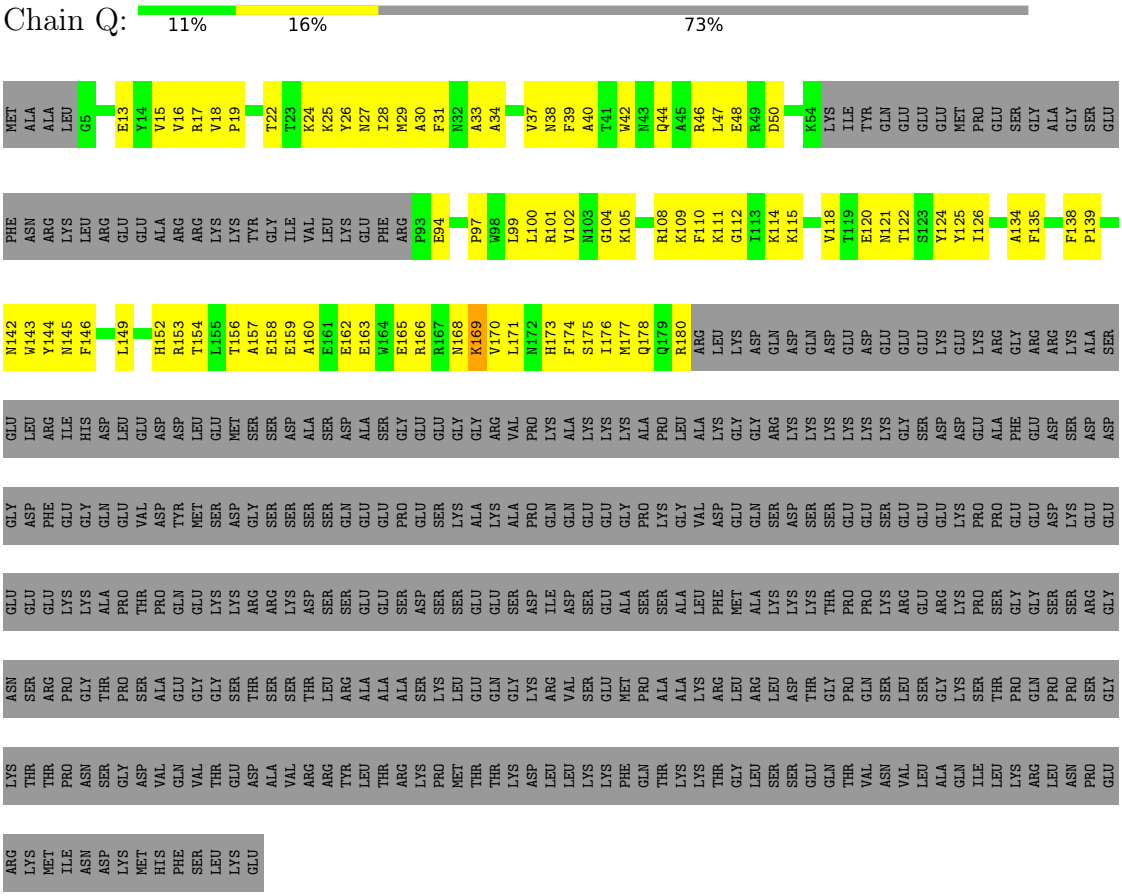
Chain P:





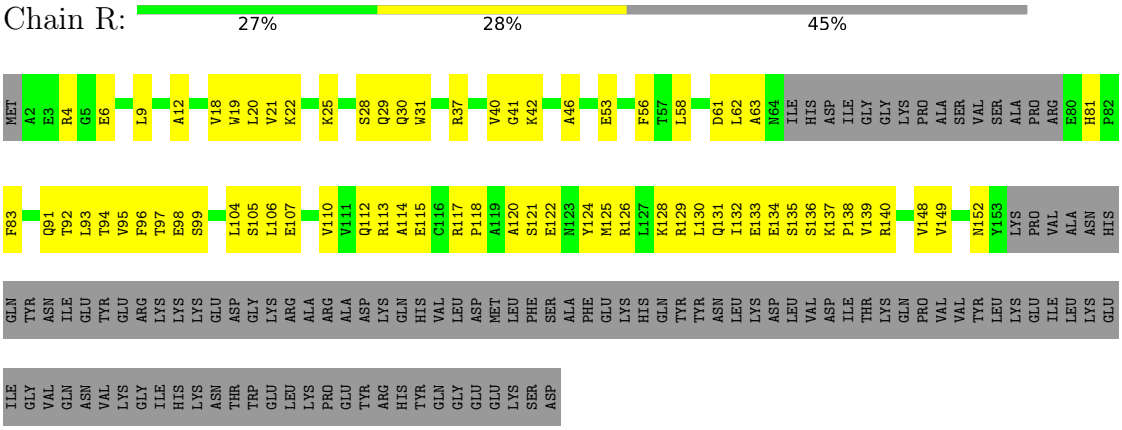
• Molecule 15: General transcription factor IIF subunit 1

Chain Q:



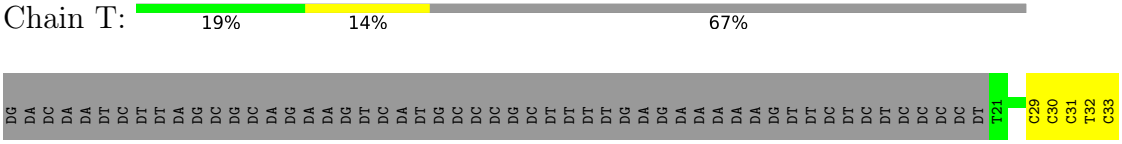
• Molecule 16: General transcription factor IIF subunit 2

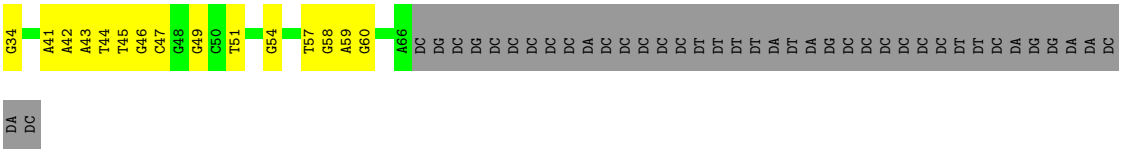
Chain R:



• Molecule 17: Template DNA

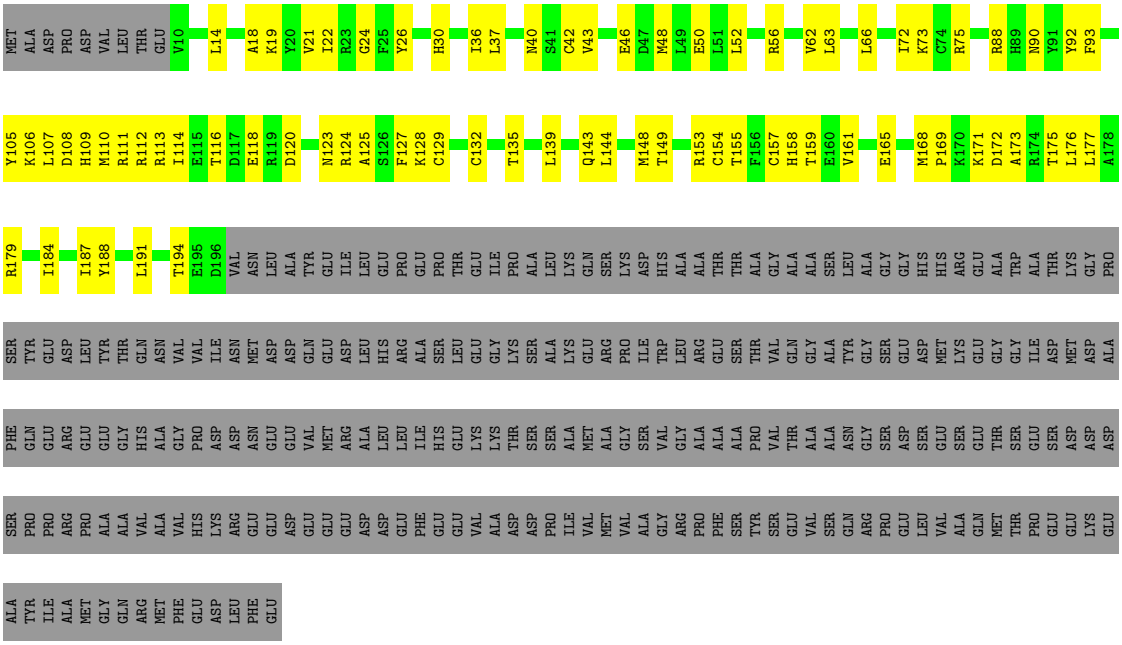
Chain T:





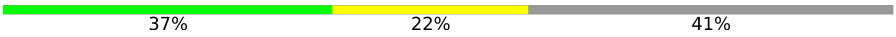
• Molecule 18: General transcription factor IIE subunit 1

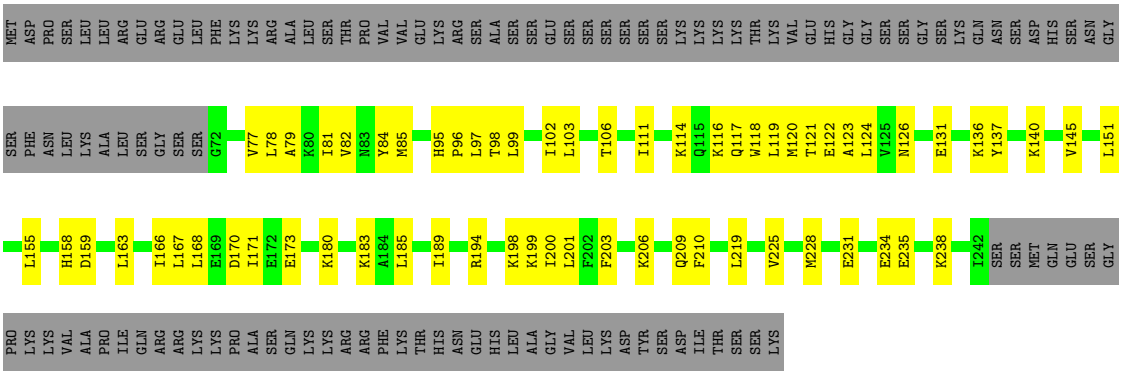
Chain W:



• Molecule 19: Transcription initiation factor IIE subunit beta

Chain X:





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	50000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	40	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	1700	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	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, MG

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	0.29	0/11439	0.58	0/15440
2	B	0.29	0/9210	0.58	0/12431
3	C	0.28	0/2102	0.54	0/2857
4	D	0.24	0/1064	0.43	0/1428
5	E	0.29	0/1751	0.59	0/2366
6	F	0.31	0/667	0.65	0/901
7	G	0.26	0/1382	0.47	0/1874
8	H	0.32	0/1207	0.59	0/1628
9	I	0.29	0/948	0.60	1/1284 (0.1%)
10	J	0.30	0/542	0.59	0/730
11	K	0.29	0/935	0.53	0/1266
12	L	0.30	0/377	0.71	0/500
13	N	0.96	0/797	1.30	0/1225
14	P	0.76	0/330	1.12	0/511
15	Q	0.25	0/1167	0.50	0/1576
16	R	0.25	0/1083	0.51	0/1459
17	T	0.89	0/1064	1.27	0/1642
18	W	0.24	0/1560	0.48	0/2097
19	X	0.23	0/1427	0.42	0/1916
All	All	0.35	0/39052	0.63	1/53131 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
2	B	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	I	20	CYS	CA-CB-SG	5.28	123.51	114.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	108	ARG	Sidechain
1	A	351	ARG	Sidechain
2	B	1085	ARG	Sidechain

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	11234	0	11366	159	0
2	B	9030	0	9073	150	0
3	C	2059	0	2007	28	0
4	D	1050	0	1033	85	0
5	E	1720	0	1737	26	0
6	F	657	0	684	10	0
7	G	1351	0	1358	96	0
8	H	1186	0	1147	15	0
9	I	927	0	859	17	0
10	J	533	0	553	9	0
11	K	916	0	939	15	0
12	L	372	0	378	1	0
13	N	712	0	395	7	0
14	P	296	0	154	20	0
15	Q	1138	0	1103	128	0
16	R	1070	0	1086	103	0
17	T	949	0	518	27	0
18	W	1535	0	1540	63	0
19	X	1403	0	1428	55	0
20	A	2	0	0	0	0
20	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
20	C	1	0	0	0	0
20	I	2	0	0	0	0
20	J	1	0	0	0	0
20	L	1	0	0	0	0
20	W	1	0	0	0	0
21	P	1	0	0	0	0
All	All	38148	0	37358	853	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (853) 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)
2:B:747:LEU:CD2	17:T:49:DG:OP1	1.77	1.32
2:B:747:LEU:HD23	17:T:49:DG:OP1	1.35	1.15
1:A:264:VAL:HG21	14:P:5:C:H4'	1.38	1.05
2:B:747:LEU:HD21	17:T:49:DG:OP1	1.58	1.03
7:G:113:ILE:HG23	7:G:117:MET:HE2	1.40	1.03
7:G:82:GLY:HA2	7:G:146:LYS:HE2	1.42	1.00
7:G:129:LYS:HG2	7:G:136:VAL:HG13	1.44	0.99
8:H:96:VAL:HG12	8:H:116:VAL:HG22	1.45	0.98
5:E:29:THR:CG2	5:E:32:GLU:HG3	1.94	0.98
18:W:184:ILE:HD12	18:W:187:ILE:HD13	1.42	0.97
15:Q:24:LYS:HD3	16:R:99:SER:HB3	1.49	0.94
4:D:60:VAL:HG13	7:G:103:PRO:HB3	1.49	0.93
7:G:52:ASP:HB3	7:G:71:LYS:HG3	1.50	0.93
19:X:99:LEU:HD22	19:X:120:MET:HE3	1.50	0.92
5:E:29:THR:CG2	5:E:32:GLU:CG	2.53	0.85
2:B:841:ARG:HH11	14:P:3:A:C1'	1.91	0.84
1:A:264:VAL:HG21	14:P:5:C:C4'	2.11	0.81
2:B:707:CYS:SG	2:B:726:SER:OG	2.39	0.80
18:W:107:LEU:HD21	18:W:184:ILE:HD11	1.63	0.80
3:C:15:THR:OG1	3:C:18:ASN:OD1	2.00	0.80
1:A:406:VAL:HG21	1:A:419:ILE:HD11	1.63	0.79
15:Q:18:VAL:HG12	16:R:40:VAL:HG11	1.64	0.79
18:W:144:LEU:HD22	18:W:155:THR:H	1.48	0.79
15:Q:110:PHE:HB3	15:Q:146:PHE:HB3	1.65	0.79
1:A:358:ARG:NH2	17:T:46:DG:OP1	2.15	0.78
2:B:783:ALA:HB2	2:B:1041:ILE:HG23	1.65	0.78
5:E:29:THR:HG22	5:E:32:GLU:CD	2.03	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:Q:30:ALA:HB1	16:R:91:GLN:HE21	1.49	0.76
1:A:79:THR:OG1	1:A:80:GLU:OE1	2.04	0.76
1:A:29:ASP:OD2	1:A:33:ARG:NH1	2.19	0.75
14:P:10:C:H2'	14:P:11:G:C8	2.22	0.75
15:Q:15:VAL:O	15:Q:135:PHE:N	2.21	0.73
1:A:1458:ILE:HD11	2:B:1107:LEU:HD12	1.70	0.73
18:W:168:MET:HG2	18:W:169:PRO:HD3	1.69	0.73
1:A:1375:ARG:NE	1:A:1403:ASP:OD1	2.22	0.73
18:W:22:ILE:HD11	18:W:30:HIS:HB3	1.69	0.73
4:D:71:PHE:O	4:D:73:ARG:NH2	2.22	0.72
2:B:952:GLU:O	3:C:36:ARG:NH2	2.23	0.72
3:C:41:GLU:OE2	3:C:255:LYS:NZ	2.18	0.72
5:E:51:GLY:O	5:E:54:ARG:NH1	2.23	0.71
2:B:1040:GLN:NE2	3:C:196:VAL:O	2.23	0.71
2:B:1032:PHE:O	3:C:32:ASN:ND2	2.24	0.71
1:A:998:PRO:O	1:A:1059:ARG:NH2	2.24	0.70
7:G:147:ILE:HG23	7:G:159:ALA:HB1	1.73	0.70
2:B:591:ARG:NH1	2:B:663:GLU:OE2	2.24	0.70
14:P:10:C:H2'	14:P:11:G:H8	1.56	0.70
2:B:1091:ARG:NH2	2:B:1092:ASP:OD1	2.25	0.70
15:Q:166:ARG:HE	15:Q:170:VAL:HG23	1.55	0.70
2:B:146:LYS:O	16:R:152:ASN:ND2	2.23	0.70
3:C:234:GLU:OE1	10:J:42:ARG:NH1	2.25	0.70
17:T:41:DA:H2'	17:T:42:DA:C8	2.27	0.70
2:B:381:GLU:OE1	2:B:381:GLU:N	2.25	0.70
3:C:190:ASN:ND2	3:C:195:THR:O	2.25	0.70
16:R:136:SER:O	16:R:139:VAL:HG12	1.92	0.70
7:G:54:ILE:HD13	7:G:70:VAL:HG13	1.74	0.70
1:A:1206:ARG:NH2	1:A:1263:ASN:O	2.25	0.69
1:A:264:VAL:HG11	14:P:5:C:O2'	1.91	0.69
6:F:53:THR:OG1	6:F:116:GLU:OE1	2.08	0.69
4:D:41:LEU:HD12	4:D:68:THR:HG21	1.74	0.69
1:A:931:ARG:NH2	1:A:936:GLU:OE1	2.25	0.69
1:A:197:GLU:N	1:A:197:GLU:OE1	2.25	0.69
4:D:30:GLU:OE2	7:G:4:HIS:ND1	2.27	0.68
2:B:841:ARG:NH1	14:P:3:A:H1'	2.07	0.68
7:G:40:GLY:HA2	7:G:157:ILE:HD11	1.76	0.68
15:Q:29:MET:HG3	16:R:96:PHE:CD1	2.27	0.68
5:E:29:THR:HG21	5:E:32:GLU:HG3	1.75	0.68
6:F:58:THR:OG1	6:F:61:GLU:OE1	2.08	0.68
5:E:29:THR:HG23	5:E:32:GLU:HG3	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:ARG:NH2	13:N:65:DC:OP1	2.23	0.67
2:B:1085:ARG:NE	17:T:45:DT:OP1	2.26	0.67
4:D:34:ASN:OD1	4:D:102:ASN:ND2	2.28	0.67
16:R:21:VAL:HG22	16:R:114:ALA:HB3	1.76	0.67
19:X:98:THR:HG22	19:X:136:LYS:HG2	1.77	0.67
9:I:40:ARG:NH2	15:Q:174:PHE:HB3	2.10	0.67
2:B:611:GLU:OE1	2:B:613:ARG:NH2	2.28	0.66
5:E:3:ASP:OD1	5:E:49:SER:OG	2.11	0.66
7:G:39:THR:O	7:G:43:GLY:N	2.27	0.66
19:X:159:ASP:HB2	19:X:203:PHE:HB3	1.76	0.66
4:D:32:LEU:HD12	4:D:36:GLU:HG2	1.77	0.66
7:G:10:GLU:N	7:G:10:GLU:OE1	2.26	0.66
1:A:948:ILE:HD12	1:A:1007:ILE:HG21	1.78	0.66
2:B:31:SER:OG	2:B:766:TYR:OH	2.00	0.66
19:X:235:GLU:HA	19:X:238:LYS:HZ3	1.61	0.65
4:D:64:THR:HG21	7:G:46:ILE:HG12	1.78	0.65
8:H:26:SER:HB2	8:H:45:ILE:HD11	1.78	0.65
1:A:457:ILE:HD11	1:A:515:ILE:HD12	1.77	0.65
4:D:131:LEU:O	4:D:135:GLN:HG2	1.97	0.65
15:Q:100:LEU:HD22	15:Q:110:PHE:HD2	1.62	0.65
4:D:57:LEU:HD13	4:D:61:PHE:CE2	2.31	0.65
2:B:543:GLU:OE1	15:Q:121:ASN:ND2	2.29	0.65
1:A:349:ARG:NH2	1:A:353:ASN:OD1	2.29	0.65
2:B:474:THR:OG1	2:B:732:ALA:O	2.07	0.65
4:D:43:HIS:O	4:D:46:GLN:HG3	1.97	0.65
1:A:876:ASP:O	1:A:890:ARG:NH2	2.29	0.65
15:Q:166:ARG:HH11	15:Q:171:LEU:H	1.45	0.65
18:W:173:ALA:O	18:W:177:LEU:HG	1.96	0.64
15:Q:157:ALA:HA	15:Q:160:ALA:HB3	1.78	0.64
15:Q:29:MET:HG3	16:R:96:PHE:HD1	1.63	0.64
14:P:6:C:H2'	14:P:7:A:C8	2.32	0.64
2:B:206:TYR:OH	2:B:390:ASN:O	2.16	0.64
7:G:165:ASP:HB3	7:G:168:LEU:HD11	1.78	0.64
18:W:36:ILE:HG21	18:W:48:MET:HG2	1.80	0.64
4:D:32:LEU:HD12	4:D:36:GLU:CG	2.27	0.64
8:H:143:LEU:HD23	8:H:144:LEU:N	2.12	0.64
4:D:135:GLN:OE1	4:D:138:ARG:NH2	2.31	0.64
4:D:37:VAL:HG21	7:G:2:PHE:HD2	1.63	0.63
15:Q:24:LYS:HD2	16:R:97:THR:HB	1.78	0.63
15:Q:42:TRP:CH2	15:Q:102:VAL:HG22	2.32	0.63
16:R:128:LYS:O	16:R:132:ILE:HG12	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:34:ASN:O	4:D:68:THR:OG1	2.15	0.63
5:E:29:THR:HG23	5:E:32:GLU:H	1.63	0.63
5:E:67:ASP:OD2	5:E:69:THR:OG1	2.08	0.63
15:Q:160:ALA:HA	15:Q:163:GLU:OE2	1.97	0.63
17:T:30:DC:H2'	17:T:31:DC:C6	2.33	0.63
18:W:36:ILE:HD11	18:W:52:LEU:HD21	1.79	0.63
15:Q:99:LEU:HD21	15:Q:101:ARG:HE	1.63	0.63
19:X:122:GLU:OE2	19:X:126:ASN:ND2	2.32	0.63
15:Q:24:LYS:HD3	16:R:99:SER:CB	2.26	0.63
19:X:119:LEU:HA	19:X:123:ALA:HB3	1.81	0.63
10:J:57:GLU:OE1	10:J:57:GLU:N	2.31	0.63
18:W:113:ARG:HG2	19:X:219:LEU:HD21	1.81	0.63
1:A:595:ILE:HD11	1:A:675:VAL:HG11	1.81	0.63
5:E:133:GLN:OE1	5:E:133:GLN:N	2.31	0.63
6:F:52:ILE:O	6:F:53:THR:OG1	2.17	0.63
2:B:841:ARG:HH11	14:P:3:A:H1'	1.62	0.63
2:B:841:ARG:NH1	14:P:3:A:C1'	2.62	0.63
15:Q:15:VAL:HB	15:Q:134:ALA:CB	2.29	0.63
7:G:52:ASP:HB3	7:G:71:LYS:CG	2.27	0.62
1:A:663:ASP:OD1	1:A:666:ARG:NH2	2.31	0.62
7:G:21:ASN:OD1	7:G:24:ASN:HB2	1.98	0.62
4:D:41:LEU:HB2	4:D:65:LEU:HD13	1.82	0.62
1:A:352:GLY:HA2	2:B:1085:ARG:HH12	1.62	0.62
2:B:223:SER:OG	2:B:350:HIS:ND1	2.28	0.62
19:X:81:ILE:HD13	19:X:106:THR:HG21	1.81	0.62
2:B:105:PRO:HB3	16:R:148:VAL:HG11	1.82	0.62
4:D:64:THR:HG21	7:G:46:ILE:HG21	1.82	0.62
15:Q:42:TRP:HH2	15:Q:100:LEU:HD21	1.65	0.62
1:A:644:SER:O	1:A:651:SER:OG	2.15	0.61
14:P:12:C:H2'	14:P:13:A:H8	1.65	0.61
15:Q:118:VAL:HG23	15:Q:120:GLU:OE1	2.00	0.61
4:D:111:SER:OG	4:D:127:LEU:HD21	2.00	0.61
15:Q:44:GLN:HE22	15:Q:46:ARG:HG3	1.65	0.61
1:A:47:THR:OG1	1:A:51:ARG:O	2.08	0.61
19:X:145:VAL:HG11	19:X:151:LEU:HB2	1.82	0.61
16:R:122:GLU:HA	16:R:125:MET:CE	2.31	0.61
2:B:59:VAL:HG23	2:B:408:PHE:CZ	2.35	0.61
7:G:114:PRO:HD2	7:G:117:MET:CE	2.30	0.61
7:G:53:ASN:OD1	7:G:54:ILE:N	2.33	0.61
15:Q:30:ALA:HB3	15:Q:145:ASN:HD22	1.66	0.61
1:A:668:PHE:CZ	1:A:672:ILE:HD11	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:45:LYS:HE3	4:D:61:PHE:CZ	2.36	0.61
15:Q:114:LYS:HZ3	15:Q:115:LYS:HB3	1.65	0.61
1:A:1343:LEU:HD12	1:A:1359:SER:OG	2.01	0.61
18:W:171:LYS:O	18:W:175:THR:HG23	2.00	0.60
2:B:929:PRO:O	2:B:948:GLN:NE2	2.34	0.60
15:Q:44:GLN:HE21	16:R:6:GLU:HG2	1.65	0.60
18:W:42:CYS:SG	18:W:75:ARG:NH1	2.74	0.60
15:Q:121:ASN:HA	16:R:25:LYS:HD3	1.83	0.60
19:X:99:LEU:CD2	19:X:120:MET:HE3	2.30	0.60
4:D:44:ARG:NH2	7:G:47:ALA:HB1	2.16	0.60
5:E:29:THR:CG2	5:E:32:GLU:CD	2.70	0.60
1:A:499:ASP:OD1	14:P:14:A:H4'	2.02	0.60
1:A:1016:LEU:HD22	1:A:1045:LEU:HD21	1.83	0.60
7:G:27:LYS:HD3	7:G:51:ILE:HD13	1.84	0.60
8:H:58:LEU:HD11	8:H:143:LEU:HD21	1.82	0.60
4:D:60:VAL:CG1	7:G:103:PRO:HB3	2.29	0.60
9:I:93:GLU:OE1	9:I:93:GLU:N	2.35	0.60
4:D:41:LEU:HD13	4:D:65:LEU:HA	1.84	0.59
1:A:378:VAL:HG21	1:A:458:PHE:CZ	2.36	0.59
7:G:129:LYS:HD2	7:G:133:GLU:OE2	2.03	0.59
4:D:45:LYS:HE2	4:D:45:LYS:HA	1.84	0.59
12:L:21:GLU:OE1	12:L:21:GLU:N	2.35	0.59
15:Q:94:GLU:N	15:Q:94:GLU:OE1	2.34	0.59
1:A:671:ASN:O	1:A:675:VAL:HG12	2.01	0.59
4:D:37:VAL:HG11	7:G:2:PHE:CE2	2.37	0.59
4:D:132:ASP:O	4:D:136:THR:HG23	2.02	0.59
9:I:116:ALA:HB3	9:I:119:CYS:HB3	1.84	0.59
1:A:1166:LEU:O	1:A:1170:THR:HG22	2.02	0.59
1:A:691:ASP:OD2	1:A:765:ASN:ND2	2.35	0.59
19:X:81:ILE:HG22	19:X:85:MET:CE	2.33	0.59
11:K:11:LEU:O	11:K:37:LYS:NZ	2.32	0.58
18:W:165:GLU:HA	18:W:168:MET:HE1	1.84	0.58
7:G:85:VAL:HA	7:G:144:ARG:HH11	1.67	0.58
2:B:743:ARG:O	2:B:922:ARG:NH2	2.37	0.58
1:A:266:MET:HE1	14:P:5:C:N3	2.18	0.58
7:G:138:GLN:HG2	7:G:141:ASP:OD2	2.03	0.58
15:Q:31:PHE:HD2	15:Q:37:VAL:HG11	1.68	0.58
1:A:590:GLN:O	1:A:593:SER:OG	2.16	0.58
1:A:853:LYS:NZ	1:A:1100:THR:O	2.35	0.58
5:E:38:GLU:OE1	5:E:38:GLU:N	2.35	0.58
11:K:45:ILE:HG21	11:K:94:LEU:HD21	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:Q:109:LYS:O	15:Q:149:LEU:N	2.37	0.58
1:A:904:GLN:NE2	1:A:981:CYS:O	2.36	0.58
4:D:33:LEU:HD13	4:D:101:ALA:HB1	1.85	0.58
7:G:45:VAL:HA	7:G:76:VAL:HG12	1.86	0.58
9:I:27:LYS:O	9:I:36:LEU:N	2.36	0.58
15:Q:33:ALA:HB2	16:R:91:GLN:HA	1.85	0.58
1:A:1307:VAL:HG22	1:A:1338:THR:HG22	1.84	0.58
1:A:564:LEU:HD13	1:A:567:LEU:HD12	1.85	0.58
11:K:100:LEU:HD21	11:K:104:ARG:HH11	1.69	0.58
15:Q:39:PHE:HA	15:Q:42:TRP:HD1	1.67	0.58
1:A:868:MET:SD	1:A:1404:THR:HG21	2.44	0.58
1:A:948:ILE:HG23	1:A:1007:ILE:CG2	2.34	0.57
10:J:31:GLU:OE1	10:J:31:GLU:N	2.36	0.57
15:Q:111:LYS:HB2	15:Q:149:LEU:HD11	1.85	0.57
2:B:161:CYS:SG	2:B:162:LEU:N	2.77	0.57
1:A:266:MET:SD	14:P:5:C:C2	2.97	0.57
15:Q:29:MET:HB2	16:R:94:THR:HG23	1.85	0.57
1:A:923:ASP:OD1	1:A:1052:ARG:NH2	2.37	0.57
1:A:1077:ASN:ND2	6:F:56:TYR:OH	2.36	0.57
2:B:282:ARG:NH2	15:Q:34:ALA:HB1	2.18	0.57
9:I:79:PRO:O	9:I:95:VAL:HG23	2.03	0.57
15:Q:163:GLU:OE1	15:Q:163:GLU:N	2.33	0.57
19:X:234:GLU:OE1	19:X:238:LYS:NZ	2.37	0.57
18:W:168:MET:HG2	18:W:169:PRO:CD	2.34	0.57
3:C:183:ALA:HB3	3:C:232:ASN:HB3	1.87	0.57
15:Q:13:GLU:OE2	16:R:42:LYS:HD3	2.05	0.57
15:Q:27:ASN:O	16:R:95:VAL:HG13	2.05	0.57
2:B:197:GLN:NE2	2:B:466:VAL:HG22	2.19	0.57
4:D:87:LEU:HB3	4:D:97:LEU:HD11	1.87	0.57
11:K:77:THR:OG1	11:K:81:TYR:O	2.17	0.57
1:A:368:THR:HG23	1:A:481:THR:HG21	1.87	0.57
7:G:120:ASP:OD1	7:G:122:ASN:ND2	2.37	0.57
7:G:38:CYS:HB3	7:G:155:ASN:O	2.05	0.57
13:N:58:DT:H1'	13:N:59:DT:H5"	1.86	0.57
15:Q:24:LYS:CD	16:R:97:THR:HB	2.34	0.57
4:D:133:ASP:O	4:D:137:LYS:HG2	2.05	0.57
7:G:118:GLU:OE1	7:G:118:GLU:N	2.37	0.57
19:X:180:LYS:O	19:X:183:LYS:HG2	2.05	0.57
1:A:542:LEU:HD21	1:A:772:SER:O	2.05	0.56
2:B:866:ILE:HD12	2:B:921:ILE:HD11	1.86	0.56
15:Q:30:ALA:HB1	16:R:91:GLN:NE2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:148:PHE:CD2	2:B:437:THR:HG21	2.40	0.56
15:Q:39:PHE:HA	15:Q:42:TRP:CD1	2.40	0.56
15:Q:163:GLU:O	15:Q:166:ARG:HG2	2.05	0.56
4:D:37:VAL:O	4:D:41:LEU:HG	2.05	0.56
7:G:150:THR:HG22	7:G:159:ALA:HB2	1.85	0.56
7:G:34:VAL:O	7:G:37:THR:HG22	2.05	0.56
19:X:131:GLU:HB2	19:X:140:LYS:CG	2.35	0.56
1:A:406:VAL:CG2	1:A:419:ILE:HD11	2.32	0.56
4:D:41:LEU:CD1	4:D:68:THR:HG21	2.36	0.56
15:Q:166:ARG:HA	15:Q:169:LYS:HE3	1.88	0.56
1:A:1138:SER:OG	1:A:1360:ASN:ND2	2.38	0.56
2:B:747:LEU:HD21	17:T:49:DG:P	2.46	0.56
2:B:1077:GLY:HA2	17:T:47:DC:OP2	2.06	0.56
15:Q:166:ARG:NE	15:Q:170:VAL:HG23	2.19	0.56
19:X:114:LYS:O	19:X:117:GLN:HG3	2.06	0.56
19:X:131:GLU:HB2	19:X:140:LYS:HG3	1.87	0.56
1:A:890:ARG:HE	1:A:1023:VAL:HG22	1.71	0.55
4:D:57:LEU:HD13	4:D:61:PHE:HE2	1.67	0.55
6:F:57:MET:HB2	6:F:123:LEU:HD13	1.88	0.55
9:I:30:LYS:O	9:I:33:ARG:NH1	2.38	0.55
1:A:1127:LEU:HD12	1:A:1378:LEU:CD1	2.37	0.55
6:F:73:ILE:HG21	6:F:92:ILE:HD11	1.87	0.55
16:R:96:PHE:HA	16:R:106:LEU:HA	1.89	0.55
3:C:56:SER:OG	3:C:158:GLU:N	2.39	0.55
11:K:45:ILE:CG2	11:K:94:LEU:HD21	2.37	0.55
18:W:176:LEU:HD11	18:W:179:ARG:HH11	1.71	0.55
1:A:460:ARG:NH1	1:A:493:ASN:O	2.39	0.55
1:A:1458:ILE:CD1	2:B:1107:LEU:HD12	2.35	0.55
5:E:29:THR:HG22	5:E:32:GLU:CG	2.32	0.55
16:R:130:LEU:HA	16:R:133:GLU:OE1	2.07	0.55
2:B:115:LEU:HD22	2:B:908:MET:HE1	1.87	0.55
15:Q:114:LYS:HD3	15:Q:142:ASN:HB3	1.89	0.55
2:B:1161:GLU:O	2:B:1164:SER:OG	2.17	0.55
8:H:8:ASP:OD1	8:H:9:ILE:N	2.39	0.55
15:Q:121:ASN:HA	16:R:25:LYS:CG	2.37	0.55
5:E:66:ASP:OD1	5:E:67:ASP:N	2.40	0.55
2:B:526:LEU:HD11	2:B:532:ILE:HD11	1.88	0.55
7:G:8:GLU:OE2	7:G:71:LYS:HE3	2.07	0.55
10:J:14:VAL:HG13	10:J:49:LEU:HD21	1.89	0.55
18:W:37:LEU:CD1	18:W:72:ILE:HG21	2.35	0.55
1:A:745:LEU:HD12	1:A:822:PHE:CE1	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:248:LYS:NZ	15:Q:171:LEU:HD23	2.21	0.55
15:Q:16:VAL:HG21	16:R:56:PHE:CE1	2.41	0.55
16:R:120:ALA:HB3	16:R:124:TYR:HB2	1.87	0.55
18:W:63:LEU:HD22	18:W:92:TYR:CE2	2.42	0.55
5:E:74:VAL:HG23	5:E:103:LEU:HB2	1.89	0.55
7:G:10:GLU:HB2	7:G:67:LEU:HD11	1.89	0.55
18:W:14:LEU:HD23	18:W:194:THR:HG22	1.89	0.55
8:H:31:GLU:OE1	8:H:31:GLU:N	2.40	0.54
15:Q:15:VAL:HG13	16:R:42:LYS:NZ	2.22	0.54
15:Q:44:GLN:NE2	16:R:6:GLU:HG2	2.22	0.54
3:C:242:GLU:OE1	3:C:242:GLU:N	2.40	0.54
16:R:93:LEU:HB2	16:R:110:VAL:HB	1.89	0.54
1:A:91:ALA:HB3	1:A:290:LEU:HD22	1.88	0.54
1:A:1232:ALA:O	1:A:1236:ASN:ND2	2.40	0.54
2:B:91:ILE:HD13	2:B:126:VAL:HG12	1.88	0.54
16:R:21:VAL:HA	16:R:114:ALA:O	2.07	0.54
18:W:129:CYS:HB2	18:W:132:CYS:SG	2.48	0.54
15:Q:46:ARG:HH22	15:Q:104:GLY:H	1.53	0.54
19:X:77:VAL:HG21	19:X:111:ILE:HD11	1.90	0.54
4:D:131:LEU:O	4:D:134:ILE:HG13	2.08	0.54
1:A:1316:ASN:OD1	1:A:1317:LYS:N	2.41	0.54
2:B:149:ILE:HA	2:B:437:THR:HG22	1.89	0.54
4:D:41:LEU:CB	4:D:65:LEU:HD13	2.38	0.54
15:Q:24:LYS:CG	16:R:97:THR:HB	2.38	0.54
16:R:97:THR:N	16:R:105:SER:O	2.40	0.54
4:D:59:GLU:N	4:D:59:GLU:OE1	2.38	0.54
1:A:266:MET:CE	14:P:5:C:N3	2.71	0.54
4:D:45:LYS:O	4:D:49:GLU:HG2	2.08	0.54
15:Q:15:VAL:HG22	16:R:42:LYS:HZ3	1.73	0.54
1:A:869:GLU:OE1	1:A:1455:SER:OG	2.23	0.54
4:D:85:SER:O	4:D:89:GLN:HG2	2.08	0.54
15:Q:30:ALA:HB3	15:Q:145:ASN:ND2	2.23	0.54
16:R:4:ARG:HA	16:R:4:ARG:NE	2.22	0.54
18:W:43:VAL:HB	18:W:48:MET:HE2	1.90	0.54
19:X:170:ASP:O	19:X:173:GLU:HG3	2.08	0.54
1:A:376:ASP:OD2	1:A:473:ARG:NE	2.40	0.53
4:D:33:LEU:HD22	4:D:101:ALA:HB3	1.90	0.53
4:D:119:GLU:N	4:D:119:GLU:OE1	2.41	0.53
19:X:168:LEU:N	19:X:199:LYS:O	2.37	0.53
19:X:185:LEU:HD12	19:X:189:ILE:HD11	1.90	0.53
19:X:231:GLU:O	19:X:234:GLU:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1365:ILE:HD11	5:E:144:LEU:HD13	1.89	0.53
2:B:939:HIS:NE2	2:B:983:GLU:OE1	2.40	0.53
4:D:71:PHE:HB3	4:D:73:ARG:HH22	1.73	0.53
1:A:606:HIS:HD2	1:A:608:THR:HG22	1.73	0.53
16:R:56:PHE:HB3	16:R:81:HIS:HB2	1.90	0.53
18:W:46:GLU:O	18:W:50:GLU:HG2	2.09	0.53
19:X:185:LEU:CD1	19:X:189:ILE:HD11	2.37	0.53
1:A:848:ILE:HD11	2:B:499:ARG:HD3	1.89	0.53
2:B:931:ILE:HG22	2:B:945:CYS:SG	2.49	0.53
1:A:91:ALA:CB	1:A:290:LEU:HD22	2.39	0.53
9:I:17:CYS:SG	9:I:22:ASN:N	2.80	0.53
1:A:145:TYR:OH	13:N:64:DG:OP1	2.23	0.53
1:A:1427:LEU:HD12	1:A:1459:MET:SD	2.49	0.53
2:B:1050:ARG:NH1	2:B:1054:MET:SD	2.82	0.53
6:F:112:ASP:OD1	6:F:113:GLY:N	2.41	0.53
16:R:135:SER:O	16:R:138:PRO:HD2	2.09	0.53
1:A:332:SER:OG	17:T:54:DG:OP2	2.27	0.53
7:G:114:PRO:HD2	7:G:117:MET:HE1	1.90	0.53
1:A:359:VAL:HG21	2:B:1106:ARG:HD3	1.90	0.53
4:D:35:SER:HB3	4:D:80:ILE:CD1	2.39	0.53
5:E:120:ASP:OD1	5:E:121:MET:N	2.41	0.53
18:W:129:CYS:SG	18:W:161:VAL:HG22	2.48	0.53
2:B:834:ARG:NH2	2:B:840:MET:SD	2.82	0.52
7:G:158:PHE:CE2	18:W:139:LEU:HD23	2.45	0.52
15:Q:18:VAL:CG1	16:R:40:VAL:HG11	2.36	0.52
16:R:58:LEU:HD23	16:R:63:ALA:HB2	1.91	0.52
4:D:48:ASN:HD22	4:D:57:LEU:HD12	1.74	0.52
4:D:87:LEU:HB3	4:D:97:LEU:CD1	2.40	0.52
16:R:31:TRP:CZ3	16:R:40:VAL:HB	2.44	0.52
16:R:130:LEU:HA	16:R:133:GLU:CD	2.30	0.52
7:G:25:THR:O	7:G:28:GLN:HG3	2.08	0.52
6:F:96:GLU:OE1	6:F:102:ILE:HD13	2.10	0.52
4:D:118:LEU:HD21	4:D:127:LEU:HD22	1.91	0.52
2:B:508:MET:HE1	2:B:621:ILE:HG23	1.91	0.52
4:D:77:ARG:HH11	4:D:80:ILE:CD1	2.22	0.52
9:I:50:ASN:OD1	9:I:51:SER:N	2.43	0.52
15:Q:111:LYS:O	15:Q:146:PHE:HA	2.10	0.52
7:G:40:GLY:HA2	7:G:157:ILE:CD1	2.40	0.52
17:T:32:DT:H2'	17:T:33:DC:C6	2.44	0.52
18:W:154:CYS:O	18:W:158:HIS:N	2.40	0.52
2:B:1060:HIS:NE2	2:B:1082:GLY:O	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:55:VAL:HG23	2:B:404:PHE:CE2	2.46	0.51
4:D:45:LYS:HE3	4:D:61:PHE:HZ	1.75	0.51
13:N:58:DT:H1'	13:N:59:DT:C5'	2.41	0.51
15:Q:44:GLN:HE22	15:Q:46:ARG:CG	2.23	0.51
1:A:596:ILE:HD13	1:A:668:PHE:CE1	2.46	0.51
7:G:82:GLY:CA	7:G:146:LYS:HE2	2.28	0.51
15:Q:46:ARG:HB3	15:Q:48:GLU:OE2	2.09	0.51
15:Q:124:TYR:HB3	16:R:20:LEU:HD11	1.92	0.51
1:A:685:HIS:NE2	1:A:769:MET:SD	2.83	0.51
16:R:21:VAL:HG11	16:R:83:PHE:CE2	2.45	0.51
2:B:146:LYS:HD3	16:R:149:VAL:HG21	1.91	0.51
7:G:4:HIS:HA	7:G:75:ILE:HD13	1.93	0.51
15:Q:26:TYR:CD1	16:R:97:THR:HG22	2.45	0.51
16:R:131:GLN:HA	16:R:134:GLU:OE1	2.10	0.51
17:T:59:DA:H2"	17:T:60:DG:C8	2.45	0.51
18:W:120:ASP:HA	18:W:123:ASN:HB2	1.93	0.51
4:D:71:PHE:HB3	4:D:73:ARG:NH2	2.25	0.51
9:I:119:CYS:SG	9:I:120:GLY:N	2.81	0.51
15:Q:163:GLU:HB2	15:Q:166:ARG:HD3	1.93	0.51
1:A:926:ASN:OD1	1:A:927:GLU:N	2.43	0.51
15:Q:18:VAL:HG13	15:Q:18:VAL:O	2.11	0.51
15:Q:29:MET:HG2	15:Q:146:PHE:CE1	2.45	0.51
15:Q:114:LYS:HD3	15:Q:144:TYR:HE1	1.76	0.51
18:W:125:ALA:O	18:W:128:LYS:NZ	2.41	0.51
1:A:492:TYR:O	1:A:494:ALA:N	2.44	0.51
9:I:36:LEU:HD12	9:I:46:GLN:O	2.11	0.51
18:W:113:ARG:HA	18:W:116:THR:OG1	2.11	0.51
1:A:201:GLU:OE2	1:A:213:LYS:NZ	2.44	0.51
18:W:172:ASP:O	18:W:176:LEU:HD23	2.10	0.51
2:B:596:ILE:HG22	16:R:129:ARG:HA	1.93	0.50
5:E:29:THR:HG22	5:E:32:GLU:OE1	2.11	0.50
15:Q:39:PHE:O	15:Q:42:TRP:HB2	2.11	0.50
19:X:103:LEU:HA	19:X:106:THR:HG1	1.76	0.50
1:A:475:ARG:NH1	11:K:68:GLU:OE2	2.43	0.50
4:D:38:HIS:HA	4:D:41:LEU:HD12	1.93	0.50
18:W:120:ASP:O	18:W:124:ARG:HG3	2.12	0.50
4:D:127:LEU:O	4:D:131:LEU:HG	2.12	0.50
17:T:57:DT:H2"	17:T:58:DG:N7	2.26	0.50
18:W:113:ARG:HD2	19:X:219:LEU:HD11	1.94	0.50
1:A:280:LEU:HA	1:A:283:ILE:HG22	1.93	0.50
1:A:360:ASP:OD1	2:B:1062:ARG:NE	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:728:MET:O	2:B:731:GLN:N	2.45	0.50
4:D:35:SER:HB3	4:D:80:ILE:HD11	1.94	0.50
5:E:48:PRO:O	5:E:49:SER:OG	2.28	0.50
1:A:1424:THR:O	1:A:1429:LYS:NZ	2.44	0.50
2:B:747:LEU:CD2	17:T:49:DG:P	2.95	0.50
4:D:37:VAL:HG11	7:G:2:PHE:HE2	1.76	0.50
5:E:71:GLN:HB2	5:E:99:ILE:HG22	1.92	0.50
9:I:82:GLU:OE1	9:I:82:GLU:N	2.39	0.50
16:R:37:ARG:NH2	16:R:61:ASP:OD2	2.45	0.50
1:A:873:VAL:HG22	1:A:879:VAL:HG22	1.92	0.50
7:G:15:PRO:HA	7:G:18:PHE:CE1	2.47	0.50
14:P:5:C:H2'	14:P:6:C:C6	2.47	0.50
15:Q:162:GLU:O	15:Q:165:GLU:HG3	2.11	0.50
16:R:124:TYR:OH	16:R:128:LYS:HD3	2.11	0.50
19:X:99:LEU:HB2	19:X:120:MET:HE3	1.93	0.50
1:A:564:LEU:HD11	1:A:594:LEU:HD13	1.93	0.50
1:A:564:LEU:HD12	1:A:570:TRP:CD2	2.46	0.50
2:B:273:PHE:HB3	2:B:284:ILE:HD12	1.94	0.50
4:D:41:LEU:HD22	4:D:61:PHE:CE1	2.47	0.50
7:G:114:PRO:HD2	7:G:117:MET:HE2	1.93	0.50
15:Q:153:ARG:HH11	15:Q:177:MET:CE	2.25	0.50
1:A:550:LYS:O	1:A:553:VAL:HG12	2.12	0.50
1:A:1474:LEU:HD23	7:G:58:VAL:HA	1.94	0.50
2:B:455:ASP:OD1	2:B:456:GLN:N	2.45	0.50
15:Q:112:GLY:HA2	15:Q:145:ASN:O	2.12	0.50
2:B:797:ASN:HB3	2:B:800:ALA:HB3	1.93	0.50
4:D:29:ALA:HB2	7:G:5:ILE:HG22	1.94	0.50
2:B:1028:LEU:HD11	2:B:1043:ILE:HD11	1.93	0.49
2:B:1132:THR:OG1	2:B:1134:THR:HG22	2.12	0.49
3:C:67:ARG:NH1	3:C:149:LEU:O	2.45	0.49
7:G:1:MET:HA	7:G:77:PHE:HD1	1.76	0.49
18:W:107:LEU:HD22	18:W:188:TYR:CE1	2.47	0.49
1:A:1095:LEU:O	1:A:1095:LEU:HD23	2.12	0.49
2:B:285:LEU:HD21	2:B:305:LEU:HD21	1.95	0.49
1:A:904:GLN:OE1	1:A:1044:HIS:NE2	2.46	0.49
1:A:469:MET:O	2:B:1097:HIS:NE2	2.44	0.49
16:R:28:SER:HA	16:R:31:TRP:CD1	2.47	0.49
19:X:103:LEU:HA	19:X:106:THR:OG1	2.13	0.49
3:C:260:GLN:HG2	11:K:91:ILE:HG21	1.95	0.49
7:G:165:ASP:HB3	7:G:168:LEU:CD1	2.43	0.49
16:R:97:THR:HG21	16:R:107:GLU:OE2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:R:97:THR:HG23	16:R:107:GLU:HG2	1.95	0.49
18:W:107:LEU:HD22	18:W:188:TYR:HE1	1.77	0.49
1:A:267:GLN:O	2:B:890:ARG:NH1	2.46	0.49
1:A:1038:THR:HG22	1:A:1038:THR:O	2.12	0.49
18:W:88:ARG:NH2	18:W:90:ASN:OD1	2.46	0.49
19:X:78:LEU:HD12	19:X:79:ALA:N	2.28	0.49
4:D:15:GLU:OE1	4:D:23:PRO:HD3	2.13	0.49
7:G:146:LYS:O	7:G:161:GLY:HA2	2.13	0.49
9:I:87:GLN:OE1	9:I:87:GLN:N	2.46	0.49
17:T:31:DC:H2'	17:T:32:DT:C6	2.47	0.49
1:A:481:THR:HG22	1:A:482:PHE:O	2.12	0.49
16:R:12:ALA:HB2	16:R:106:LEU:HD23	1.95	0.49
16:R:121:SER:O	16:R:125:MET:HE1	2.13	0.49
16:R:133:GLU:HA	16:R:136:SER:OG	2.13	0.49
1:A:1087:VAL:HG12	1:A:1400:LEU:HD22	1.95	0.48
15:Q:29:MET:HB2	16:R:94:THR:CG2	2.43	0.48
15:Q:156:THR:HG23	15:Q:158:GLU:HG2	1.95	0.48
7:G:30:LEU:O	7:G:34:VAL:HG22	2.13	0.48
8:H:91:VAL:HG22	8:H:144:LEU:HD13	1.96	0.48
11:K:65:HIS:HE1	11:K:67:LEU:HD12	1.78	0.48
1:A:930:LEU:HD23	1:A:939:VAL:HG12	1.95	0.48
2:B:547:GLU:OE2	16:R:118:PRO:HD3	2.12	0.48
15:Q:46:ARG:HG3	15:Q:46:ARG:HH21	1.77	0.48
2:B:546:GLU:OE2	15:Q:122:THR:HG21	2.13	0.48
2:B:715:ASP:N	2:B:715:ASP:OD1	2.47	0.48
2:B:834:ARG:NH2	2:B:841:ARG:O	2.44	0.48
7:G:96:GLY:HA3	7:G:108:ILE:O	2.14	0.48
7:G:105:SER:O	7:G:158:PHE:HB2	2.12	0.48
19:X:166:ILE:HG21	19:X:171:ILE:HD11	1.95	0.48
2:B:478:THR:O	2:B:482:LEU:HD23	2.14	0.48
2:B:1084:LEU:CD1	17:T:46:DG:H5"	2.44	0.48
4:D:36:GLU:O	4:D:40:LEU:HG	2.13	0.48
15:Q:42:TRP:HB3	16:R:9:LEU:HD13	1.96	0.48
19:X:155:LEU:HD22	19:X:203:PHE:N	2.28	0.48
8:H:28:LEU:N	8:H:41:LEU:O	2.43	0.48
1:A:514:GLU:OE2	2:B:1102:PHE:N	2.44	0.48
2:B:775:GLY:N	2:B:1047:TYR:OH	2.47	0.48
2:B:910:THR:N	2:B:918:PHE:O	2.40	0.48
16:R:46:ALA:HB3	16:R:53:GLU:HB3	1.95	0.48
16:R:126:ARG:O	16:R:130:LEU:HG	2.14	0.48
18:W:24:GLY:HA2	19:X:210:PHE:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:W:40:ASN:ND2	18:W:42:CYS:O	2.46	0.48
2:B:686:GLU:N	2:B:686:GLU:OE1	2.47	0.48
2:B:907:VAL:HG13	2:B:921:ILE:HG22	1.95	0.48
7:G:51:ILE:HG13	7:G:51:ILE:O	2.14	0.48
1:A:849:ASP:OD1	1:A:850:THR:N	2.46	0.48
2:B:508:MET:CE	2:B:621:ILE:HG23	2.44	0.48
2:B:149:ILE:CA	2:B:437:THR:HG22	2.44	0.47
2:B:610:ARG:CZ	9:I:69:ILE:HD12	2.44	0.47
7:G:35:GLU:OE2	7:G:47:ALA:HA	2.14	0.47
8:H:96:VAL:HG12	8:H:116:VAL:CG2	2.29	0.47
15:Q:114:LYS:NZ	15:Q:115:LYS:HB3	2.27	0.47
1:A:312:PHE:O	1:A:316:THR:HG22	2.14	0.47
15:Q:47:LEU:HD12	15:Q:99:LEU:O	2.15	0.47
15:Q:50:ASP:OD2	15:Q:97:PRO:HD2	2.14	0.47
16:R:18:VAL:HG23	16:R:110:VAL:HG22	1.94	0.47
17:T:59:DA:H2"	17:T:60:DG:H8	1.79	0.47
19:X:158:HIS:NE2	19:X:166:ILE:HD11	2.29	0.47
1:A:706:ILE:HD11	1:A:787:VAL:HG21	1.97	0.47
4:D:25:GLU:HG3	4:D:26:PHE:CD2	2.49	0.47
15:Q:175:SER:HA	15:Q:178:GLN:OE1	2.14	0.47
17:T:45:DT:H2'	17:T:46:DG:C8	2.49	0.47
18:W:110:MET:O	18:W:114:ILE:HG12	2.14	0.47
19:X:228:MET:SD	19:X:228:MET:N	2.87	0.47
1:A:1485:GLU:OE2	7:G:20:PRO:HD2	2.14	0.47
2:B:410:ASN:ND2	2:B:439:ILE:HD11	2.30	0.47
3:C:99:VAL:HG21	3:C:127:VAL:HG21	1.96	0.47
18:W:139:LEU:O	18:W:143:GLN:HG2	2.14	0.47
2:B:248:LYS:HZ1	15:Q:171:LEU:HD23	1.79	0.47
2:B:1040:GLN:OE1	2:B:1040:GLN:N	2.47	0.47
15:Q:31:PHE:HB2	16:R:92:THR:OG1	2.14	0.47
15:Q:173:HIS:O	15:Q:176:ILE:HG22	2.14	0.47
16:R:31:TRP:CE3	16:R:40:VAL:HB	2.49	0.47
17:T:45:DT:H2'	17:T:46:DG:H8	1.79	0.47
2:B:552:ASN:OD1	2:B:553:LEU:N	2.47	0.47
2:B:745:ASP:OD1	2:B:745:ASP:N	2.46	0.47
4:D:34:ASN:H	4:D:102:ASN:HD21	1.61	0.47
7:G:48:VAL:HG13	7:G:72:TYR:HD2	1.79	0.47
9:I:86:CYS:SG	9:I:87:GLN:N	2.88	0.47
1:A:935:GLN:O	1:A:939:VAL:HG13	2.14	0.47
2:B:819:SER:N	2:B:827:GLU:OE2	2.48	0.47
7:G:79:PRO:HB2	7:G:150:THR:HG21	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:P:6:C:H2'	14:P:7:A:H8	1.77	0.47
19:X:99:LEU:HB2	19:X:120:MET:CE	2.45	0.47
1:A:82:PRO:O	2:B:1160:GLN:NE2	2.45	0.47
1:A:567:LEU:HD21	1:A:595:ILE:HG12	1.97	0.47
15:Q:120:GLU:O	16:R:25:LYS:HD3	2.15	0.47
15:Q:121:ASN:HA	16:R:25:LYS:CD	2.45	0.47
2:B:114:ARG:NE	2:B:191:GLU:OE2	2.47	0.47
16:R:19:TRP:CD1	16:R:112:GLN:HB3	2.50	0.47
1:A:1049:LEU:HD13	1:A:1054:MET:CE	2.45	0.47
2:B:197:GLN:HE21	2:B:466:VAL:HG22	1.80	0.47
2:B:601:VAL:HG22	2:B:616:THR:HG23	1.96	0.47
18:W:108:ASP:O	18:W:112:ARG:HG2	2.15	0.47
2:B:433:LEU:HD23	2:B:433:LEU:O	2.15	0.46
2:B:910:THR:O	2:B:918:PHE:N	2.45	0.46
2:B:1134:THR:HG23	2:B:1134:THR:O	2.15	0.46
3:C:246:LEU:HD12	11:K:106:ARG:HH11	1.80	0.46
4:D:130:ILE:O	4:D:134:ILE:HG12	2.14	0.46
15:Q:152:HIS:CE1	15:Q:154:THR:HG22	2.50	0.46
17:T:43:DA:H2'	17:T:44:DT:C6	2.50	0.46
18:W:37:LEU:HD13	18:W:72:ILE:HG21	1.97	0.46
2:B:1085:ARG:CD	17:T:45:DT:OP1	2.63	0.46
16:R:97:THR:O	16:R:104:LEU:HA	2.14	0.46
4:D:80:ILE:HA	4:D:83:VAL:HG12	1.97	0.46
4:D:105:PRO:HG2	4:D:131:LEU:HD22	1.98	0.46
4:D:135:GLN:O	4:D:138:ARG:HB2	2.15	0.46
7:G:15:PRO:HA	7:G:18:PHE:CZ	2.51	0.46
11:K:24:ASP:OD1	11:K:26:LYS:N	2.44	0.46
19:X:82:VAL:O	19:X:85:MET:HG2	2.15	0.46
1:A:564:LEU:HD11	1:A:594:LEU:CD1	2.46	0.46
4:D:23:PRO:HG2	4:D:26:PHE:HD2	1.81	0.46
18:W:157:CYS:SG	18:W:159:THR:HG22	2.55	0.46
2:B:125:TYR:CE2	16:R:149:VAL:HB	2.51	0.46
2:B:931:ILE:HG21	2:B:947:ILE:HA	1.96	0.46
7:G:1:MET:SD	7:G:2:PHE:N	2.89	0.46
11:K:63:VAL:HG12	11:K:71:ILE:HG22	1.98	0.46
16:R:41:GLY:O	16:R:42:LYS:HE2	2.16	0.46
1:A:251:THR:HG23	1:A:252:VAL:HG23	1.97	0.46
4:D:41:LEU:O	4:D:45:LYS:HG2	2.15	0.46
7:G:1:MET:HA	7:G:77:PHE:CD1	2.50	0.46
18:W:107:LEU:HD21	18:W:184:ILE:CD1	2.40	0.46
1:A:1116:ASN:OD1	1:A:1117:VAL:N	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:78:ILE:HD11	3:C:127:VAL:CG2	2.46	0.46
4:D:63:LYS:NZ	7:G:100:GLU:OE2	2.48	0.46
4:D:112:LYS:HB3	4:D:119:GLU:OE2	2.16	0.46
7:G:43:GLY:HA2	7:G:157:ILE:HD12	1.98	0.46
7:G:80:PHE:O	7:G:147:ILE:HB	2.16	0.46
15:Q:114:LYS:CD	15:Q:142:ASN:HB3	2.45	0.46
1:A:1228:MET:HA	1:A:1231:ILE:HD12	1.98	0.46
2:B:270:ILE:HD11	2:B:305:LEU:HD22	1.97	0.46
2:B:510:CYS:SG	2:B:705:GLY:N	2.88	0.46
2:B:551:GLU:OE1	16:R:113:ARG:NH1	2.48	0.46
7:G:79:PRO:HG2	7:G:150:THR:HG21	1.98	0.46
10:J:9:THR:OG1	10:J:44:CYS:SG	2.73	0.46
13:N:58:DT:H2"	13:N:59:DT:H5'	1.96	0.46
16:R:28:SER:HA	16:R:31:TRP:HD1	1.80	0.46
19:X:114:LYS:HA	19:X:117:GLN:HG3	1.98	0.46
1:A:239:GLU:OE1	1:A:242:TYR:N	2.49	0.46
1:A:337:LYS:HE2	1:A:337:LYS:HA	1.98	0.46
1:A:1243:LEU:HD21	1:A:1259:ILE:HG23	1.97	0.46
7:G:3:TYR:O	7:G:75:ILE:HA	2.16	0.46
7:G:93:ASN:O	7:G:128:TYR:OH	2.17	0.46
15:Q:19:PRO:HG2	15:Q:138:PHE:HE1	1.81	0.46
2:B:928:ILE:HD12	2:B:928:ILE:H	1.81	0.46
6:F:73:ILE:HG21	6:F:92:ILE:CD1	2.46	0.46
2:B:912:ASN:OD1	2:B:916:TYR:N	2.45	0.45
4:D:17:ALA:O	4:D:20:LEU:HD23	2.16	0.45
4:D:127:LEU:HA	4:D:130:ILE:HG12	1.98	0.45
15:Q:101:ARG:HD3	15:Q:109:LYS:HD3	1.98	0.45
1:A:603:ILE:HD11	1:A:627:LYS:HE3	1.98	0.45
15:Q:27:ASN:ND2	16:R:98:GLU:OE2	2.50	0.45
4:D:17:ALA:HB1	4:D:95:PHE:CZ	2.51	0.45
16:R:124:TYR:CZ	16:R:128:LYS:HD3	2.50	0.45
2:B:344:GLN:O	2:B:361:LYS:NZ	2.45	0.45
3:C:105:VAL:HG11	3:C:115:VAL:HG22	1.97	0.45
4:D:44:ARG:HH22	7:G:48:VAL:N	2.15	0.45
5:E:29:THR:HG23	5:E:32:GLU:CG	2.33	0.45
13:N:41:DC:H2"	13:N:42:DA:C8	2.52	0.45
1:A:567:LEU:HD22	1:A:671:ASN:HD22	1.81	0.45
1:A:1211:LEU:HD23	1:A:1212:LEU:N	2.32	0.45
2:B:146:LYS:HD3	16:R:149:VAL:HG11	1.99	0.45
2:B:380:ARG:NH2	2:B:609:GLU:OE2	2.48	0.45
2:B:450:THR:HB	17:T:51:DT:H4'	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:115:ILE:HG22	4:D:118:LEU:H	1.81	0.45
15:Q:99:LEU:HD21	15:Q:101:ARG:NE	2.30	0.45
18:W:107:LEU:HD12	18:W:191:LEU:HD11	1.98	0.45
19:X:235:GLU:HA	19:X:238:LYS:NZ	2.31	0.45
1:A:340:LYS:HE3	1:A:1436:VAL:HG11	1.98	0.45
4:D:64:THR:HG23	7:G:2:PHE:CZ	2.52	0.45
15:Q:28:ILE:HD12	16:R:94:THR:O	2.17	0.45
15:Q:38:ASN:OD1	15:Q:40:ALA:HB3	2.17	0.45
18:W:43:VAL:HB	18:W:48:MET:CE	2.45	0.45
1:A:279:LYS:NZ	1:A:316:THR:HG23	2.31	0.45
2:B:249:LYS:NZ	15:Q:168:ASN:HD21	2.14	0.45
7:G:13:LEU:HD11	7:G:17:TYR:HB2	1.99	0.45
8:H:30:CYS:O	8:H:39:LEU:N	2.42	0.45
15:Q:100:LEU:HD22	15:Q:110:PHE:CD2	2.46	0.45
15:Q:126:ILE:CG2	16:R:18:VAL:HB	2.46	0.45
1:A:114:CYS:O	1:A:227:ARG:NH1	2.49	0.45
1:A:595:ILE:CD1	1:A:675:VAL:HG11	2.45	0.45
1:A:823:VAL:HG22	1:A:835:GLU:HB2	1.99	0.45
1:A:978:VAL:O	1:A:979:LEU:HD22	2.16	0.45
4:D:23:PRO:CD	7:G:78:ARG:HH12	2.28	0.45
15:Q:166:ARG:CA	15:Q:169:LYS:HE3	2.47	0.45
17:T:33:DC:H2'	17:T:34:DG:C8	2.52	0.45
1:A:457:ILE:HD11	1:A:515:ILE:CD1	2.45	0.45
1:A:1361:ASP:OD1	1:A:1362:ILE:N	2.49	0.45
2:B:596:ILE:CG2	16:R:132:ILE:HG13	2.47	0.45
15:Q:25:LYS:HB2	16:R:98:GLU:HG3	1.99	0.45
15:Q:101:ARG:NH2	15:Q:109:LYS:HE2	2.32	0.45
2:B:356:PHE:CE2	15:Q:114:LYS:HB3	2.52	0.45
4:D:26:PHE:CE1	7:G:5:ILE:HG21	2.51	0.45
15:Q:28:ILE:HG22	15:Q:143:TRP:HA	1.99	0.45
17:T:44:DT:H2'	17:T:45:DT:C6	2.52	0.45
1:A:478:PRO:O	1:A:483:ARG:NH1	2.50	0.44
2:B:363:TYR:CD2	2:B:553:LEU:HD11	2.52	0.44
3:C:100:GLU:OE2	3:C:162:ARG:NH2	2.50	0.44
1:A:1301:ILE:HB	1:A:1304:ILE:HD12	2.00	0.44
2:B:567:ILE:HD13	2:B:612:ILE:HB	2.00	0.44
3:C:182:VAL:HG23	3:C:233:VAL:HG22	1.98	0.44
1:A:100:LEU:HD21	1:A:193:ARG:HD2	1.99	0.44
2:B:248:LYS:NZ	15:Q:170:VAL:O	2.43	0.44
2:B:934:LYS:NZ	14:P:14:A:OP1	2.41	0.44
5:E:189:GLN:O	5:E:209:VAL:HG22	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:W:62:VAL:O	18:W:66:LEU:HD23	2.18	0.44
18:W:148:MET:SD	18:W:149:THR:N	2.90	0.44
1:A:107:LEU:HD21	1:A:216:LEU:HD21	2.00	0.44
2:B:90:GLN:CB	16:R:140:ARG:HD3	2.48	0.44
4:D:32:LEU:HD23	7:G:4:HIS:HB2	1.99	0.44
2:B:369:VAL:O	2:B:373:LEU:HD23	2.18	0.44
2:B:540:PRO:HB3	16:R:132:ILE:HD12	1.99	0.44
4:D:131:LEU:HA	4:D:134:ILE:HG12	2.00	0.44
7:G:152:VAL:HA	7:G:156:ASP:O	2.17	0.44
9:I:116:ALA:HB3	9:I:119:CYS:CB	2.46	0.44
18:W:19:LYS:O	18:W:22:ILE:HG22	2.18	0.44
18:W:113:ARG:CG	19:X:219:LEU:HD11	2.48	0.44
1:A:969:ILE:HD12	5:E:201:GLY:O	2.17	0.44
2:B:590:LEU:CD2	16:R:125:MET:HG2	2.48	0.44
5:E:173:ILE:HG23	5:E:209:VAL:HG12	1.99	0.44
2:B:590:LEU:HD21	16:R:125:MET:HG2	2.00	0.44
4:D:77:ARG:HH11	4:D:80:ILE:HD12	1.81	0.44
15:Q:26:TYR:HD2	15:Q:138:PHE:HB3	1.83	0.44
18:W:128:LYS:HE3	18:W:135:THR:HG22	1.98	0.44
2:B:97:THR:HG21	16:R:148:VAL:HB	1.99	0.44
8:H:74:GLU:N	8:H:74:GLU:OE1	2.50	0.44
19:X:171:ILE:HG13	19:X:201:LEU:HD12	1.98	0.44
19:X:194:ARG:HD2	19:X:200:ILE:HD11	2.00	0.44
8:H:26:SER:CB	8:H:45:ILE:HD11	2.47	0.44
10:J:10:CYS:SG	10:J:11:GLY:N	2.91	0.44
15:Q:125:TYR:HD1	15:Q:139:PRO:HA	1.83	0.44
15:Q:156:THR:CG2	15:Q:159:GLU:HG3	2.47	0.44
16:R:30:GLN:O	16:R:62:LEU:HD11	2.18	0.44
18:W:108:ASP:OD1	18:W:109:HIS:N	2.51	0.44
4:D:32:LEU:HD12	4:D:36:GLU:HG3	1.99	0.43
16:R:30:GLN:OE1	16:R:62:LEU:HG	2.17	0.43
1:A:354:LEU:HD13	1:A:1459:MET:HG3	2.01	0.43
2:B:896:LEU:HD23	2:B:897:ARG:O	2.18	0.43
15:Q:31:PHE:CD2	15:Q:37:VAL:HG21	2.53	0.43
1:A:1016:LEU:HD22	1:A:1045:LEU:CD2	2.47	0.43
5:E:29:THR:HG23	5:E:32:GLU:CB	2.48	0.43
18:W:105:TYR:HA	18:W:108:ASP:OD2	2.17	0.43
1:A:864:LEU:HD21	1:A:1128:ILE:HD12	2.00	0.43
1:A:936:GLU:HA	1:A:939:VAL:HG22	2.00	0.43
2:B:717:ASN:OD1	2:B:939:HIS:ND1	2.52	0.43
2:B:794:VAL:HG13	2:B:965:ILE:HG23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:Q:15:VAL:HB	15:Q:134:ALA:HB1	2.00	0.43
15:Q:33:ALA:HB2	16:R:91:GLN:CA	2.46	0.43
15:Q:158:GLU:HG2	15:Q:159:GLU:H	1.84	0.43
18:W:113:ARG:CD	19:X:219:LEU:HD11	2.48	0.43
1:A:37:THR:OG1	1:A:61:ARG:NH1	2.51	0.43
1:A:521:VAL:HG13	1:A:522:PRO:HD3	2.00	0.43
1:A:745:LEU:HD12	1:A:822:PHE:CD1	2.54	0.43
1:A:1101:GLN:N	1:A:1101:GLN:OE1	2.52	0.43
2:B:369:VAL:HG22	2:B:373:LEU:HD23	2.01	0.43
4:D:37:VAL:HG11	7:G:2:PHE:CD2	2.54	0.43
15:Q:15:VAL:HG22	16:R:42:LYS:NZ	2.34	0.43
1:A:1136:THR:O	1:A:1136:THR:OG1	2.32	0.43
3:C:37:VAL:O	3:C:42:VAL:HG23	2.19	0.43
7:G:120:ASP:HB2	7:G:129:LYS:HZ2	1.83	0.43
1:A:386:ALA:O	1:A:449:HIS:ND1	2.50	0.43
1:A:1480:CYS:SG	7:G:19:GLY:HA2	2.58	0.43
2:B:114:ARG:NH1	2:B:184:TYR:OH	2.48	0.43
2:B:262:TYR:O	2:B:263:ILE:HD13	2.19	0.43
3:C:14:LEU:HD12	3:C:19:VAL:HG22	2.00	0.43
3:C:256:LEU:HB3	11:K:95:ILE:HD11	1.99	0.43
6:F:86:GLU:N	6:F:86:GLU:OE1	2.52	0.43
1:A:587:THR:HG21	8:H:119:GLY:O	2.18	0.43
1:A:1123:ARG:NE	1:A:1381:GLU:OE2	2.51	0.43
7:G:109:SER:O	7:G:113:ILE:HG13	2.19	0.43
7:G:110:ARG:NH1	7:G:113:ILE:HG21	2.34	0.43
1:A:353:ASN:OD1	2:B:1071:ASN:ND2	2.52	0.43
1:A:1307:VAL:HG22	1:A:1338:THR:CG2	2.49	0.43
1:A:790:GLN:NE2	1:A:791:GLN:O	2.52	0.42
2:B:90:GLN:HA	16:R:140:ARG:HD3	2.00	0.42
2:B:393:LEU:HD13	2:B:485:LEU:HD22	2.01	0.42
7:G:7:LEU:HD22	7:G:72:TYR:OH	2.20	0.42
16:R:134:GLU:O	16:R:137:LYS:HB2	2.19	0.42
18:W:73:LYS:HB2	18:W:93:PHE:CE2	2.54	0.42
18:W:111:ARG:O	18:W:114:ILE:HB	2.19	0.42
19:X:82:VAL:HA	19:X:85:MET:HG2	2.00	0.42
1:A:827:TYR:O	2:B:716:HIS:ND1	2.52	0.42
19:X:97:LEU:HB2	19:X:102:ILE:HD11	2.02	0.42
19:X:111:ILE:HG23	19:X:116:LYS:HE2	2.01	0.42
19:X:131:GLU:HB2	19:X:140:LYS:HG2	2.01	0.42
1:A:340:LYS:CD	1:A:1436:VAL:HG11	2.49	0.42
15:Q:15:VAL:HG13	16:R:42:LYS:HZ1	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:Q:115:LYS:HB2	15:Q:143:TRP:HB2	2.01	0.42
18:W:184:ILE:O	18:W:184:ILE:HG13	2.20	0.42
19:X:158:HIS:ND1	19:X:163:LEU:HD12	2.34	0.42
1:A:74:CYS:SG	1:A:84:HIS:CE1	3.12	0.42
1:A:668:PHE:CE1	1:A:672:ILE:HD11	2.54	0.42
2:B:952:GLU:HA	3:C:39:ILE:HD11	2.02	0.42
15:Q:99:LEU:CD2	15:Q:101:ARG:HE	2.31	0.42
15:Q:125:TYR:HA	15:Q:138:PHE:O	2.19	0.42
18:W:165:GLU:O	18:W:168:MET:HE3	2.18	0.42
7:G:104:MET:HG2	7:G:158:PHE:HA	2.02	0.42
3:C:70:LEU:O	10:J:6:ARG:NH2	2.49	0.42
7:G:90:THR:HG21	7:G:100:GLU:HB2	2.01	0.42
18:W:106:LYS:NZ	19:X:225:VAL:HG11	2.35	0.42
1:A:606:HIS:CD2	1:A:608:THR:HG22	2.53	0.42
4:D:30:GLU:HG2	7:G:4:HIS:HB3	2.02	0.42
15:Q:124:TYR:OH	16:R:22:LYS:HE2	2.19	0.42
16:R:4:ARG:HA	16:R:4:ARG:CZ	2.50	0.42
2:B:411:LEU:HD22	2:B:440:ILE:HD11	2.02	0.42
3:C:113:ARG:NH1	3:C:119:ASP:OD2	2.51	0.42
7:G:4:HIS:NE2	7:G:73:LYS:HB3	2.34	0.42
15:Q:17:ARG:N	15:Q:135:PHE:O	2.48	0.42
16:R:29:GLN:OE1	16:R:29:GLN:HA	2.19	0.42
16:R:115:GLU:HA	16:R:115:GLU:OE1	2.19	0.42
16:R:122:GLU:HA	16:R:125:MET:HE2	2.01	0.42
2:B:105:PRO:CB	16:R:148:VAL:HG11	2.49	0.42
2:B:1062:ARG:NH2	2:B:1066:PRO:O	2.50	0.42
15:Q:42:TRP:CZ3	15:Q:102:VAL:HG22	2.54	0.42
15:Q:102:VAL:HG21	15:Q:108:ARG:HG2	2.02	0.42
19:X:95:HIS:ND1	19:X:96:PRO:O	2.51	0.42
1:A:680:LEU:O	1:A:684:GLY:N	2.53	0.42
1:A:1210:TRP:CZ3	9:I:53:ILE:HD13	2.55	0.42
1:A:1372:GLU:OE2	5:E:207:ARG:NH1	2.52	0.42
2:B:954:MET:HB3	2:B:955:PRO:HD2	2.02	0.42
4:D:75:LYS:HG2	4:D:141:GLN:NE2	2.34	0.42
15:Q:124:TYR:HD2	15:Q:143:TRP:HZ2	1.68	0.42
18:W:176:LEU:CD1	18:W:179:ARG:HH11	2.33	0.42
19:X:155:LEU:CD1	19:X:189:ILE:HG12	2.50	0.42
2:B:198:GLU:OE1	2:B:388:TYR:OH	2.31	0.41
2:B:906:GLN:OE1	2:B:922:ARG:NH1	2.53	0.41
7:G:97:LEU:HG	7:G:128:TYR:CE2	2.55	0.41
15:Q:153:ARG:HB2	15:Q:180:ARG:CZ	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:Q:169:LYS:HD2	15:Q:169:LYS:O	2.20	0.41
19:X:118:TRP:HA	19:X:121:THR:HG22	2.02	0.41
1:A:253:LEU:HD23	1:A:254:PRO:O	2.20	0.41
1:A:937:ASP:OD1	1:A:937:ASP:N	2.53	0.41
2:B:912:ASN:OD1	2:B:915:GLY:N	2.53	0.41
7:G:23:LEU:HD23	7:G:23:LEU:H	1.85	0.41
8:H:147:LYS:O	8:H:148:LEU:HD22	2.20	0.41
1:A:766:PHE:HB3	1:A:781:ILE:HD12	2.00	0.41
4:D:26:PHE:HE1	7:G:5:ILE:HD13	1.85	0.41
11:K:82:SER:OG	11:K:84:GLN:OE1	2.38	0.41
15:Q:166:ARG:HE	15:Q:170:VAL:HA	1.86	0.41
1:A:1433:GLU:OE1	17:T:41:DA:H4'	2.20	0.41
2:B:485:LEU:HD21	2:B:526:LEU:HD21	2.01	0.41
7:G:7:LEU:HD22	7:G:72:TYR:CE1	2.54	0.41
7:G:34:VAL:HG21	7:G:48:VAL:CG2	2.50	0.41
15:Q:19:PRO:HG2	15:Q:138:PHE:CE1	2.55	0.41
15:Q:28:ILE:HD13	16:R:95:VAL:HG22	2.03	0.41
16:R:18:VAL:CG2	16:R:110:VAL:HA	2.50	0.41
18:W:63:LEU:HD22	18:W:92:TYR:CZ	2.56	0.41
2:B:418:TYR:OH	2:B:433:LEU:HD22	2.21	0.41
2:B:1165:MET:O	2:B:1166:SER:OG	2.26	0.41
4:D:29:ALA:O	4:D:94:LYS:NZ	2.34	0.41
7:G:152:VAL:HB	7:G:157:ILE:HG12	2.02	0.41
15:Q:24:LYS:HD3	16:R:99:SER:CA	2.50	0.41
18:W:36:ILE:HD11	18:W:52:LEU:CD2	2.49	0.41
2:B:1035:ARG:NH2	2:B:1036:LYS:O	2.53	0.41
9:I:49:ASP:OD1	9:I:49:ASP:N	2.54	0.41
1:A:1191:GLU:O	1:A:1195:VAL:HG23	2.20	0.41
1:A:1191:GLU:OE1	1:A:1191:GLU:N	2.53	0.41
2:B:690:CYS:SG	2:B:691:SER:N	2.94	0.41
3:C:78:ILE:HD11	3:C:127:VAL:HG23	2.03	0.41
4:D:75:LYS:N	4:D:75:LYS:HD3	2.34	0.41
1:A:564:LEU:HD12	1:A:570:TRP:CE2	2.56	0.41
1:A:737:PHE:O	1:A:741:VAL:HG23	2.21	0.41
4:D:26:PHE:CE1	7:G:5:ILE:HD13	2.55	0.41
18:W:22:ILE:O	18:W:26:TYR:HB2	2.21	0.41
1:A:603:ILE:HG23	1:A:603:ILE:O	2.21	0.41
2:B:90:GLN:HB2	16:R:140:ARG:HD3	2.03	0.41
2:B:248:LYS:NZ	15:Q:171:LEU:HA	2.35	0.41
2:B:354:SER:O	2:B:356:PHE:N	2.49	0.41
2:B:596:ILE:CG2	16:R:129:ARG:HA	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:7:LEU:HD23	7:G:8:GLU:N	2.36	0.41
8:H:88:PHE:HB2	8:H:144:LEU:HD12	2.03	0.41
15:Q:105:LYS:HD3	15:Q:105:LYS:N	2.35	0.41
16:R:133:GLU:O	16:R:137:LYS:HG3	2.20	0.41
18:W:18:ALA:HA	18:W:21:VAL:HG22	2.02	0.41
18:W:168:MET:N	18:W:169:PRO:HD2	2.36	0.41
19:X:114:LYS:HA	19:X:117:GLN:CG	2.51	0.41
1:A:75:ALA:HB2	2:B:1131:ARG:NE	2.36	0.41
1:A:1026:ASP:OD1	1:A:1031:ARG:NH1	2.54	0.41
2:B:83:ARG:N	2:B:133:ILE:O	2.48	0.41
3:C:267:ILE:HD12	11:K:84:GLN:HB3	2.02	0.41
4:D:22:PHE:HB3	4:D:23:PRO:HD2	2.03	0.41
7:G:16:ARG:HB3	7:G:16:ARG:NH2	2.36	0.41
14:P:1:A:N3	14:P:1:A:H2'	2.35	0.41
15:Q:16:VAL:HA	15:Q:135:PHE:O	2.21	0.41
15:Q:22:THR:HG21	15:Q:26:TYR:CZ	2.56	0.41
15:Q:28:ILE:CD1	16:R:95:VAL:HG22	2.51	0.41
17:T:29:DC:H2"	17:T:30:DC:O4'	2.21	0.41
19:X:81:ILE:HG23	19:X:102:ILE:HG21	2.03	0.41
19:X:124:LEU:HD13	19:X:137:TYR:CD2	2.55	0.41
19:X:167:LEU:HD23	19:X:198:LYS:HD3	2.03	0.41
2:B:549:SER:HB3	16:R:117:ARG:HH21	1.86	0.40
7:G:92:VAL:HB	7:G:126:PRO:HG2	2.03	0.40
18:W:127:PHE:HD1	18:W:161:VAL:HG11	1.85	0.40
19:X:84:TYR:HD2	19:X:102:ILE:HG23	1.87	0.40
10:J:20:ALA:O	10:J:24:LEU:HD23	2.20	0.40
3:C:4:ALA:HB2	11:K:93:ASP:HB3	2.02	0.40
3:C:74:THR:HG22	3:C:128:ILE:O	2.21	0.40
4:D:30:GLU:HG2	4:D:30:GLU:O	2.21	0.40
4:D:41:LEU:HD22	4:D:61:PHE:HE1	1.86	0.40
7:G:85:VAL:HA	7:G:144:ARG:NH1	2.32	0.40
13:N:57:DT:H2"	13:N:58:DT:H5'	2.03	0.40
15:Q:156:THR:HG22	15:Q:159:GLU:HG3	2.03	0.40
18:W:118:GLU:OE2	18:W:177:LEU:HD13	2.21	0.40
19:X:206:LYS:HA	19:X:209:GLN:OE1	2.21	0.40
1:A:948:ILE:HD12	1:A:1007:ILE:HD13	2.03	0.40
1:A:1206:ARG:NH1	1:A:1262:MET:O	2.55	0.40
2:B:356:PHE:CD2	15:Q:114:LYS:HB3	2.56	0.40
2:B:653:TRP:CZ2	2:B:662:VAL:HG11	2.56	0.40
2:B:1053:HIS:NE2	14:P:13:A:H4'	2.36	0.40
7:G:152:VAL:HG23	7:G:157:ILE:HG12	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:165:ASP:O	7:G:168:LEU:HD12	2.22	0.40
1:A:930:LEU:CD2	1:A:939:VAL:HG12	2.52	0.40
1:A:1226:LEU:HD23	1:A:1226:LEU:H	1.86	0.40
1:A:1416:ARG:NH1	17:T:41:DA:O4'	2.55	0.40
7:G:27:LYS:HD3	7:G:51:ILE:CD1	2.50	0.40
10:J:14:VAL:HG13	10:J:49:LEU:CD2	2.51	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	1405/1984 (71%)	1316 (94%)	88 (6%)	1 (0%)	51 82
2	B	1122/1300 (86%)	1045 (93%)	77 (7%)	0	100 100
3	C	253/275 (92%)	238 (94%)	15 (6%)	0	100 100
4	D	126/184 (68%)	124 (98%)	2 (2%)	0	100 100
5	E	207/210 (99%)	189 (91%)	17 (8%)	1 (0%)	29 61
6	F	80/127 (63%)	74 (92%)	6 (8%)	0	100 100
7	G	169/172 (98%)	167 (99%)	2 (1%)	0	100 100
8	H	146/150 (97%)	139 (95%)	7 (5%)	0	100 100
9	I	112/125 (90%)	95 (85%)	17 (15%)	0	100 100
10	J	65/67 (97%)	60 (92%)	5 (8%)	0	100 100
11	K	112/117 (96%)	109 (97%)	3 (3%)	0	100 100
12	L	42/58 (72%)	37 (88%)	5 (12%)	0	100 100
15	Q	134/517 (26%)	126 (94%)	8 (6%)	0	100 100
16	R	133/249 (53%)	131 (98%)	2 (2%)	0	100 100
18	W	185/439 (42%)	185 (100%)	0	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
19	X	169/291 (58%)	167 (99%)	2 (1%)	0	100	100
All	All	4460/6265 (71%)	4202 (94%)	256 (6%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	345	GLY
5	E	84	ILE

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	1249/1763 (71%)	1248 (100%)	1 (0%)	93	98
2	B	990/1127 (88%)	988 (100%)	2 (0%)	93	98
3	C	234/252 (93%)	234 (100%)	0	100	100
4	D	118/160 (74%)	117 (99%)	1 (1%)	81	91
5	E	191/192 (100%)	190 (100%)	1 (0%)	88	94
6	F	71/111 (64%)	71 (100%)	0	100	100
7	G	152/153 (99%)	152 (100%)	0	100	100
8	H	129/131 (98%)	129 (100%)	0	100	100
9	I	103/112 (92%)	103 (100%)	0	100	100
10	J	56/56 (100%)	56 (100%)	0	100	100
11	K	104/106 (98%)	104 (100%)	0	100	100
12	L	41/55 (74%)	41 (100%)	0	100	100
15	Q	121/448 (27%)	120 (99%)	1 (1%)	81	91
16	R	118/218 (54%)	118 (100%)	0	100	100
18	W	169/373 (45%)	167 (99%)	2 (1%)	71	85
19	X	154/261 (59%)	154 (100%)	0	100	100
All	All	4000/5518 (72%)	3992 (100%)	8 (0%)	93	98

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	853	LYS
2	B	332	LYS
2	B	1131	ARG
4	D	24	LYS
5	E	47	LYS
15	Q	169	LYS
18	W	56	ARG
18	W	153	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	296	ASN
1	A	601	ASN
1	A	671	ASN
1	A	703	GLN
1	A	791	GLN
1	A	825	ASN
1	A	1077	ASN
1	A	1146	GLN
1	A	1360	ASN
2	B	344	GLN
2	B	503	ASN
2	B	654	GLN
2	B	699	HIS
2	B	1071	ASN
2	B	1094	GLN
4	D	34	ASN
4	D	102	ASN
7	G	122	ASN
9	I	91	HIS
11	K	2	ASN
15	Q	44	GLN
15	Q	53	ASN
15	Q	145	ASN
15	Q	168	ASN
16	R	91	GLN
18	W	109	HIS

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
14	P	13/14 (92%)	3 (23%)	2 (15%)

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
14	P	2	C
14	P	3	A
14	P	7	A

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
14	P	2	C
14	P	6	C

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

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 10 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)

There are no chain breaks in this entry.

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-19726. 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

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

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

7 Map analysis

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

7.1 Map-value distribution

This section was not generated.

7.2 Volume estimate versus contour level

This section was not generated.

7.3 Rotationally averaged power spectrum

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation (i)

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit

This section was not generated.