Full wwPDB EM Validation Report ⓘ

Dec 11, 2022 – 04:15 am GMT

PDB ID	:	6RQT
EMDB ID	:	EMD-4985
Title	:	RNA Polymerase I-tWH-Rrn3-DNA
Authors	:	Mueller, C.W.; Sadian, Y.; Tafur, L.
Deposited on	:	2019-05-16
Resolution	:	4.00 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:

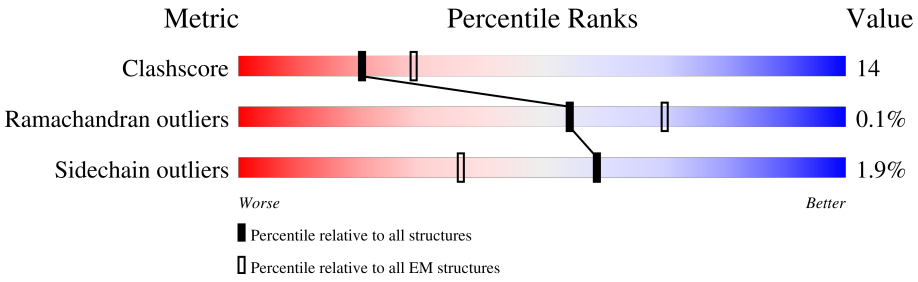
EMDB validation analysis	:	0.0.1.dev43
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

1 Overall quality at a glance

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

The reported resolution of this entry is 4.00 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

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% 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	T	70	13% 13% 13% 74%
2	U	70	11% · 16% 80%
3	A	1664	17% 62% 25% · 12%
4	B	1203	16% 64% 32% · ·
5	C	335	14% 62% 29% 9%
6	D	137	25% 28% 15% 57%
7	E	215	24% 74% 25%
8	F	155	12% 47% 17% · 35%

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Mol	Chain	Length	Quality of chain
9	G	326	32% 45% 17% 38%
10	H	146	16% 68% 23% 8%
11	I	125	18% 41% 10% 49%
12	J	70	17% 60% 37%
13	K	142	13% 49% 20% 30%
14	L	70	20% 43% 19% 39%
15	M	415	66% 69% 25% 6%
16	N	233	18% 38% 20% 42%
17	O	627	73% 60% 14% 26%

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 39820 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 DNA chain called Template strand.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	T	18	364	175	59	112	18	0	0

- Molecule 2 is a DNA chain called Nontemplate strand.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	U	14	293	139	59	81	14	0	0

- Molecule 3 is a protein called DNA-directed RNA polymerase I subunit RPA190.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	A	1466	11571	7309	2012	2188	62	0	0

- Molecule 4 is a protein called DNA-directed RNA polymerase I subunit RPA135.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	B	1170	9301	5888	1625	1737	51	0	0

- Molecule 5 is a protein called DNA-directed RNA polymerases I and III subunit RPAC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	C	304	2418	1536	414	460	8	0	0

- Molecule 6 is a protein called DNA-directed RNA polymerase I subunit RPA14.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
6	D	59	467	293	80	94	0	0

- Molecule 7 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	E	214	Total	C	N	O	S	0	0
			1751	1111	309	320	11		

- Molecule 8 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	F	100	Total	C	N	O	S	0	0
			823	522	144	154	3		

- Molecule 9 is a protein called DNA-directed RNA polymerase I subunit RPA43.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	G	202	Total	C	N	O	S	0	0
			1600	1026	276	293	5		

- Molecule 10 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	H	134	Total	C	N	O	S	0	0
			1075	677	182	212	4		

- Molecule 11 is a protein called DNA-directed RNA polymerase I subunit RPA12.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	I	64	Total	C	N	O	S	0	0
			472	295	78	95	4		

- Molecule 12 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC5.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	J	69	Total	C	N	O	S	0	0
			569	362	101	100	6		

- Molecule 13 is a protein called DNA-directed RNA polymerases I and III subunit RPAC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	K	100	Total	C	N	O	S	0	0
			785	491	129	160	5		

- Molecule 14 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC4.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	L	43	Total	C	N	O	S	0	0
			344	211	69	60	4		

- Molecule 15 is a protein called DNA-directed RNA polymerase I subunit RPA49.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	M	392	Total	C	N	O	S	0	0
			3100	1978	526	592	4		

- Molecule 16 is a protein called DNA-directed RNA polymerase I subunit RPA34.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	N	135	Total	C	N	O	S	0	0
			1070	685	175	206	4		

- Molecule 17 is a protein called RNA polymerase I-specific transcription initiation factor RRN3.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	O	463	Total	C	N	O	S	0	0
			3811	2473	623	694	21		

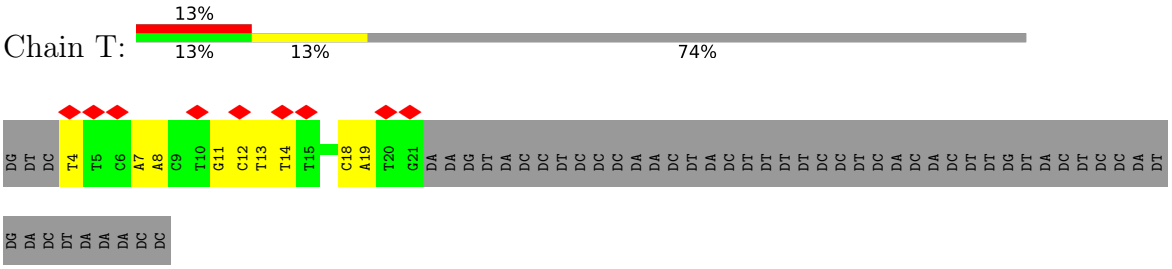
- Molecule 18 is ZINC ION (three-letter code: ZN) (formula: Zn).

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	

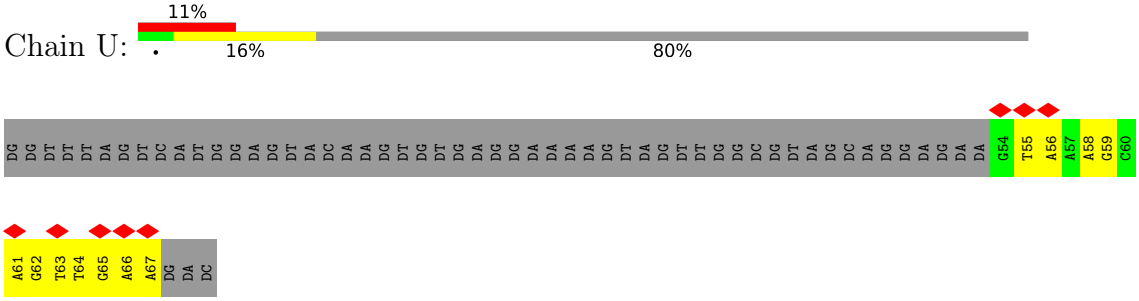
3 Residue-property plots (i)

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 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.

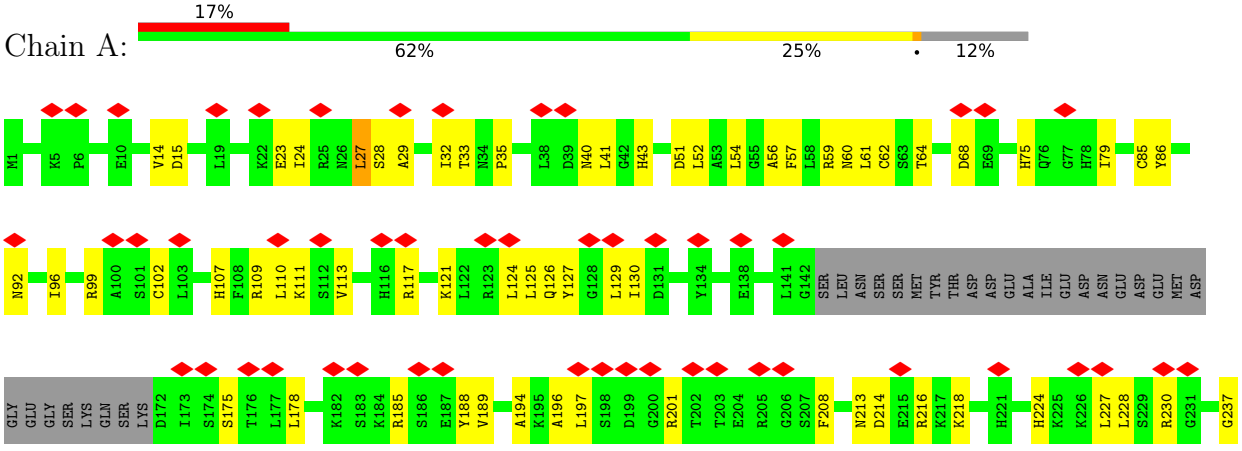
• Molecule 1: Template strand

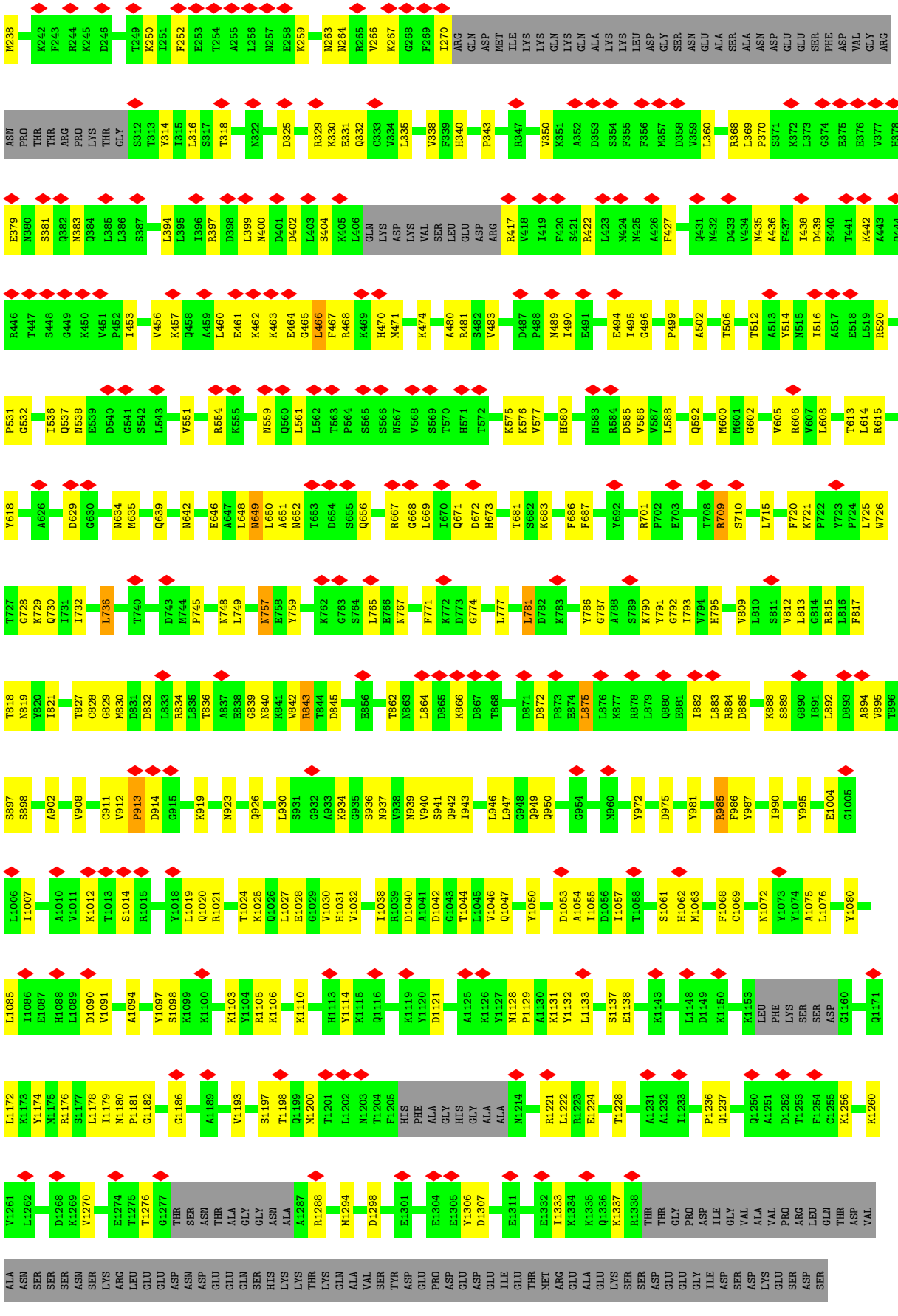


• Molecule 2: Nontemplate strand



• Molecule 3: DNA-directed RNA polymerase I subunit RPA190

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ASN PRO THR THR ARG PRO LYS THR GLY S312 T313 Y314 I315 L316 S317 T318 N322 D325 R329 K330 E331 Q332 C333 Y334 L335 V338 F339 H340 P343 R347 V350 K351 A352 D353 S354 F355 F356 M357 D358 V359 L360 R368 L369 P370 S371 K372 L373 G374 E375 E376 V377 H378
E379 N380 S381 Q382 N383 Q384 L385 L386 S387 L394 L395 I396 R397 D398 L399 N400 D401 D402 L403 S404 K405 L406 GLN LYS ASP LYS VAL SER LEU GLU ASP ARG R417 V418 I419 F420 S421 R422 L423 M424 N425 A426 F427 Q431 N432 D433 V434 N435 A436 F437 I438 D439 S440 T441 K442 A443 Q444 G445
R446 T447 S448 G449 K450 V451 P452 I453 V456 K457 Q458 A459 L460 E461 K462 K463 E464 G465 L466 F467 R468 K469 H470 M471 K474 A480 R481 S482 V483 D487 P488 N489 I490 E491 E494 I495 G496 P499 A502 T506 T512 A513 Y514 N515 I516 A517 E518 L519 R520 W530
P531 G532 I536 Q537 N538 E539 D540 G541 S542 L543 V551 R554 K555 N559 Q560 L561 L562 T563 P564 S565 S566 N567 V568 S569 T570 H571 T572 K575 K576 V577 H580 N583 R584 D585 V586 V587 L588 Q592 M600 M601 G602 V605 R606 V607 L608 T613 L614 R615
Y618 A626 D629 G630 N634 M635 Q639 N642 E646 A647 L648 N649 L650 A651 N652 T653 D654 S655 Q656 R667 G668 L669 I670 Q671 D672 H673 T681 S682 K683 F686 F687 Y692 R701 P702 E703 T708 R709 S710 L715 F720 K721 P722 Y723 P724 L725 W726
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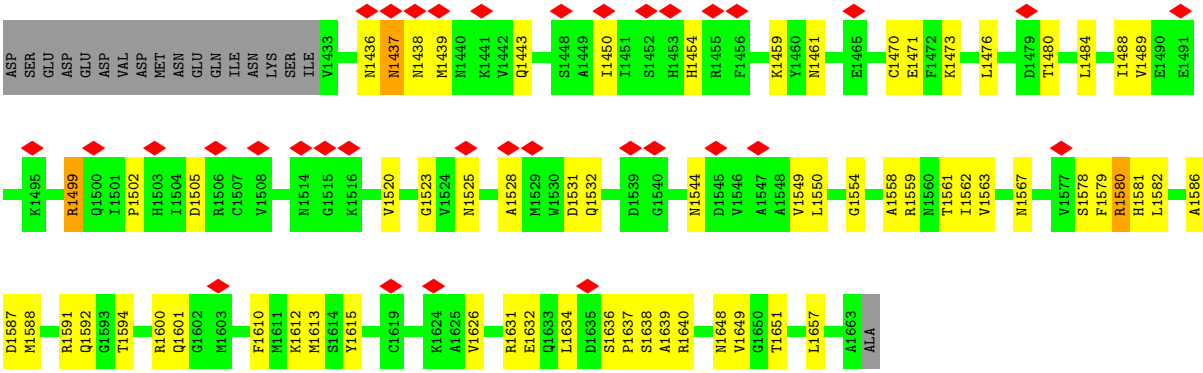
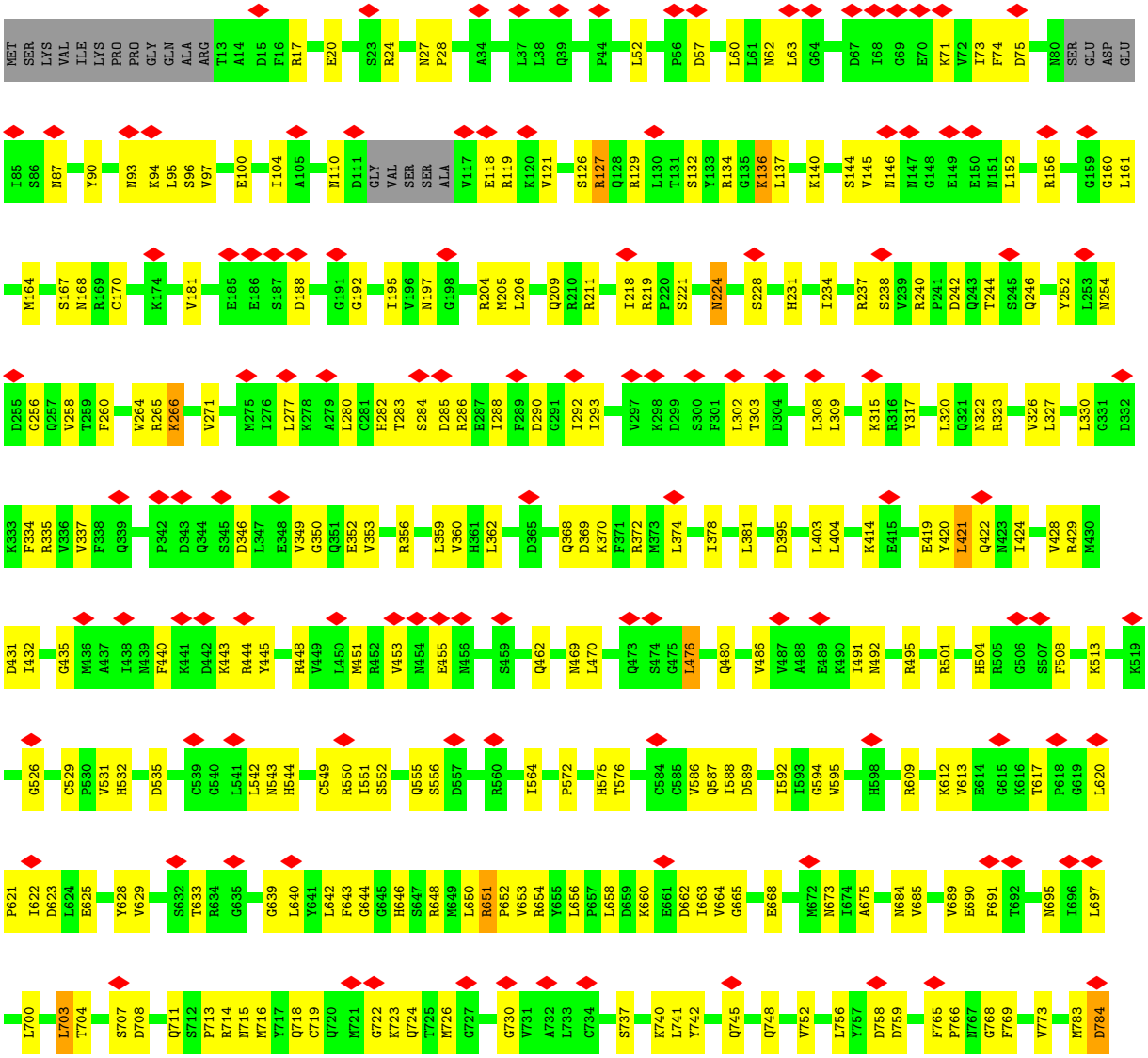
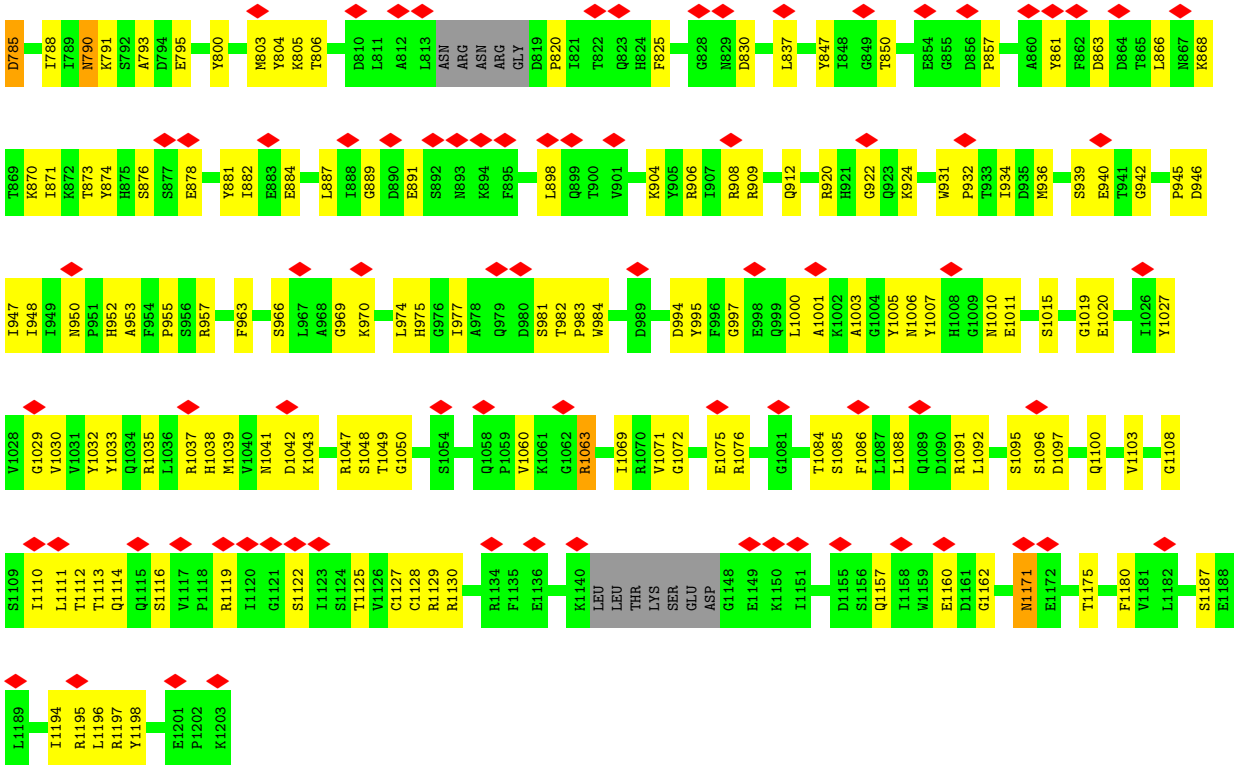
T818 N819 Y820 I821 T827 C828 G829 M830 D831 D832 L833 R834 L835 T836 A837 E838 G839 N840 K841 W842 R843 T844 D845 E856 T862 N863 L864 D865 K866 D867 T868 D871 D872 P873 E874 L875 L876 K877 R878 L879 Q880 E881 I882 L883 R884 D885 K888 S889 G890 I891 L892 D893 A894 V895 T896
S897 S898 A902 V908 C911 V912 P913 D914 G915 K919 N923 Q926 L930 S931 G932 A933 K934 G935 S936 N937 V938 N939 V940 S941 Q942 I943 L946 L947 G948 Q949 Q950 G954 M960 Y972 D975 Y981 R985 F986 Y987 I990 Y995 E1004 G1005
L1006 I1007 A1010 V1011 K1012 T1013 S1014 R1015 Y1018 L1019 Q1020 R1021 T1024 K1025 Q1026 L1027 E1028 G1029 V1030 H1031 V1032 I1038 R1039 D1040 A1041 D1042 G1043 T1044 L1045 V1046 Q1047 Y1050 D1053 A1054 I1055 D1056 I1057 T1058 S1061 H1062 M1063 F1068 C1069 N1072 Y1073 Y1074 A1075 L1076 Y1080
L1085 I1086 E1087 H1088 L1089 D1090 V1091 A1094 Y1097 S1098 K1099 K1100 K1103 Y1104 R1105 K1106 K1110 H1113 Y1114 K1115 Q1116 K1119 Y1120 D1121 A1125 K1126 Y1127 N1128 P1129 A1130 K1131 Y1132 L1133 S1137 E1138 K1143 L1148 D1149 K1150 K1153 LEU PHE LYS SER SER ASP G1160 Q1171
L1172 K1173 Y1174 M1175 R1176 S1177 L1178 I1179 N1180 P1181 G1182 G1186 A1189 V1193 S1197 T1198 Q1199 M1200 T1201 L1202 N1203 T1204 F1205 HIS PHE ALA GLY HIS GLY ALA ALA N1214 R1221 L1222 R1223 E1224 T1228 A1231 A1232 I1233 P1236 Q1237 Q1250 A1251 D1252 T1253 F1254 C1255 K1256 K1260
V1261 L1262 D1268 K1269 V1270 E1274 T1275 T1276 G1277 THR SER ASN THR ALA GLY GLY ASN ALA A1287 R1288 M1294 D1298 E1301 E1304 E1305 Y1306 D1307 E1311 E1332 I1333 K1334 K1335 Q1336 K1337 R1338 THR THR GLY PRO ASP ILE GLY VAL ALA VAL PRO ARG LEU GLN THR ASP VAL
ALA ASN SER SER SER ASN SER LYS ARG LEU GLU GLU ASP ASN ASP GLU GLN GLN SER HIS LYS LYS THR LYS GLN ALA VAL SER TYR ASP GLU PRO ASP GLU ASP GLU ILE GLU THR MET ARG GLU ALA GLU LYS SER SER ASP GLU GLU GLY ILE ASP SER ASP LYS GLU PRO SER ASP SER



• Molecule 4: DNA-directed RNA polymerase I subunit RPA135

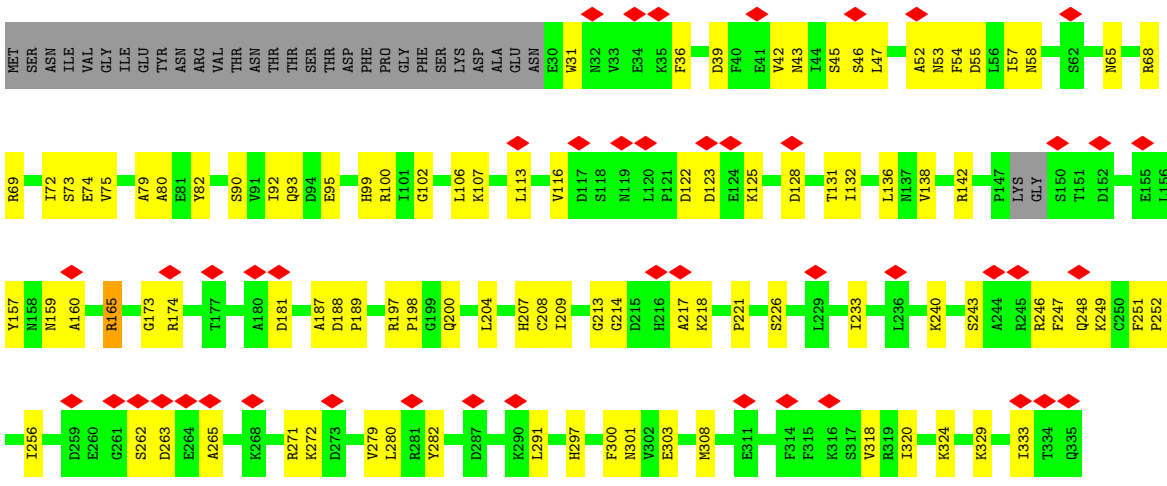
Chain B:





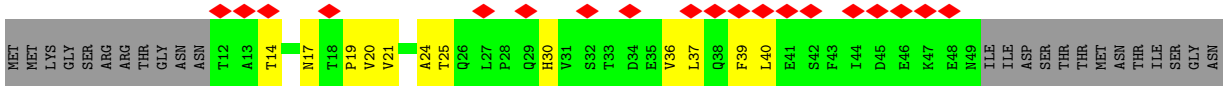
• Molecule 5: DNA-directed RNA polymerases I and III subunit RPAC1

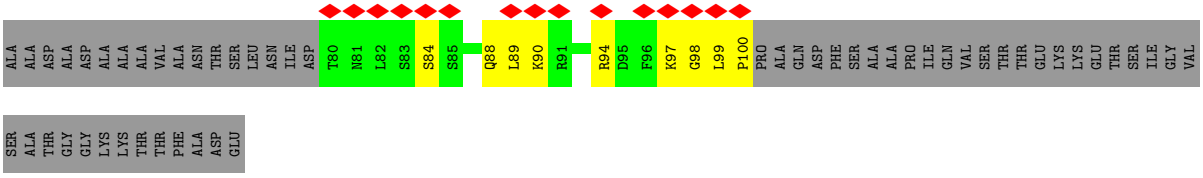
Chain C:



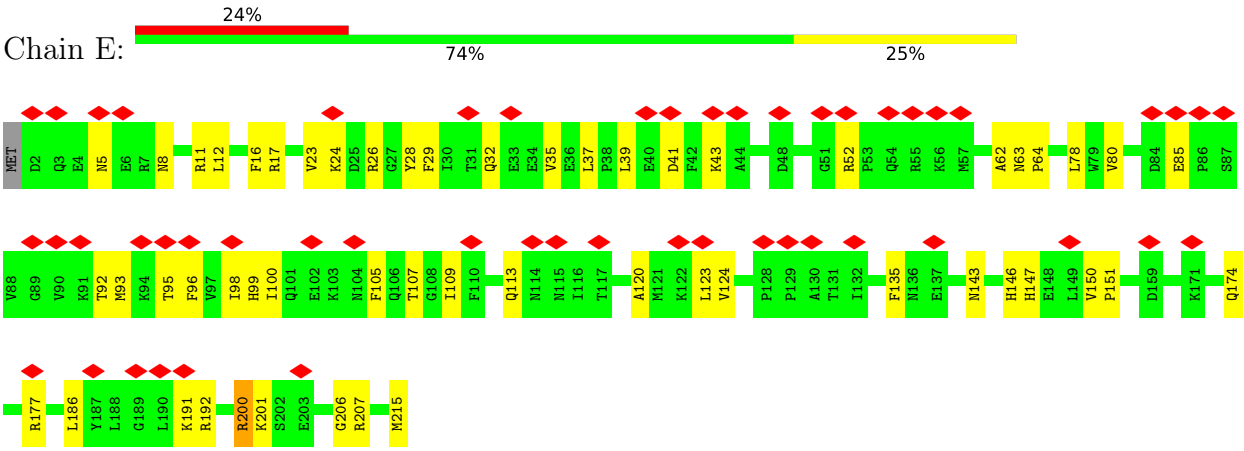
• Molecule 6: DNA-directed RNA polymerase I subunit RPA14

Chain D:

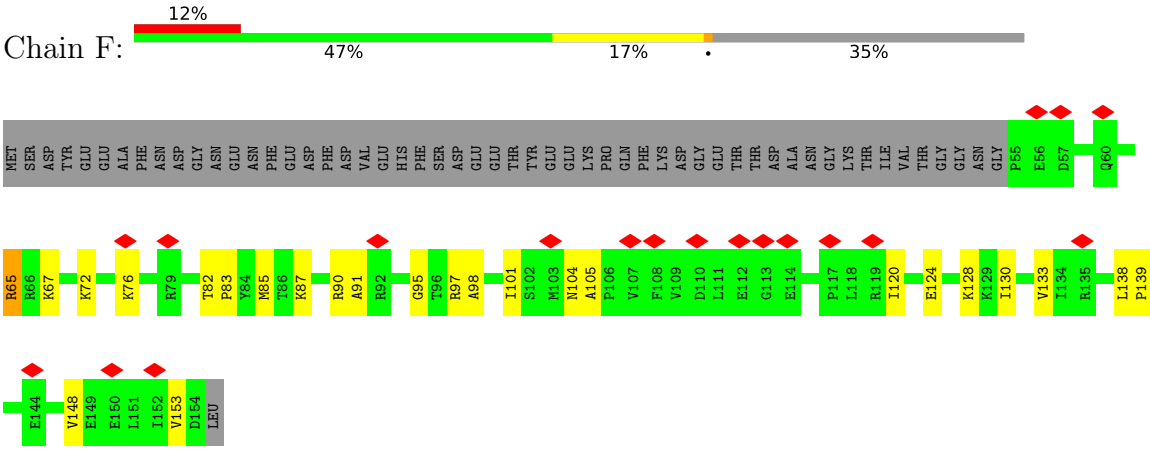




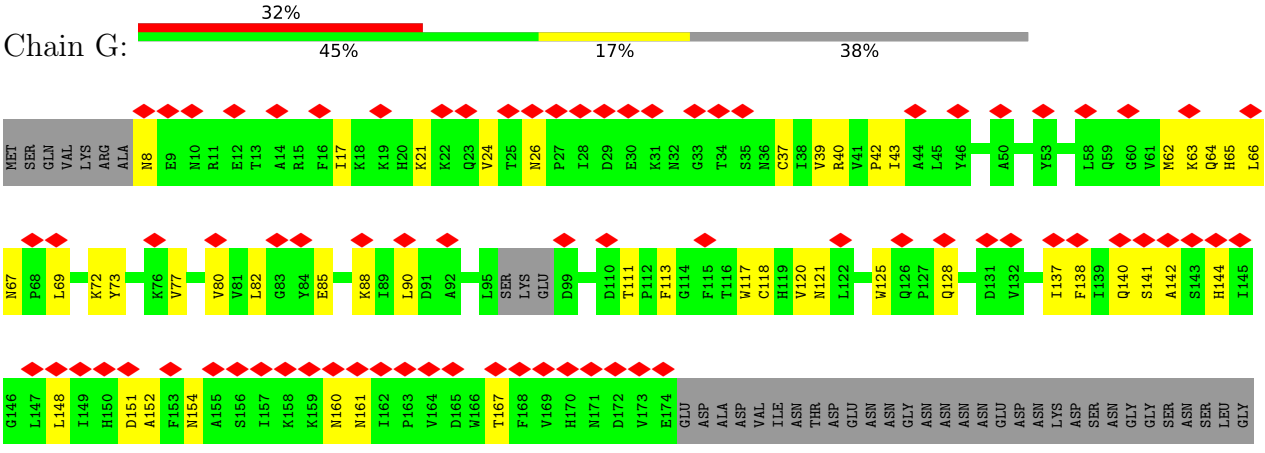
• Molecule 7: DNA-directed RNA polymerases I, II, and III subunit RPABC1

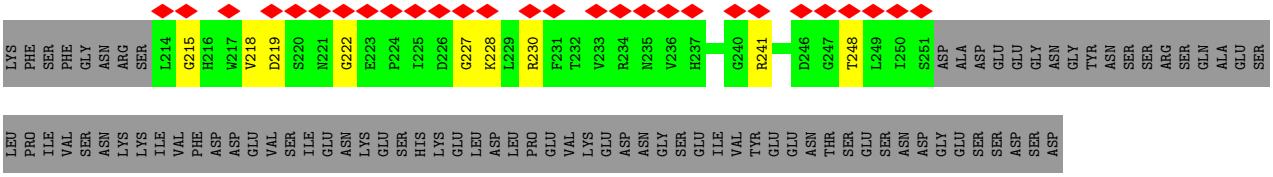


• Molecule 8: DNA-directed RNA polymerases I, II, and III subunit RPABC2



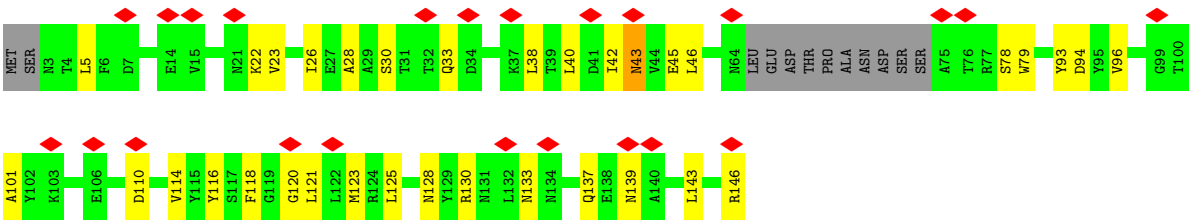
• Molecule 9: DNA-directed RNA polymerase I subunit RPA43





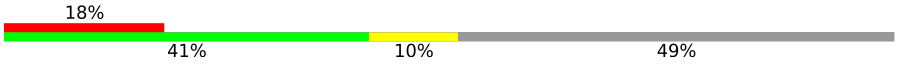
• Molecule 10: DNA-directed RNA polymerases I, II, and III subunit RPABC3

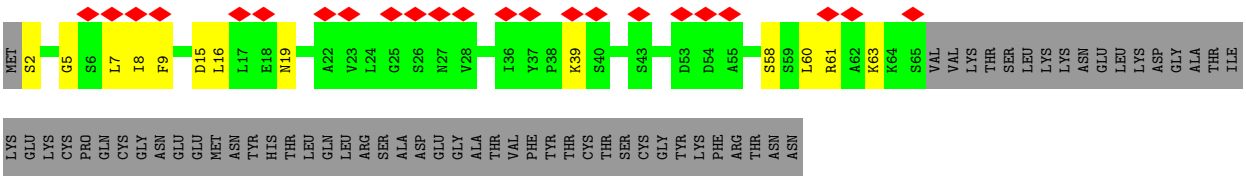
Chain H:



• Molecule 11: DNA-directed RNA polymerase I subunit RPA12

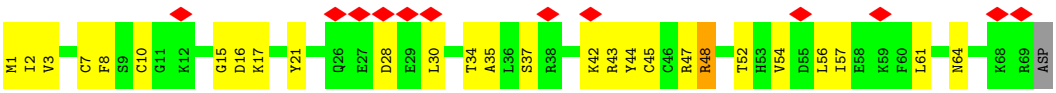
Chain I:





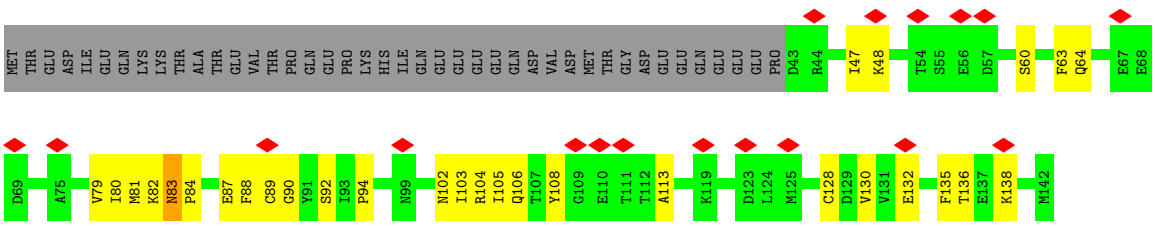
• Molecule 12: DNA-directed RNA polymerases I, II, and III subunit RPABC5

Chain J:



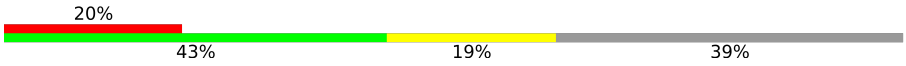
• Molecule 13: DNA-directed RNA polymerases I and III subunit RPAC2

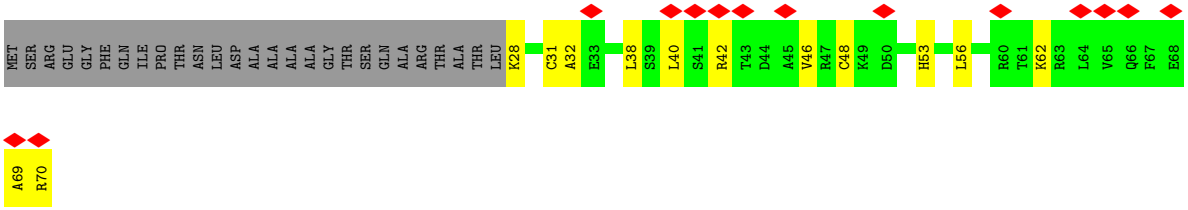
Chain K:



• Molecule 14: DNA-directed RNA polymerases I, II, and III subunit RPABC4

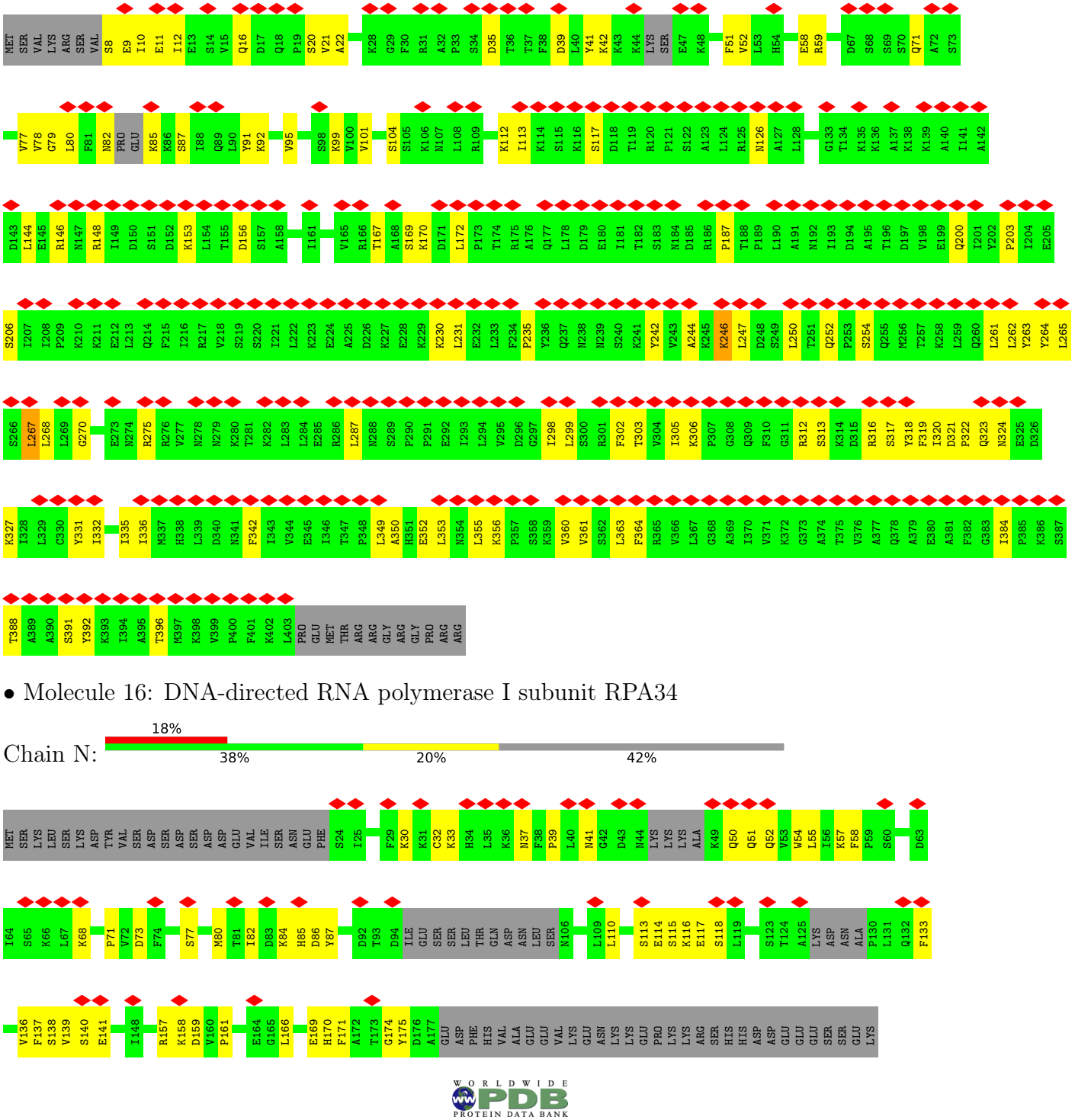
Chain L:





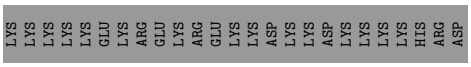
• Molecule 15: DNA-directed RNA polymerase I subunit RPA49

Chain M:

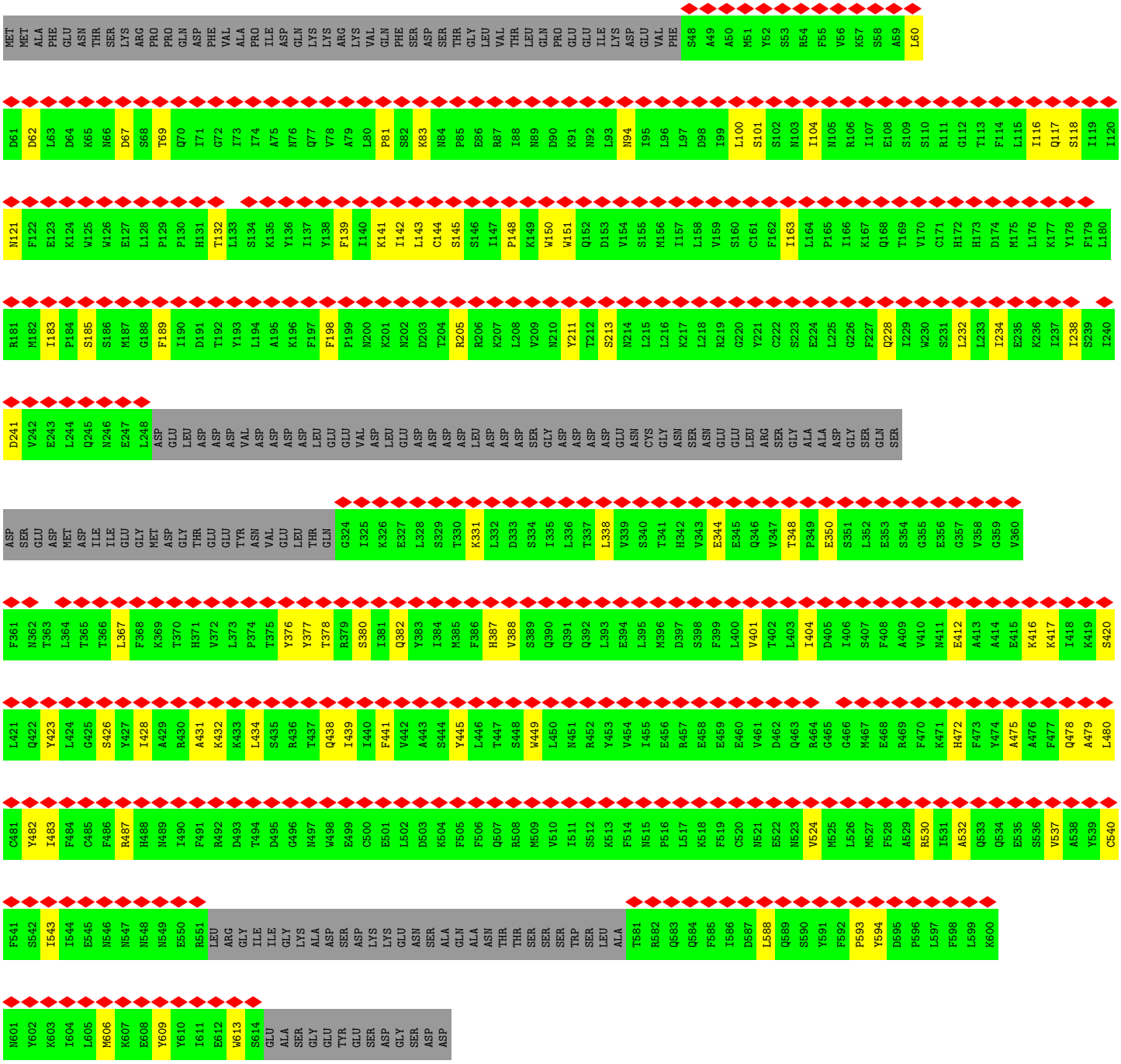


• Molecule 16: DNA-directed RNA polymerase I subunit RPA34

Chain N:



• Molecule 17: RNA polymerase I-specific transcription initiation factor RRN3



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	9789	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 ($e^-/Å^2$)	1.57175	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.286	Depositor
Minimum map value	-0.181	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.014	Depositor
Recommended contour level	0.05	Depositor
Map size (Å)	233.5168, 233.5168, 233.5168	wwPDB
Map dimensions	176, 176, 176	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.3268, 1.3268, 1.3268	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	T	0.54	0/405	0.93	1/622 (0.2%)
2	U	0.46	0/330	0.71	0/508
3	A	0.40	0/11782	0.67	0/15913
4	B	0.44	1/9506 (0.0%)	0.69	0/12847
5	C	0.39	0/2469	0.66	0/3347
6	D	0.31	0/473	0.62	0/641
7	E	0.40	0/1787	0.63	0/2406
8	F	0.38	0/838	0.64	0/1129
9	G	0.34	0/1637	0.62	0/2226
10	H	0.41	0/1093	0.66	0/1480
11	I	0.36	0/478	0.72	0/647
12	J	0.48	0/578	0.76	0/775
13	K	0.38	0/795	0.63	0/1072
14	L	0.33	0/346	0.67	0/457
15	M	0.33	0/3150	0.65	0/4247
16	N	0.38	0/1090	0.71	0/1466
17	O	0.28	0/3897	0.55	0/5268
All	All	0.39	1/40654 (0.0%)	0.66	1/55051 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	121	VAL	C-N	-5.01	1.22	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	T	11	DG	P-O3'-C3'	5.31	126.07	119.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	T	364	0	206	17	0
2	U	293	0	158	15	0
3	A	11571	0	11653	369	0
4	B	9301	0	9193	343	0
5	C	2418	0	2401	74	0
6	D	467	0	468	16	0
7	E	1751	0	1776	36	0
8	F	823	0	841	16	0
9	G	1600	0	1600	38	0
10	H	1075	0	1046	24	0
11	I	472	0	474	7	0
12	J	569	0	587	23	0
13	K	785	0	782	23	0
14	L	344	0	365	8	0
15	M	3100	0	3210	109	0
16	N	1070	0	1085	61	0
17	O	3811	0	3804	53	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	39820	0	39649	1077	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 14.

All (1077) 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)
3:A:912:VAL:HG22	3:A:913:PRO:CD	1.52	1.37
4:B:1005:TYR:HE1	16:N:171:PHE:CZ	1.54	1.26

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:618:TYR:CD1	4:B:783:MET:HE1	1.70	1.25
3:A:41:LEU:HD12	3:A:43:HIS:ND1	1.53	1.21
4:B:1005:TYR:OH	16:N:169:GLU:HG3	1.39	1.21
4:B:1005:TYR:CE1	16:N:171:PHE:CZ	2.31	1.18
4:B:1005:TYR:HE1	16:N:171:PHE:CE2	1.63	1.14
4:B:1005:TYR:CE1	16:N:171:PHE:CE2	2.37	1.12
15:M:262:LEU:CG	15:M:265:LEU:HD12	1.78	1.11
3:A:32:ILE:HD11	3:A:54:LEU:HD12	1.30	1.10
3:A:912:VAL:CG2	3:A:913:PRO:HD2	1.82	1.08
4:B:470:LEU:HD11	4:B:476:LEU:HD11	1.27	1.08
3:A:466:LEU:HD22	3:A:470:HIS:ND1	1.68	1.07
3:A:985:ARG:HD3	3:A:986:PHE:N	1.68	1.07
3:A:669:LEU:HD13	3:A:813:LEU:HD12	1.36	1.07
15:M:262:LEU:HG	15:M:265:LEU:CD1	1.88	1.04
4:B:359:LEU:HD11	4:B:370:LYS:HG3	1.37	1.03
3:A:618:TYR:CD1	4:B:783:MET:CE	2.42	1.02
16:N:169:GLU:OE2	16:N:170:HIS:N	1.94	1.01
15:M:262:LEU:HG	15:M:265:LEU:HD12	1.03	0.99
3:A:912:VAL:CG2	3:A:913:PRO:CD	2.41	0.98
3:A:912:VAL:HG22	3:A:913:PRO:HD2	0.99	0.97
3:A:912:VAL:HG22	3:A:913:PRO:HD3	1.46	0.95
3:A:618:TYR:HD1	4:B:783:MET:HE1	1.22	0.94
3:A:985:ARG:NH2	3:A:986:PHE:HB2	1.83	0.94
15:M:262:LEU:HA	15:M:265:LEU:HG	1.47	0.94
4:B:1196:LEU:HD13	4:B:1197:ARG:N	1.84	0.93
15:M:262:LEU:HA	15:M:265:LEU:CG	2.00	0.92
5:C:45:SER:HB3	5:C:53:ASN:HB3	1.53	0.91
3:A:32:ILE:HD11	3:A:54:LEU:CD1	2.00	0.90
3:A:942:GLN:O	3:A:985:ARG:CZ	2.21	0.88
3:A:466:LEU:CD2	3:A:470:HIS:ND1	2.35	0.88
4:B:470:LEU:CD1	4:B:476:LEU:HD11	2.04	0.86
4:B:359:LEU:CD1	4:B:370:LYS:HE3	2.06	0.85
3:A:1178:LEU:HD13	3:A:1179:ILE:N	1.92	0.84
4:B:703:LEU:O	4:B:703:LEU:HD22	1.77	0.84
4:B:703:LEU:HD13	4:B:704:THR:N	1.93	0.83
4:B:974:LEU:HD11	16:N:169:GLU:CG	2.09	0.83
4:B:974:LEU:HD21	16:N:169:GLU:HG2	1.60	0.83
4:B:327:LEU:HD23	4:B:330:LEU:HD12	1.61	0.82
3:A:669:LEU:HD13	3:A:813:LEU:CD1	2.11	0.80
15:M:250:LEU:HD23	15:M:250:LEU:H	1.46	0.80
3:A:943:ILE:HA	3:A:985:ARG:NH1	1.96	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1306:TYR:O	3:A:1307:ASP:OD1	2.00	0.80
4:B:820:PRO:HD2	15:M:356:LYS:NZ	1.97	0.79
3:A:669:LEU:CD1	3:A:813:LEU:HD12	2.11	0.79
4:B:974:LEU:CD2	16:N:169:GLU:HG2	2.14	0.77
2:U:62:DG:H1'	2:U:63:DT:H5'	1.67	0.76
4:B:359:LEU:HD11	4:B:370:LYS:CG	2.15	0.76
3:A:943:ILE:HA	3:A:985:ARG:HH11	1.48	0.76
2:U:66:DA:H1'	2:U:67:DA:H5'	1.68	0.76
15:M:262:LEU:CB	15:M:265:LEU:HD12	2.16	0.75
4:B:359:LEU:HD12	4:B:359:LEU:O	1.87	0.74
3:A:41:LEU:HD13	3:A:41:LEU:O	1.86	0.74
7:E:186:LEU:O	7:E:186:LEU:HD23	1.88	0.74
4:B:1005:TYR:HE1	16:N:171:PHE:CE1	2.06	0.73
4:B:1005:TYR:CD1	16:N:171:PHE:CE2	2.76	0.73
3:A:54:LEU:O	3:A:54:LEU:HD23	1.87	0.73
3:A:1030:VAL:HG12	3:A:1040:ASP:HA	1.71	0.73
15:M:262:LEU:HA	15:M:265:LEU:CD1	2.18	0.73
3:A:461:GLU:HB3	3:A:464:GLU:HA	1.70	0.73
3:A:985:ARG:HH21	3:A:986:PHE:H	1.34	0.73
4:B:714:ARG:HD2	4:B:957:ARG:HD2	1.72	0.72
1:T:13:DT:OP2	1:T:13:DT:H71	1.90	0.72
15:M:268:LEU:HG	15:M:331:TYR:CE2	2.24	0.72
7:E:135:PHE:CD1	7:E:186:LEU:HD21	2.25	0.71
5:C:165:ARG:HB3	5:C:189:PRO:HB3	1.71	0.71
10:H:116:TYR:HB3	10:H:123:MET:HB2	1.71	0.71
4:B:359:LEU:HD11	4:B:370:LYS:CE	2.21	0.70
5:C:53:ASN:ND2	16:N:174:GLY:O	2.24	0.70
2:U:61:DA:H4'	3:A:1601:GLN:HE22	1.57	0.69
3:A:985:ARG:CD	3:A:987:TYR:H	2.04	0.69
3:A:930:LEU:HD13	3:A:930:LEU:O	1.92	0.69
3:A:875:LEU:O	3:A:875:LEU:HD22	1.93	0.69
4:B:359:LEU:CD1	4:B:370:LYS:CE	2.71	0.69
4:B:820:PRO:HD2	15:M:356:LYS:HZ2	1.55	0.69
3:A:875:LEU:C	3:A:875:LEU:HD13	2.13	0.69
3:A:759:TYR:CE1	3:A:913:PRO:HG2	2.27	0.69
3:A:875:LEU:O	3:A:875:LEU:HD13	1.93	0.69
3:A:930:LEU:O	3:A:930:LEU:HD22	1.93	0.68
3:A:926:GLN:O	3:A:930:LEU:HB3	1.94	0.68
3:A:985:ARG:HD3	3:A:985:ARG:C	2.13	0.68
16:N:169:GLU:CD	16:N:170:HIS:N	2.47	0.68
15:M:363:LEU:HD13	15:M:363:LEU:O	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:M:250:LEU:HD23	15:M:250:LEU:N	2.08	0.68
3:A:1256:LYS:CB	3:A:1307:ASP:OD2	2.43	0.67
3:A:943:ILE:HD12	3:A:985:ARG:NH1	2.10	0.67
9:G:148:LEU:HB3	9:G:151:ASP:HA	1.76	0.67
3:A:1443:GLN:NE2	3:A:1461:ASN:OD1	2.28	0.67
4:B:359:LEU:HD11	4:B:370:LYS:HE3	1.76	0.67
4:B:803:MET:SD	4:B:909:ARG:NH1	2.68	0.66
4:B:531:VAL:HG23	4:B:715:ASN:HB3	1.76	0.66
4:B:1005:TYR:OH	16:N:169:GLU:CG	2.31	0.66
1:T:13:DT:H3	2:U:56:DA:H2	1.44	0.66
15:M:363:LEU:O	15:M:363:LEU:HD22	1.95	0.66
4:B:96:SER:HB3	4:B:144:SER:HB3	1.76	0.66
3:A:1436:ASN:HB2	3:A:1461:ASN:HD21	1.60	0.66
4:B:703:LEU:HD13	4:B:703:LEU:C	2.16	0.66
4:B:362:LEU:HD12	4:B:369:ASP:HB3	1.77	0.66
4:B:110:ASN:HD21	4:B:118:GLU:HA	1.61	0.65
4:B:589:ASP:HB3	4:B:643:PHE:HA	1.78	0.65
3:A:126:GLN:HB2	3:A:343:PRO:HD3	1.79	0.65
3:A:813:LEU:HD13	3:A:813:LEU:C	2.17	0.65
5:C:197:ARG:HE	12:J:61:LEU:HB3	1.61	0.65
4:B:866:LEU:HD23	4:B:868:LYS:HD3	1.77	0.65
3:A:943:ILE:HD12	3:A:985:ARG:HH11	1.61	0.65
3:A:1634:LEU:HB2	3:A:1639:ALA:HB1	1.79	0.65
3:A:828:CYS:SG	3:A:829:GLY:N	2.70	0.65
4:B:1047:ARG:HE	4:B:1050:GLY:HA3	1.61	0.65
4:B:1196:LEU:HD13	4:B:1196:LEU:C	2.16	0.65
3:A:985:ARG:HD2	3:A:987:TYR:H	1.62	0.64
4:B:942:GLY:HA2	5:C:226:SER:HB3	1.80	0.64
4:B:1006:ASN:ND2	4:B:1010:ASN:OD1	2.29	0.64
3:A:40:ASN:O	3:A:41:LEU:HB3	1.97	0.64
4:B:1005:TYR:CE1	16:N:171:PHE:CE1	2.85	0.64
3:A:813:LEU:HD13	3:A:813:LEU:O	1.98	0.64
15:M:262:LEU:HA	15:M:265:LEU:HD12	1.78	0.64
7:E:78:LEU:HD11	7:E:109:ILE:HG12	1.78	0.64
3:A:618:TYR:CG	4:B:783:MET:HE1	2.28	0.64
3:A:1085:LEU:HD11	3:A:1176:ARG:HH11	1.62	0.64
3:A:1138:GLU:HB3	7:E:206:GLY:HA3	1.80	0.64
3:A:1470:CYS:SG	3:A:1471:GLU:N	2.71	0.64
4:B:974:LEU:CD1	16:N:169:GLU:HG2	2.27	0.64
1:T:4:DT:O2	2:U:67:DA:H2	1.80	0.64
4:B:429:ARG:HA	4:B:432:ILE:HG22	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:O:387:HIS:ND1	17:O:606:MET:SD	2.72	0.63
4:B:303:THR:HG21	11:I:5:GLY:HA3	1.79	0.63
4:B:1196:LEU:HD11	4:B:1198:TYR:CE1	2.34	0.63
4:B:134:ARG:HE	4:B:160:GLY:HA3	1.63	0.63
3:A:1178:LEU:HD13	3:A:1178:LEU:C	2.18	0.63
4:B:246:GLN:HE22	4:B:360:VAL:HG12	1.64	0.63
4:B:625:GLU:HB2	4:B:648:ARG:HH11	1.63	0.63
7:E:96:PHE:HA	7:E:99:HIS:HD2	1.64	0.63
17:O:432:LYS:HB3	17:O:609:TYR:HD1	1.63	0.63
17:O:62:ASP:HB3	17:O:67:ASP:HB3	1.80	0.63
3:A:15:ASP:O	4:B:1196:LEU:HD22	1.99	0.62
12:J:10:CYS:HB3	12:J:43:ARG:HH21	1.64	0.62
4:B:703:LEU:HD22	4:B:703:LEU:C	2.19	0.62
4:B:974:LEU:HD21	16:N:169:GLU:CG	2.30	0.62
5:C:80:ALA:HA	5:C:208:CYS:HA	1.80	0.62
3:A:41:LEU:HD11	15:M:323:GLN:HA	1.82	0.62
3:A:456:VAL:O	3:A:460:LEU:N	2.33	0.62
4:B:740:LYS:HE3	4:B:805:LYS:HG2	1.80	0.62
15:M:268:LEU:C	15:M:268:LEU:HD23	2.20	0.62
3:A:872:ASP:HB2	3:A:875:LEU:HB3	1.82	0.62
4:B:327:LEU:HD23	4:B:330:LEU:CD1	2.29	0.62
7:E:28:TYR:HA	7:E:64:PRO:HA	1.82	0.62
3:A:463:LYS:O	3:A:468:ARG:NH1	2.33	0.62
4:B:424:ILE:HG22	4:B:453:VAL:HG11	1.82	0.62
3:A:490:ILE:HG23	3:A:494:GLU:HB2	1.82	0.62
3:A:32:ILE:CD1	3:A:54:LEU:CD1	2.78	0.61
3:A:439:ASP:O	3:A:442:LYS:NZ	2.33	0.61
3:A:985:ARG:HD2	3:A:987:TYR:HB3	1.82	0.61
3:A:652:ASN:O	3:A:656:GLN:NE2	2.33	0.61
4:B:745:GLN:HE21	4:B:800:TYR:HB3	1.65	0.61
4:B:974:LEU:HD11	16:N:169:GLU:HG2	1.79	0.61
6:D:37:LEU:HD12	6:D:97:LYS:HE3	1.81	0.61
5:C:122:ASP:HA	5:C:125:LYS:HB3	1.82	0.61
8:F:85:MET:HB3	8:F:153:VAL:HA	1.83	0.61
13:K:105:ILE:HD11	13:K:113:ALA:HA	1.83	0.61
5:C:246:ARG:HA	5:C:249:LYS:HE2	1.82	0.61
3:A:175:SER:HA	3:A:178:LEU:HD23	1.82	0.61
3:A:531:PRO:HB3	3:A:580:HIS:HB2	1.81	0.61
4:B:211:ARG:HH22	4:B:646:HIS:HB2	1.66	0.61
9:G:82:LEU:HD11	9:G:125:TRP:HB2	1.81	0.61
4:B:104:ILE:HG23	4:B:137:LEU:HD23	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:G:241:ARG:HG2	17:O:189:PHE:HB3	1.83	0.61
4:B:1003:ALA:HB1	4:B:1005:TYR:CD2	2.36	0.60
3:A:618:TYR:CG	4:B:783:MET:CE	2.83	0.60
4:B:285:ASP:HA	11:I:15:ASP:HB3	1.81	0.60
5:C:45:SER:HB3	5:C:53:ASN:CB	2.29	0.60
4:B:974:LEU:HD11	16:N:169:GLU:CB	2.31	0.60
17:O:116:ILE:HG23	17:O:143:LEU:HD21	1.83	0.60
3:A:912:VAL:O	3:A:914:ASP:N	2.33	0.60
4:B:63:LEU:O	4:B:63:LEU:HD23	2.01	0.60
3:A:618:TYR:HA	4:B:783:MET:HE1	1.81	0.60
17:O:213:SER:HB2	17:O:338:LEU:HD21	1.81	0.60
7:E:174:GLN:NE2	8:F:140:ASP:OD2	2.32	0.60
17:O:417:LYS:HD3	17:O:472:HIS:HB2	1.83	0.60
4:B:741:LEU:H	4:B:804:TYR:HB2	1.67	0.60
6:D:30:HIS:HA	9:G:39:VAL:HG23	1.84	0.60
3:A:1080:TYR:HB3	3:A:1172:LEU:HD22	1.84	0.59
15:M:262:LEU:CA	15:M:265:LEU:HD12	2.32	0.59
16:N:80:MET:HB2	16:N:87:TYR:HB2	1.83	0.59
17:O:118:SER:HA	17:O:121:ASN:HB2	1.82	0.59
3:A:985:ARG:CD	3:A:987:TYR:N	2.65	0.59
4:B:17:ARG:NH1	4:B:758:ASP:OD2	2.35	0.59
4:B:404:LEU:HD11	4:B:551:ILE:HG21	1.85	0.59
4:B:492:ASN:HB3	4:B:495:ARG:HG3	1.83	0.59
5:C:333:ILE:HG22	13:K:47:ILE:HG12	1.83	0.59
7:E:100:ILE:HG23	7:E:105:PHE:HB2	1.84	0.59
10:H:22:LYS:O	10:H:43:ASN:ND2	2.35	0.59
3:A:130:ILE:O	7:E:192:ARG:NH2	2.36	0.59
3:A:912:VAL:CG2	3:A:913:PRO:HD3	2.24	0.59
4:B:1039:MET:HG2	4:B:1042:ASP:H	1.65	0.59
3:A:85:CYS:SG	3:A:86:TYR:N	2.75	0.59
3:A:792:GLY:H	3:A:795:HIS:HB3	1.67	0.59
12:J:30:LEU:HB3	12:J:35:ALA:HB2	1.85	0.59
3:A:370:PRO:HB3	3:A:379:GLU:HA	1.85	0.59
15:M:268:LEU:HG	15:M:331:TYR:HE2	1.67	0.59
3:A:1459:LYS:HD3	3:A:1473:LYS:HE2	1.83	0.59
4:B:791:LYS:NZ	5:C:214:GLY:O	2.35	0.59
12:J:42:LYS:HG3	12:J:43:ARG:HG3	1.84	0.59
3:A:466:LEU:HD13	3:A:471:MET:SD	2.42	0.59
4:B:718:GLN:HB2	4:B:922:GLY:HA2	1.84	0.59
5:C:80:ALA:HB3	5:C:102:GLY:HA2	1.83	0.59
15:M:268:LEU:HG	15:M:331:TYR:CD2	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:332:GLN:HE22	3:A:350:VAL:HG22	1.68	0.59
3:A:465:GLY:HA2	3:A:468:ARG:HB3	1.84	0.59
3:A:588:LEU:HD21	3:A:600:MET:HG3	1.84	0.59
3:A:335:LEU:HA	3:A:338:VAL:HG12	1.85	0.59
4:B:100:GLU:HB3	4:B:140:LYS:HD2	1.85	0.59
6:D:24:ALA:HA	9:G:43:ILE:HG22	1.85	0.59
7:E:62:ALA:HB3	7:E:78:LEU:HB3	1.85	0.59
3:A:748:ASN:ND2	3:A:771:PHE:O	2.36	0.58
3:A:1600:ARG:HG3	3:A:1601:GLN:HG3	1.85	0.58
5:C:318:VAL:HG13	13:K:128:CYS:HB2	1.85	0.58
15:M:313:SER:O	17:O:117:GLN:NE2	2.36	0.58
3:A:1276:THR:OG1	3:A:1288:ARG:NH1	2.36	0.58
12:J:44:TYR:HA	12:J:47:ARG:HG2	1.84	0.58
3:A:892:LEU:HA	3:A:895:VAL:HG12	1.85	0.58
4:B:136:LYS:HA	4:B:160:GLY:HA2	1.83	0.58
3:A:1129:PRO:HB2	3:A:1178:LEU:HD23	1.84	0.58
5:C:248:GLN:HG3	5:C:256:ILE:HB	1.85	0.58
4:B:609:ARG:NH2	4:B:662:ASP:OD2	2.37	0.58
3:A:1180:ASN:HD22	8:F:87:LYS:HG2	1.69	0.58
15:M:42:LYS:HB2	16:N:30:LYS:HB2	1.86	0.58
3:A:41:LEU:HD13	15:M:322:PRO:O	2.03	0.58
3:A:79:ILE:HG22	3:A:360:LEU:H	1.67	0.58
4:B:277:LEU:HD23	4:B:374:LEU:HD12	1.84	0.58
15:M:187:PRO:HB3	15:M:327:LYS:HG2	1.85	0.58
17:O:144:CYS:HB3	17:O:151:TRP:HB2	1.85	0.58
17:O:431:ALA:O	17:O:487:ARG:NH2	2.36	0.58
3:A:496:GLY:HA3	3:A:608:LEU:HD13	1.84	0.58
3:A:936:SER:H	3:A:939:ASN:HB2	1.68	0.58
4:B:327:LEU:CD2	4:B:330:LEU:HD12	2.33	0.58
4:B:966:SER:HB3	4:B:1029:GLY:HA3	1.85	0.58
12:J:7:CYS:SG	12:J:8:PHE:N	2.77	0.58
4:B:205:MET:SD	4:B:206:LEU:N	2.77	0.58
5:C:157:TYR:HB2	5:C:160:ALA:HB2	1.86	0.58
17:O:532:ALA:HA	17:O:537:VAL:HB	1.85	0.58
3:A:28:SER:O	4:B:1129:ARG:NH1	2.37	0.57
3:A:671:GLN:O	4:B:952:HIS:NE2	2.37	0.57
3:A:1626:VAL:HG21	4:B:1194:ILE:HG12	1.86	0.57
5:C:142:ARG:HE	5:C:198:PRO:HG3	1.69	0.57
9:G:42:PRO:HA	9:G:121:ASN:HA	1.86	0.57
10:H:114:VAL:HG22	10:H:125:LEU:HB2	1.86	0.57
12:J:17:LYS:O	12:J:21:TYR:N	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:240:ARG:NH2	4:B:356:ARG:O	2.37	0.57
15:M:77:VAL:HG12	15:M:92:LYS:HA	1.86	0.57
4:B:372:ARG:NH1	4:B:592:ILE:O	2.35	0.57
10:H:116:TYR:OH	10:H:137:GLN:NE2	2.37	0.57
3:A:512:THR:HB	3:A:514:TYR:HB3	1.86	0.57
4:B:887:LEU:HD12	14:L:56:LEU:HB2	1.85	0.57
4:B:1003:ALA:O	4:B:1005:TYR:CD2	2.57	0.57
3:A:1028:GLU:OE2	3:A:1638:SER:N	2.36	0.57
3:A:1068:PHE:O	3:A:1072:ASN:ND2	2.37	0.57
10:H:110:ASP:OD2	10:H:128:ASN:ND2	2.37	0.57
13:K:89:CYS:SG	13:K:90:GLY:N	2.78	0.57
15:M:80:LEU:HB3	16:N:51:GLN:HE21	1.69	0.57
1:T:13:DT:OP2	3:A:464:GLU:OE2	2.22	0.57
3:A:985:ARG:HD3	3:A:987:TYR:H	1.70	0.57
15:M:11:GLU:OE2	15:M:87:SER:OG	2.22	0.57
4:B:613:VAL:HB	4:B:660:LYS:HD2	1.86	0.57
4:B:1003:ALA:HB1	4:B:1005:TYR:HD2	1.70	0.57
4:B:1005:TYR:HE1	16:N:171:PHE:CD2	2.20	0.57
4:B:675:ALA:HB3	4:B:689:VAL:HG12	1.86	0.56
4:B:791:LYS:HE3	4:B:932:PRO:HD3	1.87	0.56
4:B:352:GLU:HG3	4:B:356:ARG:HE	1.71	0.56
4:B:784:ASP:C	4:B:950:ASN:ND2	2.58	0.56
4:B:788:ILE:HG22	4:B:948:ILE:HB	1.86	0.56
2:U:56:DA:N3	4:B:513:LYS:NZ	2.49	0.56
3:A:28:SER:OG	3:A:29:ALA:N	2.39	0.56
3:A:650:LEU:HD23	4:B:1084:THR:HB	1.87	0.56
4:B:73:ILE:HD13	4:B:428:VAL:HB	1.86	0.56
4:B:994:ASP:OD1	4:B:1007:TYR:OH	2.24	0.56
9:G:88:LYS:H	9:G:120:VAL:HG12	1.70	0.56
16:N:54:TRP:HA	16:N:133:PHE:HE1	1.71	0.56
16:N:169:GLU:CD	16:N:170:HIS:H	2.07	0.56
1:T:18:DC:N3	1:T:19:DA:N6	2.53	0.56
4:B:197:ASN:HD21	4:B:462:GLN:HE21	1.53	0.56
4:B:240:ARG:HD2	4:B:244:THR:HG23	1.88	0.56
5:C:240:LYS:HD3	5:C:265:ALA:H	1.70	0.56
9:G:154:ASN:HD22	17:O:148:PRO:HG3	1.71	0.56
10:H:30:SER:OG	10:H:33:GLN:O	2.24	0.56
3:A:466:LEU:CD2	3:A:470:HIS:CE1	2.88	0.56
4:B:359:LEU:HD12	4:B:370:LYS:NZ	2.21	0.56
4:B:1157:GLN:HE22	4:B:1171:ASN:HB2	1.71	0.56
15:M:230:LYS:HD2	15:M:247:LEU:HD21	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:238:MET:SD	3:A:238:MET:N	2.78	0.56
4:B:492:ASN:ND2	4:B:722:GLY:O	2.39	0.56
3:A:629:ASP:O	4:B:924:LYS:NZ	2.38	0.56
7:E:135:PHE:CE1	7:E:186:LEU:CD2	2.89	0.56
15:M:336:ILE:HG21	15:M:349:LEU:HD11	1.88	0.56
2:U:61:DA:H4'	3:A:1601:GLN:NE2	2.20	0.56
3:A:673:HIS:HD2	3:A:817:PHE:HB2	1.71	0.56
4:B:623:ASP:HA	4:B:663:ILE:HD11	1.88	0.56
16:N:114:GLU:OE1	16:N:116:LYS:NZ	2.37	0.56
3:A:400:ASN:ND2	15:M:169:SER:OG	2.37	0.56
3:A:483:VAL:O	3:A:613:THR:OG1	2.23	0.56
3:A:818:THR:HA	3:A:821:ILE:HG22	1.88	0.56
3:A:325:ASP:OD1	3:A:329:ARG:NH2	2.39	0.55
3:A:1559:ARG:HH22	7:E:200:ARG:HG3	1.70	0.55
4:B:440:PHE:HA	4:B:445:TYR:HB3	1.88	0.55
4:B:1076:ARG:HH12	4:B:1088:LEU:HD22	1.70	0.55
5:C:100:ARG:HH12	12:J:2:ILE:HB	1.71	0.55
5:C:218:LYS:NZ	14:L:70:ARG:OXT	2.40	0.55
9:G:90:LEU:HD23	9:G:118:CYS:HA	1.88	0.55
3:A:942:GLN:O	3:A:985:ARG:NH1	2.38	0.55
5:C:90:SER:HA	5:C:200:GLN:HG2	1.88	0.55
15:M:268:LEU:HD23	15:M:268:LEU:O	2.06	0.55
3:A:1256:LYS:HB2	3:A:1307:ASP:OD2	2.05	0.55
4:B:703:LEU:HD13	4:B:704:THR:CA	2.36	0.55
3:A:729:LYS:HE2	10:H:120:GLY:HA3	1.87	0.55
4:B:652:PRO:HA	4:B:663:ILE:HA	1.88	0.55
15:M:262:LEU:O	15:M:265:LEU:HB2	2.05	0.55
3:A:1634:LEU:HD12	3:A:1634:LEU:O	2.06	0.55
9:G:161:ASN:HD22	9:G:248:THR:HG22	1.71	0.55
6:D:84:SER:O	6:D:88:GLN:NE2	2.40	0.55
14:L:31:CYS:SG	14:L:32:ALA:N	2.79	0.55
3:A:194:ALA:HA	3:A:197:LEU:HB2	1.88	0.55
4:B:97:VAL:HG21	4:B:424:ILE:HD11	1.88	0.55
9:G:218:VAL:HG13	9:G:222:GLY:HA2	1.89	0.55
3:A:250:LYS:HD2	3:A:316:LEU:HA	1.88	0.55
3:A:1450:ILE:O	3:A:1454:HIS:N	2.40	0.55
4:B:912:GLN:OE1	4:B:1041:ASN:ND2	2.39	0.55
5:C:116:VAL:HG21	5:C:209:ILE:HG13	1.87	0.55
9:G:37:CYS:SG	9:G:125:TRP:NE1	2.78	0.55
9:G:140:GLN:HE21	9:G:215:GLY:H	1.52	0.55
3:A:532:GLY:H	3:A:580:HIS:CE1	2.25	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1005:TYR:CE1	16:N:171:PHE:CD2	2.95	0.55
7:E:5:ASN:OD1	7:E:52:ARG:NH2	2.40	0.55
3:A:559:ASN:ND2	17:O:376:TYR:O	2.40	0.54
3:A:1053:ASP:OD1	3:A:1174:TYR:OH	2.24	0.54
4:B:444:ARG:HG3	4:B:448:ARG:HH11	1.72	0.54
4:B:673:ASN:ND2	4:B:685:VAL:O	2.40	0.54
4:B:766:PRO:HG3	12:J:54:VAL:HG21	1.89	0.54
4:B:997:GLY:O	4:B:1001:ALA:N	2.41	0.54
14:L:48:CYS:N	14:L:53:HIS:O	2.39	0.54
15:M:267:LEU:HG	15:M:270:GLY:HA3	1.88	0.54
3:A:15:ASP:HB3	4:B:1197:ARG:HB3	1.89	0.54
3:A:1047:GLN:HE21	3:A:1587:ASP:HB2	1.72	0.54
4:B:359:LEU:CD1	4:B:370:LYS:NZ	2.70	0.54
15:M:261:LEU:O	15:M:265:LEU:HG	2.07	0.54
3:A:438:ILE:HG22	3:A:457:LYS:HE2	1.89	0.54
3:A:1591:ARG:HE	3:A:1592:GLN:HE21	1.55	0.54
5:C:54:PHE:HB3	5:C:300:PHE:HB2	1.89	0.54
16:N:82:ILE:HG13	16:N:85:HIS:H	1.73	0.54
3:A:1256:LYS:O	3:A:1499:ARG:NH2	2.36	0.54
3:A:1306:TYR:C	3:A:1307:ASP:OD1	2.45	0.54
4:B:857:PRO:HB3	4:B:871:ILE:HG23	1.89	0.54
7:E:135:PHE:CE1	7:E:186:LEU:HD21	2.42	0.54
14:L:38:LEU:HD13	14:L:40:LEU:HD13	1.89	0.54
15:M:79:GLY:N	16:N:54:TRP:O	2.36	0.54
3:A:196:ALA:HB2	3:A:201:ARG:HH21	1.73	0.54
3:A:883:LEU:HD12	3:A:884:ARG:HE	1.73	0.54
4:B:974:LEU:HD11	16:N:169:GLU:HB3	1.89	0.54
5:C:128:ASP:OD2	5:C:174:ARG:N	2.40	0.54
17:O:382:GLN:NE2	17:O:593:PRO:O	2.40	0.54
3:A:516:ILE:HD11	17:O:588:LEU:HD11	1.89	0.54
3:A:592:GLN:NE2	3:A:634:ASN:OD1	2.35	0.54
4:B:1000:LEU:HA	4:B:1003:ALA:HB3	1.90	0.54
9:G:26:ASN:OD1	9:G:128:GLN:NE2	2.40	0.54
1:T:13:DT:O4	2:U:56:DA:N1	2.40	0.54
4:B:280:LEU:HG	4:B:370:LYS:HB3	1.89	0.54
4:B:317:TYR:HD2	4:B:320:LEU:HB2	1.72	0.54
4:B:543:ASN:OD1	4:B:543:ASN:N	2.41	0.54
6:D:94:ARG:HA	6:D:98:GLY:H	1.73	0.54
11:I:2:SER:O	11:I:9:PHE:N	2.40	0.54
1:T:12:DC:H3'	1:T:12:DC:H6	1.72	0.54
4:B:555:GLN:OE1	4:B:646:HIS:ND1	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:588:ILE:HA	4:B:642:LEU:HB2	1.89	0.54
4:B:650:LEU:HD12	4:B:663:ILE:HB	1.89	0.54
6:D:94:ARG:HD3	6:D:100:PRO:HD3	1.90	0.54
15:M:299:LEU:O	15:M:303:THR:OG1	2.26	0.54
3:A:499:PRO:HA	3:A:502:ALA:HB3	1.89	0.53
3:A:985:ARG:HD3	3:A:987:TYR:N	2.23	0.53
15:M:35:ASP:OD1	15:M:35:ASP:N	2.40	0.53
3:A:668:GLY:HA3	3:A:787:GLY:HA2	1.89	0.53
3:A:686:PHE:HB3	3:A:725:LEU:HD11	1.90	0.53
3:A:1055:ILE:O	3:A:1580:ARG:NH1	2.41	0.53
6:D:84:SER:OG	17:O:228:GLN:NE2	2.41	0.53
3:A:1046:VAL:HG22	3:A:1591:ARG:HD3	1.91	0.53
4:B:129:ARG:NH2	4:B:891:GLU:H	2.06	0.53
4:B:876:SER:OG	4:B:878:GLU:OE1	2.25	0.53
10:H:118:PHE:HB2	10:H:121:LEU:HB3	1.89	0.53
15:M:20:SER:OG	15:M:91:TYR:OH	2.26	0.53
15:M:261:LEU:HD13	15:M:335:ILE:HG12	1.90	0.53
3:A:480:ALA:HB3	3:A:635:MET:HB3	1.91	0.53
4:B:265:ARG:O	4:B:266:LYS:HG2	2.09	0.53
5:C:303:GLU:OE2	12:J:43:ARG:NH1	2.41	0.53
3:A:832:ASP:OD1	3:A:923:ASN:ND2	2.41	0.53
3:A:1562:ILE:HD11	3:A:1586:ALA:HA	1.89	0.53
4:B:221:SER:HA	4:B:224:ASN:HB2	1.90	0.53
4:B:906:ARG:O	4:B:908:ARG:NH2	2.39	0.53
2:U:65:DG:O5'	2:U:65:DG:H8	1.92	0.53
3:A:1528:ALA:O	3:A:1532:GLN:NE2	2.41	0.53
4:B:690:GLU:OE2	4:B:695:ASN:ND2	2.42	0.53
5:C:39:ASP:O	5:C:58:ASN:ND2	2.39	0.53
9:G:24:VAL:O	9:G:128:GLN:NE2	2.42	0.53
11:I:58:SER:OG	11:I:61:ARG:N	2.40	0.53
3:A:815:ARG:O	3:A:819:ASN:ND2	2.41	0.53
13:K:60:SER:OG	13:K:104:ARG:NH2	2.42	0.53
3:A:1636:SER:O	3:A:1640:ARG:N	2.42	0.53
3:A:113:VAL:HG21	3:A:178:LEU:HD12	1.89	0.53
4:B:359:LEU:CD1	4:B:370:LYS:HG3	2.27	0.53
12:J:34:THR:HA	12:J:37:SER:HB2	1.91	0.53
4:B:908:ARG:NH1	5:C:93:GLN:OE1	2.41	0.52
10:H:28:ALA:HB3	10:H:38:LEU:HB3	1.91	0.52
13:K:79:VAL:HG23	13:K:82:LYS:HD2	1.91	0.52
15:M:360:VAL:HA	15:M:363:LEU:HB3	1.91	0.52
3:A:1439:MET:SD	3:A:1439:MET:N	2.82	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:95:LEU:HA	4:B:145:VAL:HA	1.90	0.52
4:B:748:GLN:HB2	4:B:769:PHE:HA	1.91	0.52
7:E:85:GLU:O	7:E:113:GLN:NE2	2.42	0.52
1:T:19:DA:OP1	4:B:1063:ARG:NH1	2.42	0.52
3:A:830:MET:O	3:A:834:ARG:NH2	2.43	0.52
4:B:256:GLY:HA3	4:B:308:LEU:HD22	1.92	0.52
4:B:586:VAL:N	4:B:594:GLY:O	2.41	0.52
4:B:589:ASP:OD1	4:B:589:ASP:N	2.42	0.52
6:D:88:GLN:HG3	6:D:89:LEU:HD12	1.90	0.52
3:A:646:GLU:O	3:A:650:LEU:N	2.42	0.52
3:A:721:LYS:HE2	10:H:46:LEU:HD22	1.89	0.52
3:A:1097:TYR:OH	3:A:1121:ASP:OD1	2.27	0.52
4:B:127:ARG:HB3	4:B:195:ILE:HD13	1.91	0.52
3:A:41:LEU:HD12	3:A:43:HIS:CE1	2.38	0.52
4:B:167:SER:N	4:B:170:CYS:SG	2.76	0.52
4:B:244:THR:HG21	4:B:414:LYS:HD2	1.91	0.52
4:B:532:HIS:ND1	4:B:719:CYS:SG	2.70	0.52
4:B:1111:LEU:HD13	4:B:1187:SER:HB2	1.92	0.52
9:G:63:LYS:HD3	9:G:67:ASN:HD22	1.74	0.52
3:A:1038:ILE:HD11	3:A:1047:GLN:HB2	1.92	0.52
4:B:60:LEU:HD12	4:B:242:ASP:HA	1.92	0.52
8:F:101:ILE:HA	8:F:105:ALA:HB3	1.91	0.52
10:H:43:ASN:ND2	10:H:45:GLU:OE1	2.43	0.52
15:M:250:LEU:HG	15:M:250:LEU:O	2.10	0.52
17:O:480:LEU:HD12	17:O:483:ILE:HD12	1.92	0.52
3:A:985:ARG:HH21	3:A:986:PHE:N	2.04	0.52
4:B:219:ARG:HG2	4:B:221:SER:H	1.75	0.52
3:A:216:ARG:NH2	3:A:340:HIS:O	2.43	0.52
3:A:435:ASN:OD1	3:A:436:ALA:N	2.42	0.52
4:B:889:GLY:HA3	4:B:898:LEU:HD22	1.91	0.52
4:B:820:PRO:HD2	15:M:356:LYS:HZ1	1.73	0.52
4:B:228:SER:O	4:B:254:ASN:ND2	2.42	0.52
3:A:99:ARG:O	3:A:109:ARG:NH2	2.38	0.51
3:A:252:PHE:HZ	15:M:144:LEU:HG	1.74	0.51
3:A:462:LYS:NZ	3:A:465:GLY:O	2.42	0.51
3:A:470:HIS:O	3:A:474:LYS:NZ	2.43	0.51
4:B:837:LEU:O	4:B:847:TYR:OH	2.26	0.51
5:C:301:ASN:HD21	16:N:174:GLY:HA2	1.75	0.51
9:G:73:TYR:HB2	9:G:80:VAL:HG22	1.92	0.51
17:O:163:ILE:HA	17:O:211:TYR:HB2	1.90	0.51
3:A:52:LEU:HD21	3:A:60:ASN:HD22	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1333:ILE:HG12	3:A:1480:THR:HG21	1.92	0.51
4:B:784:ASP:OD1	4:B:785:ASP:N	2.43	0.51
10:H:40:LEU:HD21	10:H:42:ILE:HD13	1.92	0.51
17:O:428:ILE:HA	17:O:434:LEU:HD21	1.92	0.51
3:A:898:SER:O	3:A:902:ALA:N	2.43	0.51
4:B:576:THR:HB	4:B:595:TRP:HZ2	1.75	0.51
15:M:113:ILE:HG23	15:M:117:SER:HA	1.92	0.51
16:N:58:PHE:HB3	16:N:139:VAL:HB	1.91	0.51
3:A:908:VAL:HG23	3:A:941:SER:HB3	1.92	0.51
3:A:1042:ASP:OD1	3:A:1042:ASP:N	2.37	0.51
3:A:1588:MET:SD	3:A:1591:ARG:NH1	2.82	0.51
5:C:45:SER:CB	5:C:53:ASN:HB3	2.35	0.51
15:M:250:LEU:N	15:M:250:LEU:CD2	2.73	0.51
17:O:412:GLU:OE1	17:O:416:LYS:NZ	2.35	0.51
3:A:117:ARG:O	3:A:121:LYS:N	2.40	0.51
3:A:1072:ASN:HB3	3:A:1075:ALA:HB3	1.92	0.51
4:B:431:ASP:O	4:B:435:GLY:N	2.43	0.51
7:E:23:VAL:HG23	7:E:28:TYR:HD2	1.76	0.51
13:K:88:PHE:O	13:K:106:GLN:N	2.41	0.51
3:A:1053:ASP:O	3:A:1174:TYR:OH	2.29	0.51
3:A:1484:LEU:O	3:A:1488:ILE:N	2.42	0.51
4:B:238:SER:OG	4:B:246:GLN:NE2	2.44	0.51
4:B:327:LEU:HD22	4:B:350:GLY:HA3	1.93	0.51
4:B:995:TYR:OH	16:N:161:PRO:O	2.26	0.51
7:E:147:HIS:HB3	7:E:150:VAL:HG22	1.93	0.51
16:N:50:GLN:HE22	16:N:52:GLN:HB3	1.76	0.51
3:A:314:TYR:OH	15:M:146:ARG:O	2.26	0.51
3:A:463:LYS:HG3	3:A:468:ARG:HH12	1.75	0.51
3:A:1237:GLN:HE22	3:A:1520:VAL:HB	1.76	0.51
10:H:26:ILE:HB	10:H:40:LEU:HB3	1.92	0.51
3:A:41:LEU:HD13	3:A:41:LEU:C	2.31	0.51
3:A:214:ASP:O	3:A:218:LYS:N	2.42	0.51
4:B:346:ASP:HA	4:B:349:VAL:HG12	1.93	0.51
4:B:420:TYR:HE1	4:B:455:GLU:HG2	1.74	0.51
4:B:714:ARG:CD	4:B:957:ARG:HD2	2.38	0.51
3:A:613:THR:O	3:A:615:ARG:NH2	2.44	0.51
3:A:673:HIS:CD2	3:A:817:PHE:HB2	2.46	0.51
3:A:842:TRP:HA	3:A:845:ASP:HB2	1.93	0.51
4:B:863:ASP:OD1	4:B:870:LYS:NZ	2.36	0.51
5:C:320:ILE:O	5:C:324:LYS:N	2.44	0.51
3:A:912:VAL:CB	3:A:913:PRO:HD2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:656:LEU:HD11	4:B:689:VAL:HG13	1.92	0.50
4:B:711:GLN:HB3	4:B:713:PRO:HD2	1.93	0.50
8:F:97:ARG:O	8:F:101:ILE:N	2.35	0.50
9:G:17:ILE:O	9:G:21:LYS:N	2.43	0.50
17:O:139:PHE:HA	17:O:142:ILE:HG22	1.93	0.50
3:A:368:ARG:HH11	3:A:383:ASN:HD21	1.59	0.50
15:M:350:ALA:O	15:M:355:LEU:O	2.29	0.50
3:A:843:ARG:HH22	3:A:985:ARG:HB3	1.77	0.50
3:A:937:ASN:HA	3:A:940:VAL:HG22	1.93	0.50
3:A:1054:ALA:O	3:A:1178:LEU:HD22	2.11	0.50
3:A:1103:LYS:HA	3:A:1106:LYS:HG2	1.94	0.50
17:O:67:ASP:OD2	17:O:69:THR:OG1	2.29	0.50
3:A:368:ARG:HD2	3:A:383:ASN:HD21	1.76	0.50
5:C:57:ILE:HG23	5:C:297:HIS:HA	1.92	0.50
13:K:63:PHE:N	13:K:103:ILE:O	2.44	0.50
3:A:481:ARG:HB2	4:B:1069:ILE:HD12	1.92	0.50
3:A:1610:PHE:HB2	3:A:1639:ALA:HB2	1.94	0.50
4:B:651:ARG:N	4:B:664:VAL:O	2.44	0.50
4:B:1112:THR:HG21	4:B:1130:ARG:HB2	1.94	0.50
5:C:100:ARG:NH2	12:J:3:VAL:O	2.45	0.50
15:M:42:LYS:HD2	16:N:32:CYS:HA	1.93	0.50
3:A:1091:VAL:HA	3:A:1133:LEU:HB2	1.92	0.50
3:A:1549:VAL:HG11	3:A:1561:THR:HG21	1.92	0.50
4:B:785:ASP:N	4:B:785:ASP:OD1	2.44	0.50
4:B:898:LEU:HD11	14:L:46:VAL:HG11	1.93	0.50
12:J:1:MET:HA	12:J:56:LEU:H	1.76	0.50
4:B:322:ASN:ND2	15:M:104:SER:O	2.45	0.50
4:B:806:THR:HA	4:B:904:LYS:HA	1.93	0.50
6:D:99:LEU:HD22	6:D:100:PRO:HD2	1.94	0.50
9:G:137:ILE:HB	9:G:227:GLY:HA2	1.94	0.50
3:A:836:THR:O	3:A:840:ASN:N	2.37	0.50
3:A:1270:VAL:HG11	3:A:1489:VAL:HG11	1.93	0.50
3:A:1634:LEU:HD13	3:A:1639:ALA:O	2.11	0.50
7:E:8:ASN:HA	7:E:11:ARG:HG2	1.93	0.50
4:B:334:PHE:HA	4:B:337:VAL:HG12	1.93	0.50
13:K:83:ASN:HD22	13:K:84:PRO:HD2	1.77	0.50
3:A:237:GLY:O	3:A:263:ASN:ND2	2.44	0.49
3:A:732:ILE:O	3:A:736:LEU:HB2	2.12	0.49
3:A:843:ARG:NH2	3:A:985:ARG:HB3	2.26	0.49
3:A:872:ASP:HB2	3:A:875:LEU:CB	2.42	0.49
3:A:1031:HIS:HB3	3:A:1649:VAL:HG13	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:529:CYS:HB3	4:B:532:HIS:H	1.77	0.49
5:C:233:ILE:HD12	5:C:291:LEU:HB3	1.94	0.49
16:N:73:ASP:HB2	16:N:77:SER:HB3	1.93	0.49
3:A:125:LEU:N	3:A:126:GLN:OE1	2.40	0.49
4:B:335:ARG:HH22	15:M:112:LYS:HD2	1.77	0.49
4:B:726:MET:SD	4:B:1035:ARG:NH1	2.85	0.49
5:C:197:ARG:HG2	12:J:61:LEU:HD22	1.94	0.49
17:O:438:GLN:HA	17:O:441:PHE:HB3	1.94	0.49
3:A:809:VAL:HA	3:A:812:VAL:HG22	1.95	0.49
3:A:947:LEU:HD11	3:A:950:GLN:HE21	1.77	0.49
3:A:1651:THR:HG22	4:B:1085:SER:H	1.78	0.49
4:B:1088:LEU:HA	4:B:1091:ARG:HB3	1.95	0.49
9:G:64:GLN:HG3	9:G:65:HIS:CD2	2.48	0.49
3:A:506:THR:O	3:A:639:GLN:NE2	2.33	0.49
3:A:1032:VAL:O	3:A:1182:GLY:N	2.40	0.49
7:E:80:VAL:HG22	7:E:109:ILE:HB	1.95	0.49
12:J:16:ASP:OD1	12:J:16:ASP:N	2.42	0.49
15:M:58:GLU:OE2	15:M:59:ARG:NH1	2.46	0.49
3:A:1437:ASN:HD22	3:A:1438:ASN:H	1.60	0.49
7:E:95:THR:HA	7:E:98:ILE:HD12	1.94	0.49
12:J:45:CYS:O	12:J:48:ARG:NE	2.39	0.49
3:A:715:LEU:HD22	3:A:730:GLN:HE22	1.77	0.49
3:A:975:ASP:N	3:A:975:ASP:OD1	2.44	0.49
3:A:1256:LYS:HB3	3:A:1307:ASP:OD2	2.12	0.49
4:B:286:ARG:O	4:B:290:ASP:N	2.35	0.49
4:B:572:PRO:HB2	4:B:575:HIS:HD2	1.77	0.49
5:C:65:ASN:HD22	5:C:68:ARG:HE	1.61	0.49
15:M:321:ASP:OD1	15:M:324:ASN:ND2	2.46	0.49
3:A:466:LEU:HD23	3:A:470:HIS:CE1	2.48	0.49
3:A:972:TYR:OH	4:B:633:THR:OG1	2.29	0.49
8:F:65:ARG:HH12	9:G:117:TRP:HE3	1.61	0.49
3:A:24:ILE:HA	3:A:27:LEU:HD22	1.95	0.49
3:A:720:PHE:HB2	10:H:96:VAL:HB	1.95	0.49
3:A:1634:LEU:CD1	3:A:1639:ALA:O	2.61	0.49
4:B:947:ILE:HG21	4:B:1033:TYR:HE2	1.77	0.49
11:I:58:SER:HB2	11:I:60:LEU:HD23	1.95	0.49
13:K:92:SER:OG	13:K:102:ASN:O	2.31	0.49
16:N:57:LYS:HB3	16:N:138:SER:HA	1.93	0.49
17:O:81:PRO:HB2	17:O:83:LYS:HG2	1.94	0.49
4:B:974:LEU:CD1	16:N:169:GLU:CG	2.84	0.48
17:O:241:ASP:OD2	17:O:380:SER:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:O:445:TYR:O	17:O:449:TRP:N	2.46	0.48
3:A:1648:ASN:OD1	3:A:1648:ASN:N	2.46	0.48
4:B:152:LEU:HD23	4:B:443:LYS:HG3	1.95	0.48
4:B:653:VAL:N	4:B:662:ASP:O	2.40	0.48
17:O:198:PHE:HB2	17:O:232:LEU:HD21	1.95	0.48
4:B:556:SER:HB2	4:B:621:PRO:HG3	1.95	0.48
4:B:1116:SER:OG	4:B:1160:GLU:O	2.31	0.48
8:F:72:LYS:HE3	8:F:141:GLY:HA2	1.94	0.48
3:A:110:LEU:HA	3:A:227:LEU:HD13	1.95	0.48
4:B:218:ILE:HD12	4:B:231:HIS:HD2	1.78	0.48
4:B:756:LEU:HD13	4:B:759:ASP:HB3	1.95	0.48
4:B:784:ASP:C	4:B:785:ASP:OD1	2.52	0.48
7:E:143:ASN:HD22	7:E:146:HIS:CE1	2.31	0.48
3:A:862:THR:HG23	3:A:864:LEU:H	1.78	0.48
4:B:27:ASN:N	4:B:27:ASN:OD1	2.47	0.48
4:B:969:GLY:HA3	4:B:1030:VAL:HG23	1.94	0.48
4:B:204:ARG:HD2	4:B:486:VAL:HG13	1.95	0.48
4:B:283:THR:OG1	4:B:323:ARG:NH2	2.47	0.48
5:C:200:GLN:NE2	12:J:64:ASN:OD1	2.42	0.48
17:O:101:SER:HA	17:O:104:ILE:HD13	1.95	0.48
3:A:224:HIS:HA	3:A:228:LEU:HD22	1.94	0.48
3:A:399:LEU:HD23	3:A:422:ARG:HD3	1.95	0.48
4:B:742:TYR:HE2	4:B:1037:ARG:HG3	1.79	0.48
6:D:19:PRO:HD3	9:G:65:HIS:HB3	1.96	0.48
3:A:381:SER:HB2	3:A:453:ILE:HD11	1.94	0.48
3:A:1550:LEU:HA	3:A:1554:GLY:H	1.79	0.48
3:A:1578:SER:OG	3:A:1579:PHE:N	2.46	0.48
15:M:22:ALA:HB3	16:N:110:LEU:HD21	1.95	0.48
3:A:912:VAL:CB	3:A:913:PRO:CD	2.92	0.48
4:B:126:SER:OG	4:B:132:SER:O	2.31	0.48
4:B:931:TRP:CD2	4:B:932:PRO:HD2	2.49	0.48
3:A:911:CYS:O	3:A:911:CYS:SG	2.72	0.48
3:A:1032:VAL:HB	3:A:1181:PRO:HA	1.95	0.48
4:B:564:ILE:HD11	4:B:620:LEU:HD21	1.96	0.48
7:E:12:LEU:O	7:E:16:PHE:N	2.41	0.48
8:F:87:LYS:O	8:F:91:ALA:N	2.45	0.48
3:A:736:LEU:HD22	3:A:736:LEU:HA	1.60	0.47
3:A:1657:LEU:HB2	8:F:133:VAL:HB	1.96	0.47
3:A:52:LEU:HD11	3:A:60:ASN:HB2	1.95	0.47
3:A:885:ASP:N	3:A:885:ASP:OD1	2.42	0.47
15:M:78:VAL:HA	16:N:55:LEU:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1097:ASP:OD1	4:B:1180:PHE:N	2.43	0.47
4:B:1100:GLN:HB3	4:B:1175:THR:HB	1.96	0.47
4:B:1116:SER:O	4:B:1125:THR:OG1	2.30	0.47
8:F:130:ILE:HG23	8:F:148:VAL:HG11	1.95	0.47
3:A:33:THR:HG22	3:A:394:LEU:HD11	1.96	0.47
3:A:41:LEU:CD1	15:M:323:GLN:HA	2.43	0.47
3:A:41:LEU:CD1	15:M:322:PRO:O	2.62	0.47
3:A:96:ILE:HG21	3:A:224:HIS:HD2	1.78	0.47
3:A:726:TRP:HE3	3:A:730:GLN:HG2	1.80	0.47
3:A:781:LEU:HD21	3:A:786:TYR:HE1	1.80	0.47
3:A:1027:LEU:HD13	3:A:1186:GLY:HA2	1.97	0.47
3:A:1531:ASP:OD1	3:A:1532:GLN:NE2	2.47	0.47
4:B:830:ASP:N	4:B:830:ASP:OD1	2.43	0.47
10:H:28:ALA:O	10:H:38:LEU:N	2.48	0.47
17:O:478:GLN:O	17:O:482:TYR:N	2.47	0.47
3:A:400:ASN:HD22	15:M:167:THR:HG23	1.78	0.47
3:A:672:ASP:OD1	3:A:673:HIS:ND1	2.48	0.47
3:A:1044:THR:HA	7:E:174:GLN:HE21	1.79	0.47
4:B:52:LEU:HA	4:B:60:LEU:HD23	1.96	0.47
4:B:939:SER:OG	4:B:1011:GLU:OE2	2.26	0.47
4:B:1015:SER:HG	4:B:1020:GLU:H	1.63	0.47
5:C:73:SER:OG	5:C:74:GLU:OE1	2.28	0.47
17:O:344:GLU:HG3	17:O:388:VAL:HG23	1.96	0.47
3:A:14:VAL:O	3:A:1631:ARG:NH1	2.48	0.47
4:B:104:ILE:HG21	4:B:161:LEU:HB3	1.97	0.47
6:D:39:PHE:HD2	9:G:40:ARG:HH22	1.60	0.47
15:M:153:LYS:HB2	15:M:156:ASP:HB2	1.96	0.47
15:M:305:ILE:HG21	15:M:316:ARG:HH21	1.80	0.47
3:A:618:TYR:CD1	4:B:783:MET:HE3	2.42	0.47
3:A:618:TYR:HA	4:B:783:MET:CE	2.44	0.47
3:A:1178:LEU:CD1	3:A:1179:ILE:O	2.63	0.47
4:B:359:LEU:CG	4:B:370:LYS:HE3	2.44	0.47
4:B:623:ASP:OD1	4:B:623:ASP:N	2.35	0.47
4:B:707:SER:H	4:B:983:PRO:HG3	1.79	0.47
4:B:975:HIS:CE1	16:N:166:LEU:HB3	2.50	0.47
5:C:107:LYS:HE3	5:C:187:ALA:HA	1.96	0.47
3:A:23:GLU:OE1	4:B:1130:ARG:NH1	2.47	0.47
3:A:985:ARG:HD2	3:A:987:TYR:N	2.26	0.47
3:A:1057:ILE:HG13	3:A:1581:HIS:HE1	1.80	0.47
4:B:825:PHE:HA	4:B:861:TYR:HB2	1.97	0.47
5:C:93:GLN:HE22	5:C:95:GLU:HB3	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:E:151:PRO:HB3	7:E:200:ARG:HA	1.95	0.47
10:H:137:GLN:OE1	10:H:139:ASN:N	2.48	0.47
15:M:8:SER:OG	15:M:9:GLU:N	2.46	0.47
17:O:238:ILE:HG23	17:O:378:THR:HG21	1.97	0.47
3:A:43:HIS:HA	15:M:322:PRO:HG2	1.97	0.47
3:A:64:THR:HA	4:B:1162:GLY:HA3	1.96	0.47
3:A:759:TYR:CD1	3:A:913:PRO:HG2	2.50	0.47
4:B:784:ASP:C	4:B:950:ASN:HD22	2.18	0.47
6:D:36:VAL:HG22	6:D:40:LEU:HG	1.96	0.47
14:L:32:ALA:HB3	14:L:53:HIS:HE1	1.80	0.47
15:M:306:LYS:HG3	15:M:319:PHE:HB2	1.96	0.47
3:A:836:THR:OG1	3:A:839:GLY:N	2.46	0.47
4:B:395:ASP:OD1	4:B:395:ASP:N	2.39	0.47
4:B:724:GLN:HE21	4:B:1037:ARG:HB3	1.80	0.47
4:B:790:ASN:HB3	4:B:793:ALA:HB3	1.96	0.47
5:C:272:LYS:HG2	16:N:175:TYR:CZ	2.50	0.47
3:A:942:GLN:HG2	3:A:947:LEU:HB2	1.96	0.46
5:C:42:VAL:HB	13:K:138:LYS:HG3	1.98	0.46
9:G:80:VAL:N	9:G:125:TRP:O	2.48	0.46
4:B:20:GLU:O	4:B:24:ARG:NE	2.47	0.46
4:B:939:SER:O	5:C:226:SER:OG	2.33	0.46
10:H:101:ALA:HB3	10:H:137:GLN:H	1.79	0.46
12:J:48:ARG:O	12:J:52:THR:N	2.47	0.46
17:O:420:SER:HA	17:O:423:TYR:HD2	1.79	0.46
3:A:561:LEU:HA	3:A:575:LYS:HD2	1.96	0.46
4:B:1113:THR:OG1	4:B:1127:CYS:O	2.28	0.46
5:C:181:ASP:OD1	5:C:181:ASP:N	2.46	0.46
3:A:949:GLN:HG2	3:A:981:TYR:HA	1.97	0.46
4:B:258:VAL:HG11	4:B:378:ILE:HD11	1.98	0.46
4:B:480:GLN:OE1	4:B:508:PHE:N	2.46	0.46
4:B:963:PHE:O	4:B:1027:TYR:OH	2.24	0.46
7:E:39:LEU:O	7:E:43:LYS:N	2.34	0.46
15:M:42:LYS:HA	15:M:51:PHE:HD1	1.79	0.46
16:N:140:SER:OG	16:N:141:GLU:N	2.48	0.46
3:A:61:LEU:HG	15:M:306:LYS:HG2	1.97	0.46
4:B:1047:ARG:HH12	4:B:1060:VAL:H	1.63	0.46
7:E:93:MET:HG3	7:E:120:ALA:HB1	1.96	0.46
9:G:138:PHE:HE2	17:O:183:ILE:HG12	1.81	0.46
16:N:39:PRO:HB2	16:N:51:GLN:HE22	1.81	0.46
3:A:127:TYR:HB2	3:A:129:LEU:HD23	1.98	0.46
3:A:1114:TYR:HB2	7:E:146:HIS:HD2	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:36:PHE:HE2	13:K:130:VAL:HG11	1.81	0.46
15:M:231:LEU:HD11	15:M:244:ALA:HB2	1.96	0.46
15:M:363:LEU:HD13	15:M:363:LEU:C	2.35	0.46
15:M:388:THR:OG1	15:M:391:SER:OG	2.33	0.46
3:A:267:LYS:HB2	3:A:270:ILE:HD11	1.96	0.46
3:A:514:TYR:CZ	8:F:120:ILE:HD11	2.50	0.46
3:A:759:TYR:CZ	3:A:913:PRO:HG2	2.51	0.46
4:B:629:VAL:HG13	4:B:639:GLY:H	1.81	0.46
4:B:730:GLY:HA2	4:B:765:PHE:HE1	1.80	0.46
16:N:33:LYS:O	16:N:115:SER:OG	2.33	0.46
4:B:850:THR:H	4:B:882:ILE:HG13	1.81	0.46
4:B:881:TYR:HB2	4:B:908:ARG:NH2	2.31	0.46
10:H:5:LEU:HD12	10:H:133:ASN:HB2	1.96	0.46
15:M:9:GLU:HA	16:N:71:PRO:HA	1.97	0.46
3:A:489:ASN:ND2	13:K:94:PRO:O	2.49	0.46
3:A:1337:LYS:HZ1	3:A:1480:THR:HA	1.81	0.46
4:B:71:LYS:HZ3	4:B:421:LEU:HD12	1.79	0.46
4:B:96:SER:OG	4:B:97:VAL:N	2.49	0.46
4:B:532:HIS:HE1	4:B:723:LYS:HD3	1.81	0.46
4:B:586:VAL:HG22	4:B:640:LEU:HB3	1.97	0.46
4:B:1015:SER:OG	4:B:1020:GLU:N	2.49	0.46
4:B:1103:VAL:HG22	4:B:1110:ILE:HG13	1.97	0.46
17:O:401:VAL:HA	17:O:404:ILE:HD12	1.97	0.46
2:U:65:DG:H2'	2:U:66:DA:C8	2.50	0.46
3:A:14:VAL:HG12	3:A:1632:GLU:HB2	1.97	0.46
3:A:252:PHE:HB2	3:A:314:TYR:HA	1.98	0.45
3:A:883:LEU:HA	3:A:889:SER:HB2	1.97	0.45
3:A:1178:LEU:HD13	3:A:1179:ILE:O	2.16	0.45
4:B:977:ILE:H	4:B:977:ILE:HG13	1.52	0.45
4:B:970:LYS:NZ	4:B:1005:TYR:CD1	2.76	0.45
13:K:88:PHE:HB3	13:K:106:GLN:HB2	1.98	0.45
3:A:1085:LEU:HD12	3:A:1172:LEU:HD11	1.98	0.45
4:B:1003:ALA:CB	4:B:1005:TYR:HD2	2.29	0.45
5:C:252:PRO:HA	5:C:253:PRO:HD3	1.84	0.45
15:M:82:ASN:HB2	15:M:85:LYS:HB3	1.99	0.45
16:N:169:GLU:OE2	16:N:170:HIS:C	2.54	0.45
3:A:56:ALA:HB2	3:A:62:CYS:HB2	1.97	0.45
3:A:602:GLY:N	3:A:651:ALA:O	2.49	0.45
4:B:359:LEU:HG	4:B:370:LYS:HE3	1.99	0.45
4:B:974:LEU:HD23	12:J:44:TYR:HB3	1.99	0.45
5:C:53:ASN:HD21	5:C:271:ARG:NH1	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:7:DA:H2"	1:T:8:DA:C8	2.51	0.45
4:B:1039:MET:HG3	4:B:1043:LYS:HG2	1.98	0.45
4:B:1119:ARG:H	4:B:1122:SER:HB2	1.81	0.45
15:M:235:PRO:HG2	15:M:263:TYR:HE1	1.81	0.45
3:A:538:ASN:HA	3:A:575:LYS:HG2	1.98	0.45
4:B:644:GLY:HA2	4:B:648:ARG:HG3	1.99	0.45
9:G:228:LYS:HE2	9:G:230:ARG:HH12	1.82	0.45
10:H:38:LEU:HD12	10:H:125:LEU:HD21	1.97	0.45
13:K:102:ASN:N	13:K:102:ASN:OD1	2.48	0.45
3:A:1197:SER:HA	3:A:1200:MET:HG2	1.99	0.45
9:G:62:MET:HG3	9:G:66:LEU:HD23	1.99	0.45
17:O:540:CYS:HA	17:O:543:ILE:HD12	1.98	0.45
3:A:397:ARG:HD3	15:M:172:LEU:HB2	1.99	0.45
3:A:757:ASN:HD21	3:A:767:ASN:H	1.63	0.45
3:A:1221:ARG:NH2	3:A:1544:ASN:OD1	2.49	0.45
4:B:350:GLY:HA2	4:B:353:VAL:HG22	1.99	0.45
4:B:419:GLU:HA	4:B:422:GLN:HB2	1.97	0.45
4:B:773:VAL:N	4:B:1029:GLY:O	2.50	0.45
4:B:939:SER:OG	4:B:940:GLU:N	2.50	0.45
4:B:1072:GLY:H	4:B:1075:GLU:HB3	1.81	0.45
6:D:14:THR:H	6:D:17:ASN:HB2	1.81	0.45
15:M:41:TYR:HB3	15:M:52:VAL:HG13	1.99	0.45
17:O:348:THR:OG1	17:O:350:GLU:OE1	2.31	0.45
3:A:536:ILE:HD12	3:A:577:VAL:HG22	1.98	0.45
4:B:535:ASP:OD1	4:B:535:ASP:N	2.44	0.45
5:C:69:ARG:HA	5:C:72:ILE:HG22	1.99	0.45
15:M:246:LYS:O	15:M:250:LEU:CD2	2.64	0.45
3:A:1193:VAL:HG22	3:A:1222:LEU:HD11	1.99	0.45
3:A:1502:PRO:HD2	3:A:1525:ASN:HD21	1.81	0.45
4:B:654:ARG:NE	4:B:691:PHE:HB3	2.31	0.45
4:B:790:ASN:HB2	4:B:946:ASP:HA	1.99	0.44
4:B:884:GLU:HB2	4:B:904:LYS:HB3	1.99	0.44
5:C:46:SER:OG	5:C:53:ASN:HB2	2.17	0.44
5:C:240:LYS:HD3	5:C:265:ALA:N	2.32	0.44
11:I:9:PHE:HA	11:I:16:LEU:HA	1.98	0.44
2:U:65:DG:O5'	2:U:65:DG:C8	2.71	0.44
3:A:618:TYR:CA	4:B:783:MET:HE1	2.47	0.44
3:A:687:PHE:O	3:A:726:TRP:N	2.46	0.44
3:A:813:LEU:CD1	3:A:813:LEU:C	2.83	0.44
3:A:827:THR:OG1	3:A:828:CYS:N	2.50	0.44
3:A:894:ALA:HA	3:A:897:SER:HB3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1025:LYS:HZ2	3:A:1615:TYR:HB2	1.82	0.44
4:B:280:LEU:HD11	4:B:370:LYS:HD2	1.99	0.44
4:B:282:HIS:CE1	15:M:101:VAL:HA	2.52	0.44
1:T:12:DC:C5	1:T:12:DC:OP2	2.70	0.44
2:U:58:DA:H2"	2:U:59:DG:H5"	1.98	0.44
3:A:1236:PRO:HG2	3:A:1523:GLY:HA2	1.99	0.44
3:A:1298:ASP:OD1	3:A:1298:ASP:N	2.50	0.44
4:B:359:LEU:HD12	4:B:370:LYS:CE	2.48	0.44
4:B:526:GLY:HA3	4:B:651:ARG:HH22	1.82	0.44
4:B:768:GLY:N	4:B:1032:TYR:OH	2.50	0.44
5:C:136:LEU:HB3	5:C:204:LEU:HG	1.99	0.44
16:N:39:PRO:O	16:N:51:GLN:NE2	2.50	0.44
3:A:467:PHE:HA	3:A:471:MET:HB2	2.00	0.44
3:A:882:ILE:HA	3:A:888:LYS:HB3	1.99	0.44
4:B:334:PHE:HB3	4:B:353:VAL:HG11	2.00	0.44
4:B:658:LEU:HD22	4:B:660:LYS:HE2	2.00	0.44
4:B:1113:THR:HA	4:B:1128:CYS:HA	1.98	0.44
4:B:1128:CYS:SG	4:B:1129:ARG:N	2.90	0.44
16:N:55:LEU:HG	16:N:136:VAL:HG23	2.00	0.44
3:A:404:SER:HB3	15:M:170:LYS:HE3	2.00	0.44
4:B:75:ASP:HA	4:B:432:ILE:HG12	1.99	0.44
6:D:94:ARG:NH1	9:G:151:ASP:OD2	2.51	0.44
15:M:22:ALA:H	16:N:110:LEU:HD11	1.83	0.44
3:A:397:ARG:NH2	15:M:172:LEU:H	2.16	0.44
4:B:1071:VAL:HG23	4:B:1091:ARG:HH21	1.82	0.44
5:C:128:ASP:HB2	5:C:173:GLY:HA3	1.99	0.44
8:F:82:THR:HA	8:F:83:PRO:HD3	1.77	0.44
13:K:104:ARG:HE	13:K:106:GLN:HE22	1.65	0.44
3:A:107:HIS:CE1	3:A:330:LYS:HD3	2.53	0.44
3:A:316:LEU:HG	3:A:318:THR:H	1.81	0.44
3:A:1333:ILE:HD11	3:A:1476:LEU:HD21	2.00	0.44
4:B:936:MET:HB3	4:B:945:PRO:HD2	2.00	0.44
5:C:188:ASP:OD1	5:C:188:ASP:N	2.43	0.44
10:H:93:TYR:HD2	10:H:143:LEU:HD23	1.82	0.44
12:J:57:ILE:O	12:J:61:LEU:N	2.48	0.44
3:A:264:ASN:HA	3:A:267:LYS:HG2	2.00	0.44
3:A:399:LEU:HD22	3:A:402:ASP:HB3	1.99	0.44
3:A:790:LYS:HE3	3:A:791:TYR:CZ	2.53	0.44
3:A:985:ARG:HD3	3:A:986:PHE:H	1.72	0.44
4:B:282:HIS:CG	15:M:101:VAL:HG12	2.53	0.44
4:B:1116:SER:N	4:B:1125:THR:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1195:ARG:HH12	4:B:1197:ARG:NE	2.16	0.44
17:O:117:GLN:HE21	17:O:150:TRP:HH2	1.64	0.44
17:O:482:TYR:HD1	17:O:524:VAL:HG11	1.83	0.44
3:A:124:LEU:HD12	3:A:189:VAL:HG13	2.00	0.43
3:A:520:ARG:HH22	17:O:377:TYR:HB3	1.82	0.43
3:A:551:VAL:HA	3:A:554:ARG:HB2	2.00	0.43
3:A:1131:LYS:HG3	3:A:1132:TYR:CD2	2.53	0.43
3:A:1558:ALA:HA	3:A:1561:THR:HG22	1.99	0.43
5:C:113:LEU:HD21	5:C:132:ILE:HD11	2.00	0.43
6:D:20:VAL:HG22	6:D:21:VAL:HG23	2.00	0.43
15:M:235:PRO:HB3	15:M:287:LEU:HB3	2.00	0.43
2:U:55:DT:H72	2:U:56:DA:H61	1.82	0.43
3:A:41:LEU:CD1	3:A:43:HIS:ND1	2.49	0.43
3:A:537:GLN:HE21	3:A:576:LYS:H	1.66	0.43
3:A:683:LYS:HB3	10:H:23:VAL:HG21	2.00	0.43
3:A:1012:LYS:HB3	3:A:1198:THR:HG23	1.99	0.43
5:C:221:PRO:HB2	5:C:308:MET:HG3	2.00	0.43
7:E:17:ARG:NH2	7:E:35:VAL:O	2.51	0.43
1:T:12:DC:C3'	1:T:12:DC:C6	3.01	0.43
3:A:912:VAL:C	3:A:914:ASP:H	2.20	0.43
3:A:1069:CYS:HB2	3:A:1076:LEU:HD21	1.99	0.43
3:A:1090:ASP:HB3	3:A:1132:TYR:HD1	1.83	0.43
4:B:252:TYR:HB2	4:B:381:LEU:HD11	1.99	0.43
15:M:252:GLN:HG3	15:M:254:SER:H	1.83	0.43
15:M:384:ILE:HG12	15:M:392:TYR:CE1	2.54	0.43
17:O:60:LEU:HD11	17:O:100:LEU:HA	2.00	0.43
17:O:475:ALA:O	17:O:479:ALA:N	2.48	0.43
1:T:7:DA:H2"	1:T:8:DA:N7	2.33	0.43
3:A:1040:ASP:HB3	3:A:1044:THR:HG23	2.01	0.43
16:N:157:ARG:NH1	16:N:159:ASP:OD1	2.51	0.43
4:B:654:ARG:HE	4:B:691:PHE:HB3	1.84	0.43
4:B:663:ILE:H	4:B:663:ILE:HG13	1.62	0.43
4:B:1038:HIS:HB3	4:B:1039:MET:H	1.65	0.43
5:C:243:SER:O	5:C:247:PHE:N	2.48	0.43
9:G:141:SER:HB2	9:G:144:HIS:HB3	2.01	0.43
12:J:3:VAL:HG11	12:J:15:GLY:HA2	2.00	0.43
3:A:102:CYS:HB2	3:A:109:ARG:HG2	2.00	0.43
3:A:1128:ASN:OD1	3:A:1131:LYS:N	2.43	0.43
4:B:658:LEU:CD2	4:B:660:LYS:HE2	2.49	0.43
4:B:953:ALA:O	4:B:957:ARG:N	2.43	0.43
5:C:57:ILE:HD12	5:C:297:HIS:CG	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:K:132:GLU:O	13:K:136:THR:OG1	2.28	0.43
15:M:264:TYR:O	15:M:268:LEU:N	2.51	0.43
16:N:80:MET:H	16:N:87:TYR:H	1.67	0.43
17:O:141:LYS:O	17:O:145:SER:OG	2.36	0.43
3:A:749:LEU:H	3:A:771:PHE:HB2	1.84	0.43
4:B:292:ILE:HG22	4:B:293:ILE:N	2.34	0.43
4:B:445:TYR:HA	4:B:448:ARG:HE	1.83	0.43
7:E:24:LYS:NZ	7:E:32:GLN:OE1	2.51	0.43
15:M:16:GLN:HG3	15:M:92:LYS:HZ2	1.83	0.43
15:M:361:VAL:HA	15:M:364:PHE:HD2	1.83	0.43
3:A:1559:ARG:HG2	3:A:1586:ALA:HB1	2.01	0.43
4:B:612:LYS:NZ	4:B:622:ILE:O	2.39	0.43
4:B:795:GLU:HG3	5:C:217:ALA:H	1.83	0.43
15:M:268:LEU:C	15:M:268:LEU:CD2	2.86	0.43
4:B:234:ILE:HG23	4:B:381:LEU:HD23	2.01	0.43
4:B:742:TYR:CE2	4:B:1037:ARG:HG3	2.54	0.43
7:E:37:LEU:HG	7:E:41:ASP:HB3	2.01	0.43
9:G:111:THR:HG23	9:G:113:PHE:H	1.84	0.43
1:T:12:DC:H2'	1:T:13:DT:C7	2.49	0.43
3:A:495:ILE:HB	3:A:605:VAL:HG12	2.01	0.43
3:A:745:PRO:HD2	3:A:1075:ALA:HA	2.01	0.43
4:B:703:LEU:HD21	4:B:920:ARG:HH22	1.83	0.43
4:B:1196:LEU:HD13	4:B:1197:ARG:CA	2.48	0.43
7:E:100:ILE:HA	7:E:105:PHE:HD2	1.84	0.43
7:E:151:PRO:HA	7:E:201:LYS:HZ3	1.84	0.43
8:F:95:GLY:HA2	8:F:98:ALA:HB3	2.00	0.43
13:K:104:ARG:HE	13:K:106:GLN:NE2	2.17	0.43
3:A:681:THR:HA	3:A:728:GLY:HA3	2.01	0.42
3:A:1094:ALA:O	3:A:1098:SER:N	2.51	0.42
3:A:1260:LYS:N	3:A:1505:ASP:O	2.40	0.42
3:A:1525:ASN:HD21	3:A:1528:ALA:HB3	1.84	0.42
4:B:617:THR:HB	4:B:620:LEU:HB2	2.01	0.42
8:F:124:GLU:O	8:F:128:LYS:N	2.44	0.42
17:O:62:ASP:O	17:O:67:ASP:N	2.46	0.42
3:A:68:ASP:OD1	3:A:68:ASP:N	2.51	0.42
3:A:771:PHE:HB3	3:A:774:GLY:HA2	2.01	0.42
3:A:1004:GLU:HA	3:A:1007:ILE:HB	2.01	0.42
4:B:700:LEU:O	4:B:703:LEU:HD12	2.19	0.42
4:B:752:VAL:HG22	4:B:981:SER:HB3	2.00	0.42
4:B:785:ASP:OD2	4:B:957:ARG:NH2	2.53	0.42
4:B:847:TYR:HD2	4:B:850:THR:HG21	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:279:VAL:HG23	5:C:280:LEU:HD12	2.01	0.42
15:M:12:ILE:HD12	16:N:68:LYS:HA	2.02	0.42
4:B:188:ASP:OD1	4:B:188:ASP:N	2.52	0.42
5:C:82:TYR:O	5:C:207:HIS:N	2.50	0.42
7:E:29:PHE:N	7:E:63:ASN:O	2.49	0.42
11:I:7:LEU:HD22	11:I:8:ILE:H	1.83	0.42
4:B:264:TRP:CD1	4:B:265:ARG:HD3	2.54	0.42
4:B:284:SER:HA	4:B:288:ILE:HD11	2.02	0.42
5:C:79:ALA:HA	5:C:106:LEU:HD12	2.00	0.42
5:C:262:SER:OG	5:C:263:ASP:N	2.53	0.42
15:M:39:ASP:HA	16:N:118:SER:HB2	2.02	0.42
3:A:369:LEU:HD22	3:A:370:PRO:HD2	2.01	0.42
3:A:1256:LYS:C	3:A:1307:ASP:OD2	2.57	0.42
4:B:258:VAL:HG12	4:B:309:LEU:HG	2.01	0.42
4:B:491:ILE:HD11	4:B:742:TYR:CG	2.55	0.42
4:B:552:SER:HB3	4:B:648:ARG:HB3	2.01	0.42
4:B:708:ASP:OD2	4:B:984:TRP:N	2.51	0.42
4:B:874:TYR:CZ	4:B:876:SER:HB3	2.55	0.42
5:C:42:VAL:HG11	13:K:135:PHE:HD1	1.84	0.42
7:E:92:THR:HA	7:E:95:THR:HG22	2.00	0.42
9:G:167:THR:HG23	9:G:218:VAL:HB	2.02	0.42
10:H:78:SER:OG	10:H:79:TRP:N	2.51	0.42
12:J:28:ASP:HB3	12:J:30:LEU:HD23	2.01	0.42
15:M:71:GLN:NE2	15:M:95:VAL:O	2.52	0.42
3:A:985:ARG:HH21	3:A:986:PHE:HB2	1.79	0.42
4:B:164:MET:HG3	4:B:192:GLY:HA2	2.02	0.42
16:N:41:ASN:HD21	16:N:51:GLN:H	1.68	0.42
3:A:516:ILE:HG21	17:O:376:TYR:HE2	1.85	0.42
4:B:368:GLN:HE22	15:M:99:LYS:NZ	2.17	0.42
5:C:123:ASP:N	5:C:123:ASP:OD1	2.53	0.42
5:C:131:THR:HG23	5:C:208:CYS:H	1.85	0.42
5:C:279:VAL:HA	5:C:282:TYR:HD2	1.85	0.42
15:M:360:VAL:O	15:M:364:PHE:N	2.52	0.42
1:T:13:DT:H6	1:T:13:DT:O5'	2.03	0.42
3:A:61:LEU:HD23	15:M:306:LYS:HZ1	1.85	0.42
3:A:489:ASN:OD1	3:A:701:ARG:NH2	2.44	0.42
3:A:642:ASN:HB3	4:B:1086:PHE:CE2	2.54	0.42
3:A:1637:PRO:HA	3:A:1640:ARG:HB3	2.00	0.42
4:B:156:ARG:HH12	4:B:451:MET:HB3	1.84	0.42
4:B:504:HIS:ND1	4:B:542:LEU:HB3	2.35	0.42
5:C:138:VAL:HG13	5:C:159:ASN:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:N:86:ASP:OD1	16:N:86:ASP:N	2.53	0.42
2:U:64:DT:H4'	2:U:64:DT:OP1	2.19	0.42
4:B:470:LEU:HD11	4:B:476:LEU:CD1	2.19	0.42
4:B:981:SER:OG	4:B:982:THR:N	2.52	0.42
7:E:26:ARG:HH12	7:E:107:THR:HG21	1.83	0.42
9:G:219:ASP:OD1	9:G:219:ASP:N	2.46	0.42
3:A:614:LEU:HD13	3:A:614:LEU:HA	1.92	0.42
3:A:985:ARG:HD3	3:A:986:PHE:CA	2.46	0.42
3:A:1294:MET:HB3	3:A:1470:CYS:HB3	2.02	0.42
4:B:57:ASP:O	4:B:62:ASN:ND2	2.53	0.42
4:B:90:TYR:OH	4:B:146:ASN:OD1	2.38	0.42
15:M:332:ILE:HG23	15:M:336:ILE:HB	2.01	0.42
15:M:342:PHE:O	15:M:396:THR:OG1	2.27	0.42
3:A:771:PHE:HE2	3:A:793:ILE:HD13	1.84	0.41
3:A:985:ARG:CD	3:A:985:ARG:C	2.86	0.41
3:A:1224:GLU:O	3:A:1228:THR:OG1	2.29	0.41
4:B:469:ASN:OD1	4:B:469:ASN:N	2.52	0.41
4:B:947:ILE:HG21	4:B:1033:TYR:CE2	2.54	0.41
15:M:77:VAL:HG22	16:N:58:PHE:HE1	1.84	0.41
16:N:113:SER:N	16:N:117:GLU:OE2	2.51	0.41
3:A:649:ASN:OD1	8:F:90:ARG:NH1	2.53	0.41
3:A:990:ILE:HD11	3:A:995:TYR:HB2	2.02	0.41
4:B:28:PRO:HG2	4:B:181:VAL:HG21	2.03	0.41
4:B:206:LEU:HD12	4:B:403:LEU:HD12	2.02	0.41
4:B:501:ARG:HH12	4:B:550:ARG:HH12	1.66	0.41
3:A:75:HIS:HE1	4:B:1114:GLN:HG2	1.85	0.41
4:B:327:LEU:O	4:B:330:LEU:HG	2.20	0.41
4:B:1015:SER:OG	4:B:1019:GLY:N	2.53	0.41
5:C:43:ASN:HB2	5:C:55:ASP:HB2	2.01	0.41
5:C:92:ILE:HD12	12:J:2:ILE:HD11	2.02	0.41
15:M:247:LEU:HA	15:M:250:LEU:HD21	2.01	0.41
15:M:332:ILE:HD13	15:M:353:LEU:HD21	2.01	0.41
3:A:1061:SER:OG	3:A:1062:HIS:ND1	2.54	0.41
3:A:1105:ARG:HH21	7:E:207:ARG:NH1	2.19	0.41
4:B:260:PHE:HB2	4:B:271:VAL:HG12	2.02	0.41
4:B:549:CYS:O	4:B:550:ARG:NE	2.49	0.41
5:C:65:ASN:HA	5:C:68:ARG:HB3	2.02	0.41
5:C:99:HIS:HB3	14:L:69:ALA:HA	2.02	0.41
15:M:21:VAL:HB	15:M:91:TYR:HE2	1.84	0.41
15:M:305:ILE:HA	15:M:318:TYR:HA	2.01	0.41
1:T:14:DT:H4'	3:A:1014:SER:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:51:ASP:OD1	3:A:52:LEU:N	2.54	0.41
3:A:263:ASN:HA	3:A:266:VAL:HG12	2.02	0.41
3:A:1121:ASP:OD1	3:A:1121:ASP:N	2.48	0.41
3:A:1256:LYS:HB3	3:A:1307:ASP:CG	2.41	0.41
3:A:1563:VAL:HG23	3:A:1582:LEU:HB3	2.02	0.41
4:B:934:ILE:HD12	5:C:69:ARG:HG3	2.02	0.41
15:M:361:VAL:HA	15:M:364:PHE:CD2	2.55	0.41
16:N:82:ILE:HG13	16:N:84:LYS:H	1.86	0.41
3:A:54:LEU:HD22	3:A:368:ARG:NH2	2.35	0.41
3:A:107:HIS:ND1	3:A:331:GLU:OE2	2.38	0.41
4:B:93:ASN:OD1	4:B:93:ASN:N	2.53	0.41
4:B:820:PRO:CD	15:M:356:LYS:HZ2	2.30	0.41
5:C:122:ASP:HA	5:C:125:LYS:HE3	2.03	0.41
3:A:462:LYS:HG3	3:A:463:LYS:H	1.84	0.41
3:A:1612:LYS:HE3	3:A:1612:LYS:HB2	1.89	0.41
5:C:73:SER:O	5:C:213:GLY:N	2.54	0.41
9:G:142:ALA:N	17:O:142:ILE:O	2.54	0.41
13:K:80:ILE:HD13	13:K:80:ILE:HA	1.94	0.41
15:M:112:LYS:HE2	15:M:112:LYS:HB2	1.85	0.41
17:O:234:ILE:HG13	17:O:367:LEU:HD21	2.01	0.41
3:A:57:PHE:HE1	3:A:370:PRO:HD2	1.86	0.41
4:B:320:LEU:HG	4:B:326:VAL:HG22	2.02	0.41
4:B:501:ARG:HG3	4:B:544:HIS:HB2	2.01	0.41
4:B:703:LEU:HD13	4:B:704:THR:HA	2.03	0.41
4:B:714:ARG:CZ	4:B:957:ARG:HD2	2.51	0.41
8:F:138:LEU:HA	8:F:139:PRO:HD3	1.80	0.41
17:O:428:ILE:HD12	17:O:439:ILE:HD12	2.02	0.41
3:A:481:ARG:NH2	3:A:592:GLN:HE22	2.19	0.41
3:A:934:LYS:HG2	4:B:955:PRO:HB2	2.03	0.41
3:A:1137:SER:OG	3:A:1138:GLU:N	2.53	0.41
4:B:73:ILE:HD11	4:B:429:ARG:HB3	2.02	0.41
4:B:74:PHE:HE1	4:B:94:LYS:HA	1.86	0.41
4:B:359:LEU:HD12	4:B:370:LYS:HE3	1.95	0.41
4:B:480:GLN:HB3	4:B:508:PHE:HB2	2.02	0.41
4:B:564:ILE:H	4:B:564:ILE:HG13	1.75	0.41
13:K:87:GLU:HB3	13:K:108:TYR:CE2	2.56	0.41
15:M:10:ILE:HD12	15:M:10:ILE:HA	1.83	0.41
15:M:320:ILE:HG23	15:M:324:ASN:HB2	2.03	0.41
3:A:185:ARG:HA	3:A:188:TYR:HB3	2.03	0.41
3:A:721:LYS:NZ	10:H:94:ASP:HB2	2.36	0.41
4:B:529:CYS:SG	4:B:532:HIS:HB3	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:873:THR:OG1	4:B:874:TYR:N	2.52	0.41
4:B:1048:SER:OG	4:B:1049:THR:N	2.54	0.41
5:C:31:TRP:CD2	13:K:82:LYS:HD3	2.56	0.41
13:K:48:LYS:HB3	13:K:64:GLN:HB3	2.02	0.41
15:M:268:LEU:CG	15:M:331:TYR:HE2	2.33	0.41
17:O:426:SER:HB3	17:O:594:TYR:HB2	2.03	0.41
3:A:208:PHE:CZ	3:A:213:ASN:HB3	2.56	0.40
3:A:749:LEU:HA	3:A:749:LEU:HD13	1.89	0.40
3:A:1050:TYR:HB3	3:A:1054:ALA:HA	2.02	0.40
4:B:628:TYR:HD1	4:B:640:LEU:HD23	1.85	0.40
4:B:665:GLY:N	4:B:668:GLU:OE1	2.39	0.40
6:D:25:THR:HB	9:G:42:PRO:HG2	2.02	0.40
16:N:57:LYS:N	16:N:137:PHE:O	2.49	0.40
16:N:133:PHE:HE2	16:N:136:VAL:HB	1.86	0.40
1:T:12:DC:H3'	1:T:12:DC:C6	2.55	0.40
1:T:13:DT:O5'	1:T:13:DT:C6	2.74	0.40
3:A:585:ASP:OD1	3:A:585:ASP:N	2.54	0.40
3:A:1019:LEU:HD11	3:A:1193:VAL:HG13	2.04	0.40
3:A:1020:GLN:O	3:A:1024:THR:OG1	2.30	0.40
3:A:1178:LEU:C	3:A:1178:LEU:CD1	2.86	0.40
4:B:587:GLN:HA	4:B:592:ILE:HG22	2.03	0.40
9:G:69:LEU:HA	9:G:72:LYS:HE3	2.03	0.40
15:M:262:LEU:CG	15:M:265:LEU:CD1	2.69	0.40
15:M:298:ILE:HA	15:M:302:PHE:HD2	1.85	0.40
17:O:94:ASN:HD21	17:O:132:THR:HA	1.85	0.40
3:A:757:ASN:HD22	3:A:765:LEU:HD13	1.84	0.40
3:A:942:GLN:HA	3:A:947:LEU:HA	2.04	0.40
4:B:359:LEU:CD1	4:B:370:LYS:HZ2	2.34	0.40
4:B:737:SER:HA	4:B:904:LYS:NZ	2.36	0.40
4:B:1108:GLY:HA3	4:B:1198:TYR:H	1.86	0.40
5:C:47:LEU:HA	5:C:52:ALA:HA	2.03	0.40
15:M:312:ARG:NH1	15:M:317:SER:H	2.19	0.40
17:O:530:ARG:HH22	17:O:613:TRP:HE1	1.68	0.40
3:A:586:VAL:HG21	3:A:648:LEU:HB2	2.02	0.40
3:A:995:TYR:HE2	4:B:697:LEU:HD11	1.87	0.40
9:G:85:GLU:OE2	9:G:121:ASN:ND2	2.54	0.40
9:G:152:ALA:O	17:O:185:SER:OG	2.38	0.40
10:H:26:ILE:N	10:H:40:LEU:O	2.42	0.40
15:M:203:PRO:O	15:M:206:SER:OG	2.32	0.40
17:O:205:ARG:HG3	17:O:331:LYS:HE2	2.04	0.40
2:U:63:DT:H2"	2:U:64:DT:C6	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:537:GLN:NE2	3:A:576:LYS:HG2	2.36	0.40
3:A:709:ARG:HG2	3:A:710:SER:H	1.86	0.40
3:A:1550:LEU:HD11	3:A:1594:THR:HA	2.03	0.40
4:B:209:GLN:HE22	4:B:237:ARG:HB2	1.85	0.40
4:B:240:ARG:HG2	4:B:360:VAL:HG11	2.02	0.40
4:B:292:ILE:HG22	4:B:302:LEU:HD11	2.03	0.40
4:B:1095:SER:OG	4:B:1096:SER:N	2.54	0.40
9:G:137:ILE:N	9:G:227:GLY:O	2.42	0.40
15:M:242:TYR:CE1	15:M:352:GLU:HB2	2.57	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	
3	A	1450/1664 (87%)	1234 (85%)	212 (15%)	4 (0%)	41	75
4	B	1160/1203 (96%)	980 (84%)	177 (15%)	3 (0%)	41	75
5	C	300/335 (90%)	269 (90%)	31 (10%)	0	100	100
6	D	55/137 (40%)	50 (91%)	5 (9%)	0	100	100
7	E	212/215 (99%)	192 (91%)	20 (9%)	0	100	100
8	F	98/155 (63%)	87 (89%)	11 (11%)	0	100	100
9	G	196/326 (60%)	168 (86%)	28 (14%)	0	100	100
10	H	130/146 (89%)	112 (86%)	18 (14%)	0	100	100
11	I	62/125 (50%)	49 (79%)	13 (21%)	0	100	100
12	J	67/70 (96%)	59 (88%)	8 (12%)	0	100	100
13	K	98/142 (69%)	87 (89%)	11 (11%)	0	100	100
14	L	41/70 (59%)	36 (88%)	5 (12%)	0	100	100
15	M	386/415 (93%)	303 (78%)	83 (22%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
16	N	127/233 (54%)	96 (76%)	31 (24%)	0	100 100
17	O	457/627 (73%)	392 (86%)	65 (14%)	0	100 100
All	All	4839/5863 (82%)	4114 (85%)	718 (15%)	7 (0%)	54 84

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	913	PRO
3	A	757	ASN
4	B	784	ASP
4	B	684	ASN
3	A	530	TRP
4	B	266	LYS
3	A	35	PRO

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
3	A	1293/1465 (88%)	1263 (98%)	30 (2%)	50 70
4	B	1025/1053 (97%)	1008 (98%)	17 (2%)	60 78
5	C	269/296 (91%)	265 (98%)	4 (2%)	65 80
6	D	56/116 (48%)	55 (98%)	1 (2%)	59 77
7	E	196/197 (100%)	190 (97%)	6 (3%)	40 63
8	F	90/137 (66%)	86 (96%)	4 (4%)	28 55
9	G	180/291 (62%)	177 (98%)	3 (2%)	60 78
10	H	117/128 (91%)	114 (97%)	3 (3%)	46 67
11	I	56/110 (51%)	53 (95%)	3 (5%)	22 50
12	J	64/65 (98%)	63 (98%)	1 (2%)	62 79
13	K	90/130 (69%)	88 (98%)	2 (2%)	52 71
14	L	38/57 (67%)	35 (92%)	3 (8%)	12 39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
15	M	350/371 (94%)	344 (98%)	6 (2%)	60 78
16	N	125/220 (57%)	123 (98%)	2 (2%)	62 79
17	O	427/576 (74%)	427 (100%)	0	100 100
All	All	4376/5212 (84%)	4291 (98%)	85 (2%)	59 75

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	27	LEU
3	A	59	ARG
3	A	92	ASN
3	A	111	LYS
3	A	230	ARG
3	A	259	LYS
3	A	417	ARG
3	A	427	PHE
3	A	466	LEU
3	A	606	ARG
3	A	649	ASN
3	A	667	ARG
3	A	709	ARG
3	A	736	LEU
3	A	777	LEU
3	A	781	LEU
3	A	843	ARG
3	A	866	LYS
3	A	875	LEU
3	A	919	LYS
3	A	946	LEU
3	A	985	ARG
3	A	1021	ARG
3	A	1063	MET
3	A	1110	LYS
3	A	1437	ASN
3	A	1499	ARG
3	A	1567	ASN
3	A	1580	ARG
3	A	1613	MET
4	B	87	ASN
4	B	119	ARG
4	B	127	ARG

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Mol	Chain	Res	Type
4	B	136	LYS
4	B	168	ASN
4	B	224	ASN
4	B	315	LYS
4	B	421	LEU
4	B	476	LEU
4	B	651	ARG
4	B	703	LEU
4	B	716	MET
4	B	785	ASP
4	B	790	ASN
4	B	1063	ARG
4	B	1092	LEU
4	B	1171	ASN
5	C	75	VAL
5	C	165	ARG
5	C	251	PHE
5	C	329	LYS
6	D	90	LYS
7	E	123	LEU
7	E	124	VAL
7	E	177	ARG
7	E	191	LYS
7	E	200	ARG
7	E	215	MET
8	F	65	ARG
8	F	67	LYS
8	F	76	LYS
8	F	104	ASN
9	G	8	ASN
9	G	77	VAL
9	G	160	ASN
10	H	43	ASN
10	H	130	ARG
10	H	146	ARG
11	I	19	ASN
11	I	39	LYS
11	I	63	LYS
12	J	48	ARG
13	K	81	MET
13	K	83	ASN
14	L	28	LYS

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Mol	Chain	Res	Type
14	L	42	ARG
14	L	62	LYS
15	M	126	ASN
15	M	148	ARG
15	M	200	GLN
15	M	246	LYS
15	M	267	LEU
15	M	275	ARG
16	N	37	ASN
16	N	158	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (81) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	26	ASN
3	A	60	ASN
3	A	92	ASN
3	A	116	HIS
3	A	224	HIS
3	A	332	GLN
3	A	336	GLN
3	A	400	ASN
3	A	515	ASN
3	A	521	GLN
3	A	537	GLN
3	A	592	GLN
3	A	634	ASN
3	A	640	ASN
3	A	730	GLN
3	A	795	HIS
3	A	798	HIS
3	A	937	ASN
3	A	939	ASN
3	A	950	GLN
3	A	1047	GLN
3	A	1072	ASN
3	A	1113	HIS
3	A	1237	GLN
3	A	1436	ASN
3	A	1437	ASN
3	A	1532	GLN
3	A	1567	ASN

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Mol	Chain	Res	Type
3	A	1592	GLN
4	B	45	HIS
4	B	87	ASN
4	B	110	ASN
4	B	128	GLN
4	B	168	ASN
4	B	197	ASN
4	B	224	ASN
4	B	231	HIS
4	B	246	GLN
4	B	267	ASN
4	B	361	HIS
4	B	368	GLN
4	B	422	GLN
4	B	547	HIS
4	B	575	HIS
4	B	598	HIS
4	B	715	ASN
4	B	745	GLN
4	B	790	ASN
4	B	975	HIS
4	B	1058	GLN
4	B	1157	GLN
4	B	1171	ASN
5	C	53	ASN
5	C	232	GLN
7	E	99	HIS
7	E	146	HIS
7	E	147	HIS
7	E	153	HIS
8	F	104	ASN
9	G	8	ASN
9	G	65	HIS
9	G	67	ASN
9	G	75	ASN
9	G	154	ASN
9	G	160	ASN
10	H	43	ASN
11	I	19	ASN
13	K	83	ASN
14	L	53	HIS
15	M	126	ASN

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Mol	Chain	Res	Type
15	M	274	ASN
15	M	351	HIS
16	N	37	ASN
16	N	50	GLN
16	N	51	GLN
16	N	52	GLN
16	N	85	HIS
17	O	94	ASN
17	O	228	GLN
17	O	411	ASN
17	O	478	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

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

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

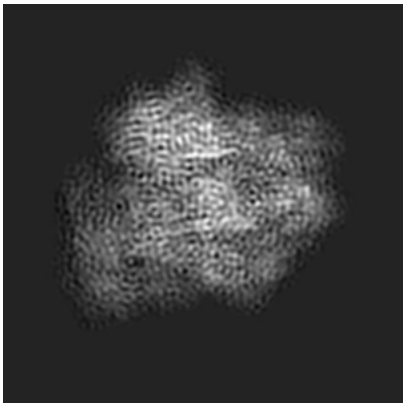
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-4985. These allow visual inspection of the internal detail of the map and identification of artifacts.

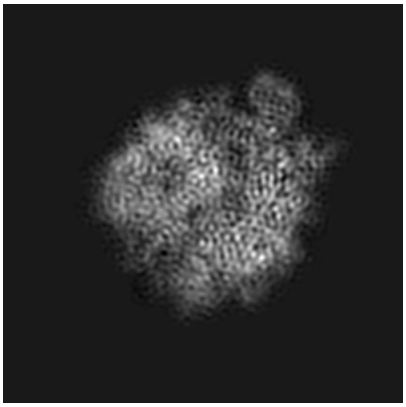
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

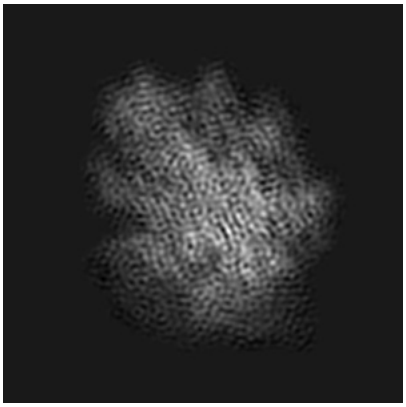
6.1.1 Primary map



X



Y

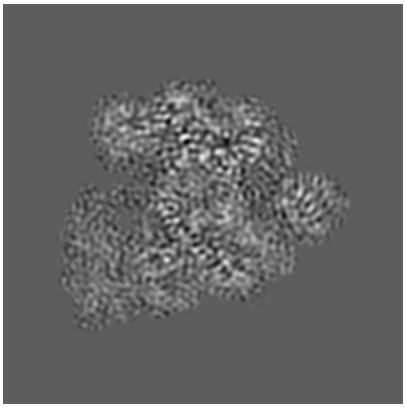


Z

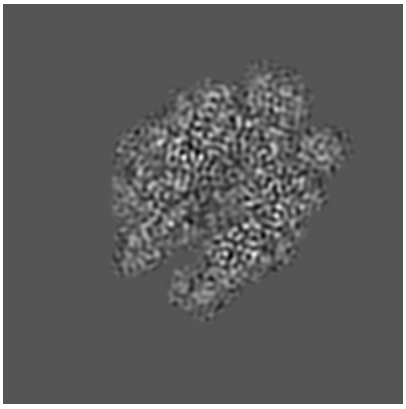
The images above show the map projected in three orthogonal directions.

6.2 Central slices

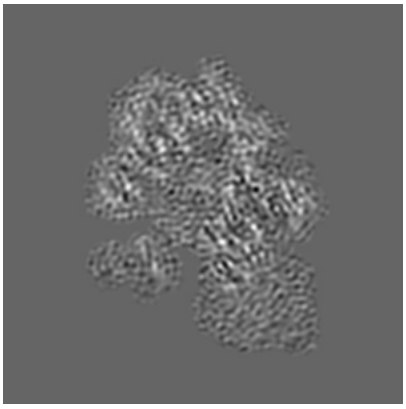
6.2.1 Primary map



X Index: 88



Y Index: 88

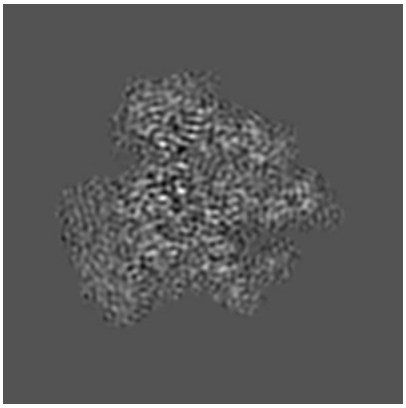


Z Index: 88

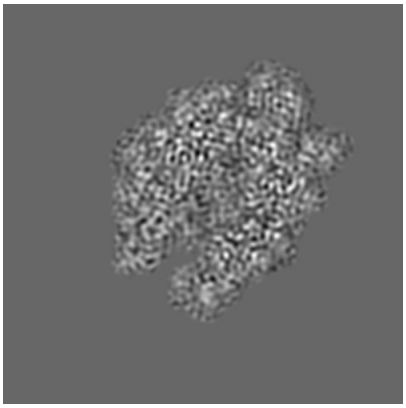
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

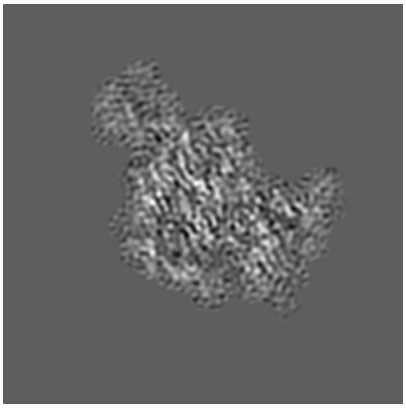
6.3.1 Primary map



X Index: 101



Y Index: 89



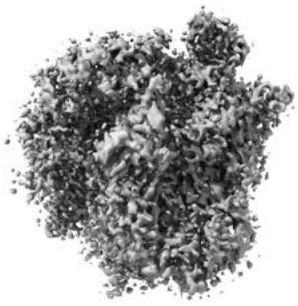
Z Index: 110

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

6.4 Orthogonal surface views (i)

6.4.1 Primary map

X



Y

Z

The images above show the 3D surface view of the map at the recommended contour level 0.05. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

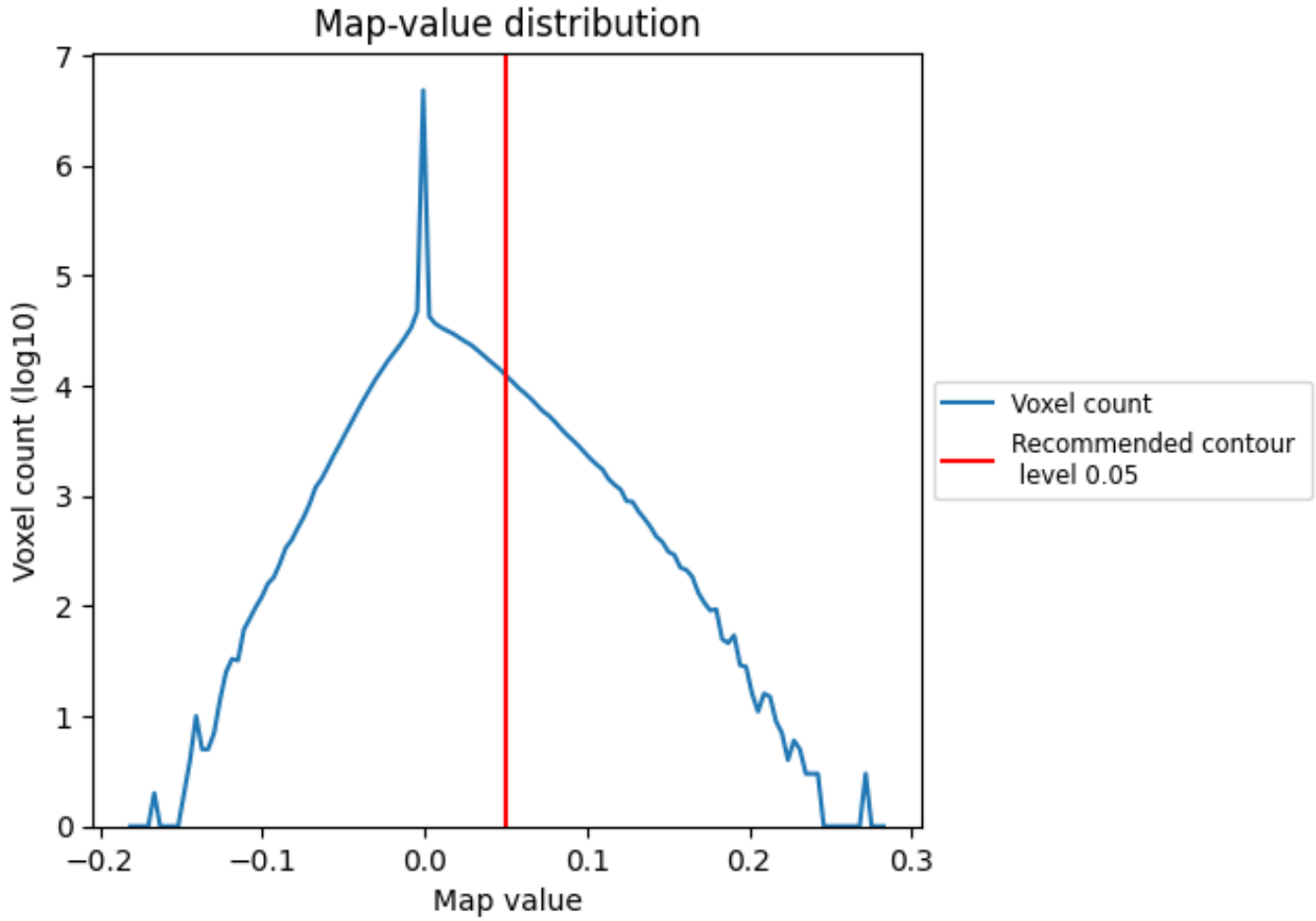
6.5 Mask visualisation (i)

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

7 Map analysis

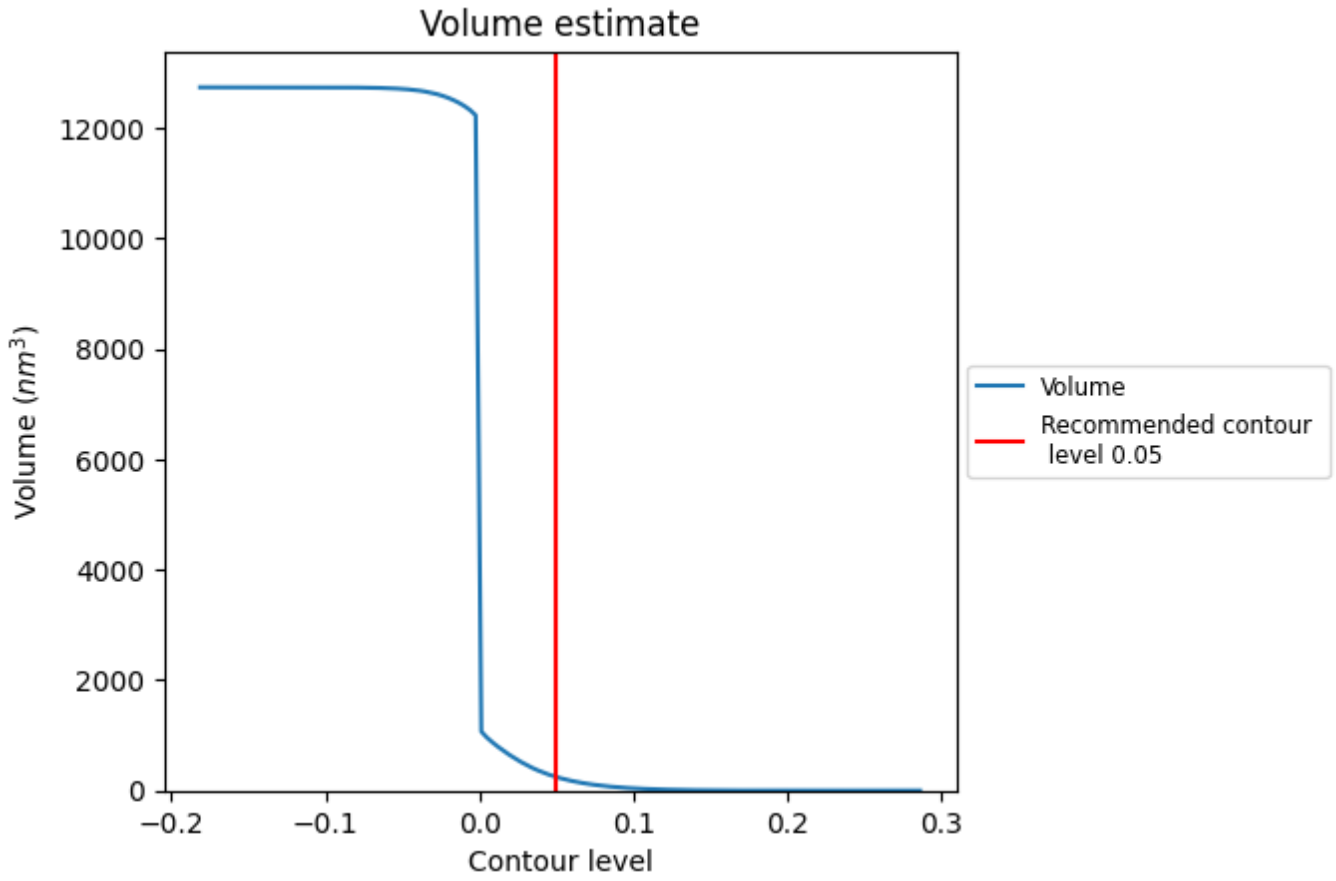
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

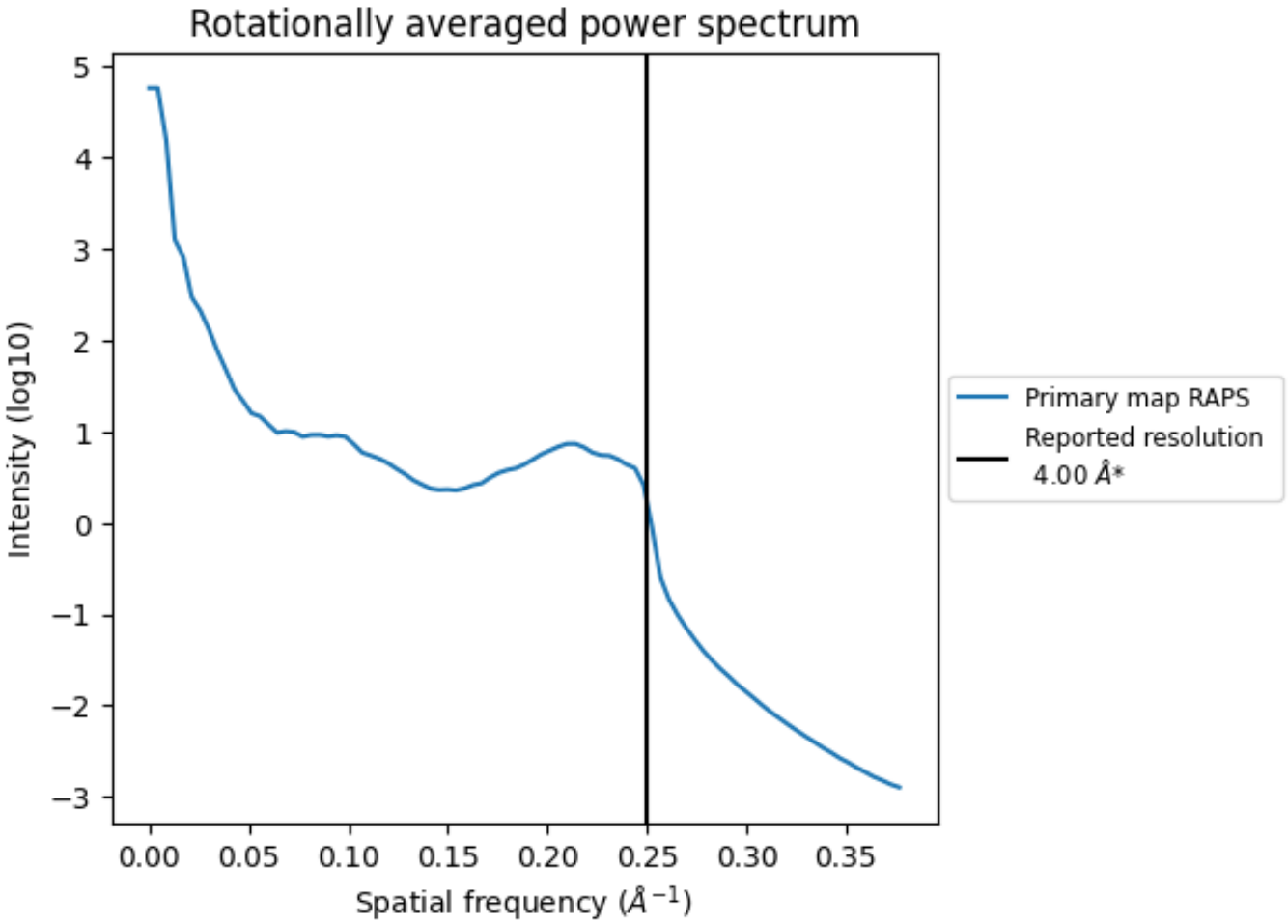
7.2 Volume estimate



The volume at the recommended contour level is 247 nm^3; this corresponds to an approximate mass of 223 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 (i)



*Reported resolution corresponds to spatial frequency of 0.250 Å^{-1}

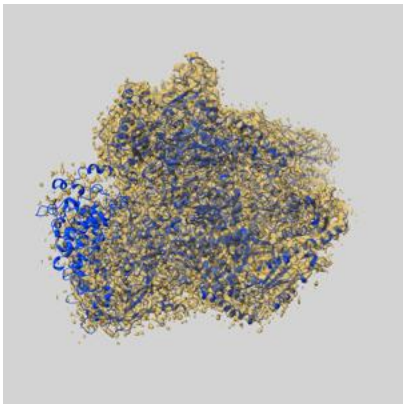
8 Fourier-Shell correlation

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

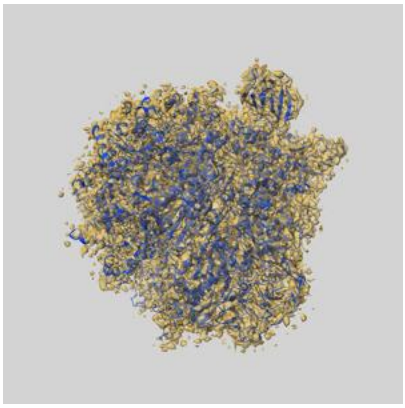
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-4985 and PDB model 6RQT. Per-residue inclusion information can be found in section 3 on page 7.

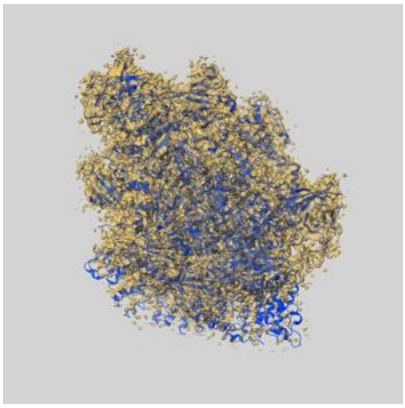
9.1 Map-model overlay



X



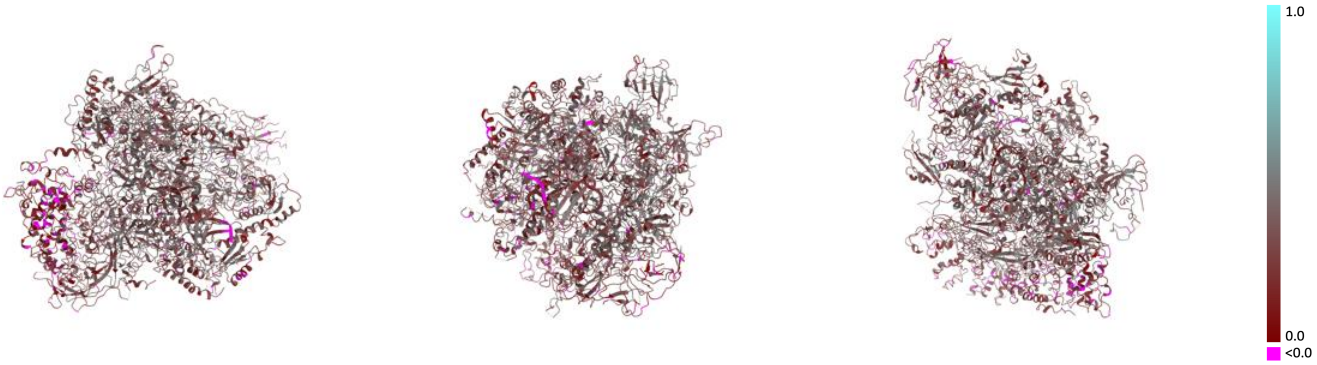
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.05 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.

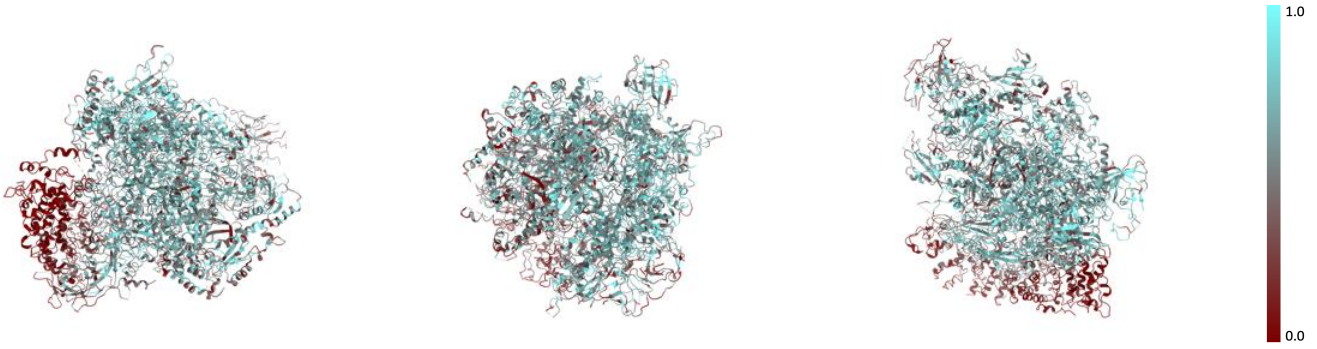
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

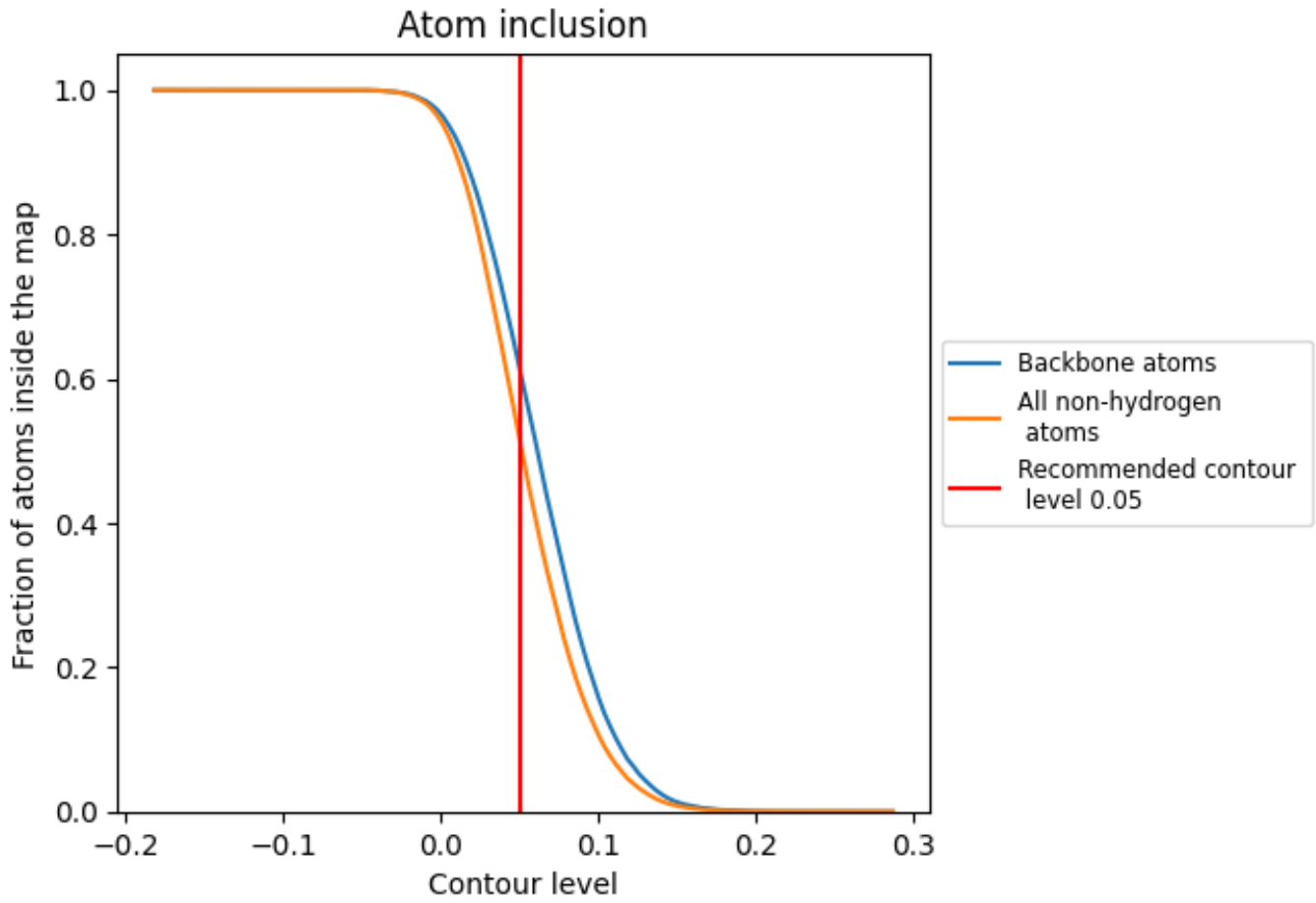
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.05).

9.4 Atom inclusion (i)



At the recommended contour level, 61% of all backbone atoms, 52% 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 (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5159	0.2950
A	0.6031	0.3270
B	0.6298	0.3340
C	0.6382	0.3400
D	0.4039	0.2660
E	0.5855	0.3040
F	0.6117	0.3580
G	0.4049	0.2720
H	0.6294	0.3260
I	0.5128	0.3020
J	0.6739	0.3220
K	0.6134	0.3220
L	0.5271	0.2890
M	0.2761	0.2050
N	0.4868	0.2550
O	0.0588	0.1670
T	0.3819	0.1360
U	0.3481	0.1270