



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:39 PM UTC

PDB ID : 5UH8 / pdb_00005uh8
Title : Crystal structure of Mycobacterium tuberculosis transcription initiation complex containing 4nt RNA
Authors : Lin, W.; Das, K.; Feng, Y.; Ebright, R.H.
Deposited on : 2017-01-11
Resolution : 4.18 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

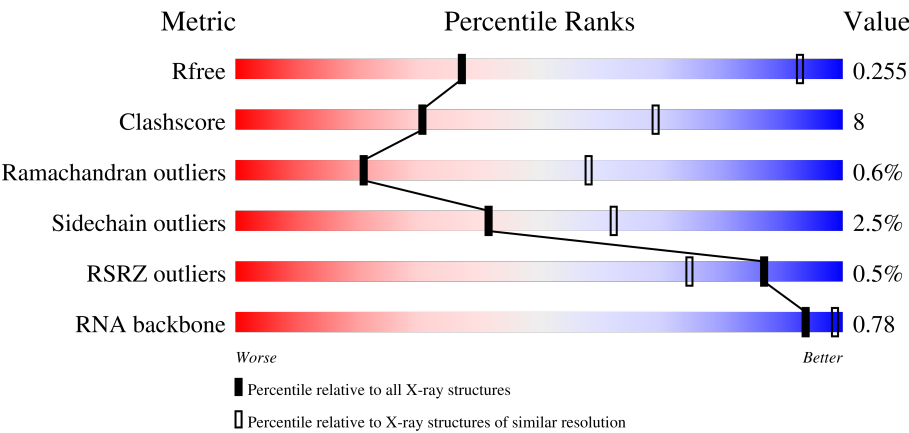
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.18 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)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1026 (4.50-3.86)
Clashscore	190562	1071 (4.50-3.86)
Ramachandran outliers	187476	1014 (4.52-3.84)
Sidechain outliers	187428	1000 (4.52-3.84)
RSRZ outliers	180081	1023 (4.50-3.86)
RNA backbone	3983	1026 (5.04-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	347	54%10%35%
1	B	347	53%12%35%
2	C	1178	73%22%
3	D	1316	74%21%

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Mol	Chain	Length	Quality of chain
4	E	110	% 55%18%26%
5	F	528	% 44%16%39%
6	H	23	35%65%
7	G	16	75%25%
8	I	4	100%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 26101 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	224	Total	C	N	O	S	0	0	0
			1704	1072	295	335	2			
1	B	227	Total	C	N	O	S	0	0	0
			1715	1080	291	342	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1126	Total	C	N	O	S	0	0	0
			8714	5454	1528	1693	39			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1265	Total	C	N	O	S	0	0	0
			9887	6188	1793	1866	40			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	81	Total	C	N	O	0	0	0
			637	408	106	123			

- Molecule 5 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	322	Total	C	N	O	S	0	0	0
			2555	1589	461	496	9			

- Molecule 6 is a DNA chain called DNA (5'-D(*TP*AP*TP*AP*AP*TP*GP*GP*GP*AP*GP*CP*TP*GP*TP*CP*AP*CP*GP*GP*AP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	H	23	Total	C	N	O	P	0	0	0
			476	227	91	136	22			

- Molecule 7 is a DNA chain called DNA (5'-D(*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*CP*GP*AP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	16	Total	C	N	O	P	0	0	0
			328	156	63	94	15			

- Molecule 8 is a RNA chain called RNA (5'-R(*UP*CP*GP*A)-3').

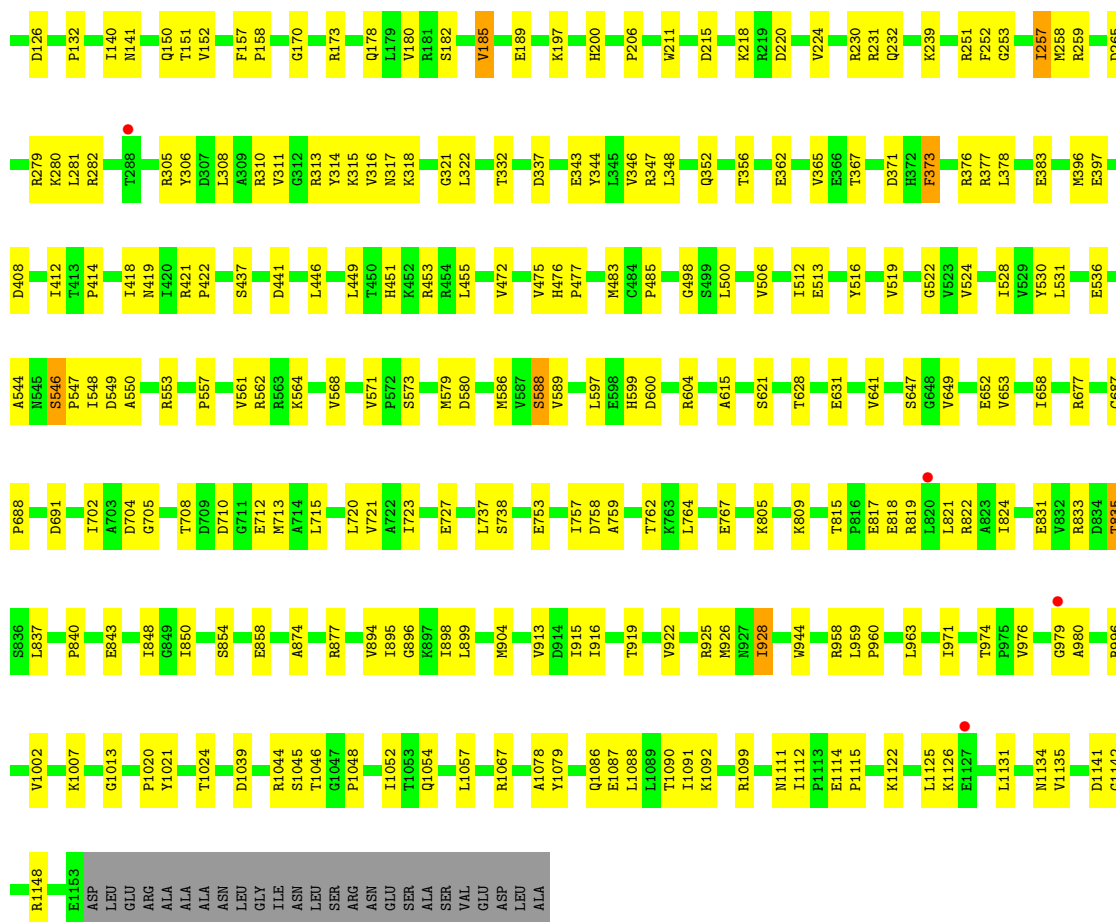
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	I	4	Total	C	N	O	P	0	0	0
			82	38	15	26	3			

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	D	2	Total	Zn	0	0
			2	2		

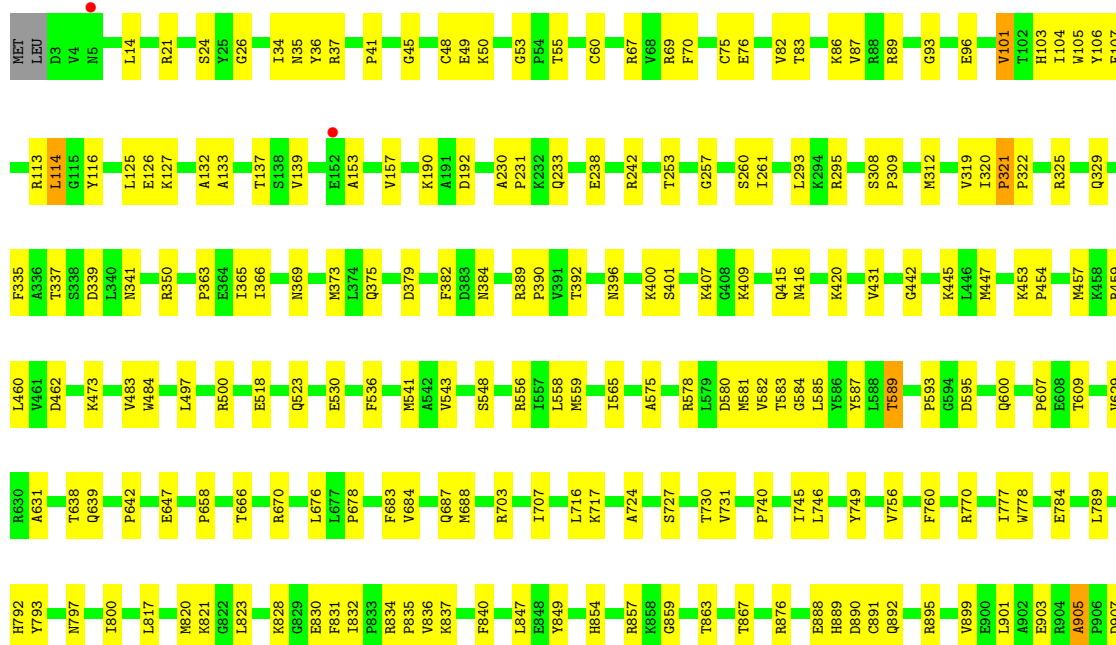
- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

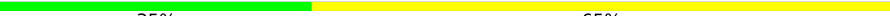
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	1	Total	Mg	0	0
			1	1		

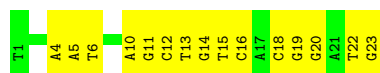


• Molecule 3: DNA-directed RNA polymerase subunit beta'

Chain D: 74% 21%



Chain H:  35% 65%



- Molecule 7: DNA (5'-D(*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*CP*GP*AP*GP*G)-3')

Chain G:  75% 25%



- Molecule 8: RNA (5'-R(*UP*CP*GP*A)-3')

Chain I:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	152.00Å 163.67Å 197.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.64 – 4.18 48.64 – 4.18	Depositor EDS
% Data completeness (in resolution range)	80.2 (48.64-4.18) 80.2 (48.64-4.18)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 4.14Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.201 , 0.257 0.206 , 0.255	Depositor DCC
R_{free} test set	1988 reflections (6.64%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.525	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	26101	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.33	0/1730	0.80	0/2354
1	B	0.30	0/1741	0.79	2/2371 (0.1%)
2	C	0.33	0/8873	0.76	5/12031 (0.0%)
3	D	0.37	2/10052 (0.0%)	0.81	15/13591 (0.1%)
4	E	0.36	0/650	0.75	0/886
5	F	0.36	0/2585	0.76	2/3485 (0.1%)
6	H	0.25	0/535	0.39	0/826
7	G	0.26	0/368	0.44	0/567
8	I	0.25	0/91	0.33	0/140
All	All	0.34	2/26625 (0.0%)	0.77	24/36251 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	910	LEU	C-N	14.18	1.51	1.33
3	D	903	GLU	C-N	7.72	1.44	1.33

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	903	GLU	O-C-N	-24.03	94.66	123.01
3	D	903	GLU	CA-C-N	12.96	141.87	121.66
3	D	903	GLU	C-N-CA	12.96	141.87	121.66
3	D	1126	GLY	CA-C-N	6.69	126.19	119.24
3	D	1126	GLY	C-N-CA	6.69	126.19	119.24
3	D	910	LEU	CA-C-N	-6.37	113.52	122.75
3	D	910	LEU	C-N-CA	-6.37	113.52	122.75
3	D	321	PRO	CA-C-N	6.10	125.32	118.97
3	D	321	PRO	C-N-CA	6.10	125.32	118.97
2	C	723	ILE	N-CA-C	5.93	112.56	106.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	797	ASN	CA-C-N	5.67	125.52	119.28
3	D	797	ASN	C-N-CA	5.67	125.52	119.28
5	F	406	THR	CA-C-N	5.54	125.06	119.19
5	F	406	THR	C-N-CA	5.54	125.06	119.19
2	C	546	SER	CA-C-N	5.44	125.45	119.90
2	C	546	SER	C-N-CA	5.44	125.45	119.90
3	D	905	ALA	CA-C-N	5.41	124.59	118.97
3	D	905	ALA	C-N-CA	5.41	124.59	118.97
2	C	46	GLU	CA-C-N	-5.40	113.09	119.84
2	C	46	GLU	C-N-CA	-5.40	113.09	119.84
1	B	184	GLU	CB-CA-C	-5.28	110.04	117.23
1	B	14	LEU	N-CA-C	5.09	118.03	110.14
3	D	53	GLY	CA-C-N	5.04	124.80	119.76
3	D	53	GLY	C-N-CA	5.04	124.80	119.76

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1704	0	1741	23	0
1	B	1715	0	1739	23	0
2	C	8714	0	8636	161	0
3	D	9887	0	9943	172	0
4	E	637	0	635	13	0
5	F	2555	0	2579	69	0
6	H	476	0	261	20	0
7	G	328	0	181	5	0
8	I	82	0	44	0	0
9	D	2	0	0	0	0
10	D	1	0	0	0	0
All	All	26101	0	25759	426	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 8.

All (426) 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)
6:H:16:DC:O2	7:G:12:DG:N2	2.09	0.83
3:D:69:ARG:HE	5:F:485:GLN:HB2	1.50	0.74
1:B:75:GLU:O	1:B:79:ASN:ND2	2.21	0.73
2:C:815:THR:HG22	2:C:817:GLU:H	1.52	0.73
2:C:1122:LYS:HE2	2:C:1148:ARG:HG2	1.69	0.73
3:D:442:GLY:HA3	3:D:523:GLN:HB2	1.69	0.73
3:D:1090:LYS:HB3	3:D:1092:GLU:HG2	1.71	0.72
3:D:891:CYS:SG	3:D:970:THR:OG1	2.48	0.71
2:C:396:MET:HE1	2:C:418:ILE:HG23	1.72	0.71
3:D:589:THR:HG22	3:D:670:ARG:HG2	1.73	0.70
5:F:468:SER:HB3	5:F:471:GLU:HG3	1.74	0.70
2:C:899:LEU:HB2	2:C:904:MET:HE1	1.75	0.69
2:C:513:GLU:HB3	2:C:530:TYR:HB3	1.75	0.69
2:C:974:THR:HG23	2:C:980:ALA:H	1.57	0.69
3:D:104:ILE:HD12	3:D:379:ASP:HB3	1.74	0.68
5:F:522:VAL:HG23	5:F:523:LEU:HD12	1.77	0.67
2:C:113:ASP:HB2	2:C:132:PRO:HG2	1.77	0.67
2:C:1024:THR:H	3:D:730:THR:HG21	1.60	0.66
3:D:965:VAL:HG13	3:D:974:VAL:HG11	1.76	0.66
1:A:107:ALA:HB2	1:A:123:MET:HE2	1.77	0.66
2:C:1045:SER:OG	2:C:1046:THR:N	2.28	0.66
4:E:70:GLN:O	4:E:74:GLY:N	2.28	0.66
1:A:40:ARG:NH1	2:C:1013:GLY:O	2.31	0.64
3:D:638:THR:HG23	3:D:639:GLN:HG2	1.80	0.64
2:C:738:SER:HA	2:C:904:MET:HE3	1.79	0.64
2:C:485:PRO:O	3:D:857:ARG:NH2	2.30	0.64
3:D:890:ASP:OD1	3:D:963:ARG:NH2	2.30	0.64
1:A:213:LYS:HD3	1:B:227:VAL:HG23	1.79	0.64
1:A:214:THR:HA	1:B:230:GLU:HG2	1.79	0.64
5:F:474:VAL:HB	5:F:503:ILE:HG13	1.79	0.64
2:C:279:ARG:HD3	5:F:215:ALA:HB1	1.79	0.63
3:D:137:THR:OG1	3:D:253:THR:O	2.17	0.63
5:F:470:ARG:HB3	5:F:506:ILE:HD13	1.81	0.62
5:F:477:LEU:HD13	5:F:492:ILE:HG23	1.80	0.62
3:D:1164:ARG:NH2	3:D:1216:ALA:O	2.32	0.62
1:A:62:GLU:HG3	1:A:77:ILE:HD12	1.80	0.62
2:C:239:LYS:NZ	2:C:265:ASP:OD2	2.32	0.62
3:D:1085:ARG:HA	3:D:1112:MET:HA	1.81	0.62
2:C:1054:GLN:OE1	2:C:1099:ARG:NH2	2.33	0.62
1:A:197:GLU:OE1	2:C:996:ARG:NH1	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:211:TRP:NE1	6:H:13:DT:O2	2.31	0.61
3:D:914:PRO:O	3:D:1143:ARG:NH1	2.33	0.61
2:C:1111:ASN:OD1	4:E:65:ASN:ND2	2.33	0.61
2:C:305:ARG:HH12	6:H:10:DA:H62	1.49	0.61
3:D:21:ARG:NH2	3:D:96:GLU:OE2	2.33	0.61
2:C:536:GLU:OE2	2:C:562:ARG:NH1	2.30	0.61
3:D:578:ARG:H	3:D:581:MET:HE3	1.66	0.60
2:C:104:SER:HB3	2:C:140:ILE:HB	1.82	0.60
2:C:189:GLU:HB2	2:C:367:THR:HG21	1.83	0.60
3:D:930:VAL:HG22	3:D:936:VAL:HG12	1.82	0.60
3:D:107:PHE:HZ	3:D:126:GLU:HG2	1.67	0.60
3:D:190:LYS:HE3	3:D:192:ASP:HB3	1.82	0.60
4:E:42:GLU:O	4:E:46:ARG:NH1	2.34	0.60
2:C:737:LEU:HB2	2:C:898:ILE:HG12	1.83	0.60
3:D:1165:VAL:HG12	3:D:1205:PRO:HA	1.84	0.60
2:C:658:ILE:HD11	2:C:688:PRO:HB3	1.84	0.60
2:C:561:VAL:HG21	2:C:571:VAL:HB	1.84	0.59
3:D:257:GLY:O	3:D:260:SER:OG	2.19	0.59
3:D:589:THR:HG21	3:D:688:MET:HG2	1.82	0.59
2:C:453:ARG:NH1	2:C:500:LEU:O	2.35	0.59
3:D:1089:PHE:HA	3:D:1095:SER:HA	1.83	0.59
6:H:15:DT:H2''	6:H:16:DC:H5'	1.85	0.59
2:C:102:SER:O	2:C:141:ASN:ND2	2.34	0.59
2:C:1148:ARG:NH1	3:D:86:LYS:O	2.35	0.59
2:C:41:PHE:HB2	2:C:979:GLY:HA2	1.85	0.59
2:C:588:SER:OG	2:C:589:VAL:N	2.37	0.58
3:D:93:GLY:O	3:D:319:VAL:N	2.28	0.58
3:D:1038:ARG:NH1	6:H:18:DC:O3'	2.36	0.58
1:B:72:ASP:OD1	1:B:73:VAL:N	2.36	0.58
2:C:455:LEU:N	2:C:498:GLY:O	2.36	0.58
3:D:293:LEU:HD21	3:D:1177:PRO:HG2	1.84	0.58
2:C:348:LEU:HD13	2:C:365:VAL:HG12	1.84	0.58
3:D:1090:LYS:HG2	3:D:1091:HIS:H	1.69	0.57
2:C:1148:ARG:NH1	3:D:86:LYS:HG3	2.20	0.57
3:D:859:GLY:O	3:D:863:THR:OG1	2.17	0.57
3:D:407:LYS:HE2	3:D:1230:THR:HG21	1.86	0.57
3:D:1055:LEU:HB2	3:D:1101:ASP:HB3	1.86	0.57
1:A:152:ASN:HB3	1:A:163:PRO:HB3	1.85	0.57
2:C:877:ARG:HH12	2:C:1039:ASP:CG	2.13	0.57
2:C:96:ILE:HD13	2:C:397:GLU:HG3	1.86	0.57
2:C:126:ASP:HA	2:C:170:GLY:HA3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:746:LEU:HD13	3:D:837:LYS:HB3	1.87	0.57
3:D:1245:LEU:HD13	3:D:1254:ILE:HD13	1.86	0.57
3:D:585:LEU:O	3:D:589:THR:OG1	2.22	0.56
5:F:360:ALA:HB1	5:F:373:VAL:HG21	1.86	0.56
3:D:373:MET:SD	5:F:318:LEU:HB3	2.45	0.56
5:F:242:ASN:OD1	5:F:243:ALA:N	2.39	0.56
2:C:959:LEU:HD12	2:C:960:PRO:HD2	1.87	0.56
5:F:378:LYS:HD3	5:F:381:ARG:HH11	1.71	0.56
2:C:758:ASP:N	2:C:758:ASP:OD1	2.38	0.56
5:F:506:ILE:HA	5:F:509:LYS:HD2	1.87	0.56
1:B:27:GLU:HG3	1:B:28:PRO:HD2	1.88	0.55
5:F:315:MET:HE2	5:F:359:MET:HE3	1.89	0.55
5:F:499:THR:OG1	5:F:500:ARG:N	2.37	0.55
2:C:48:LEU:HD12	2:C:528:ILE:HD13	1.89	0.55
5:F:364:ARG:HG3	5:F:368:ILE:HG12	1.89	0.55
5:F:345:THR:HB	6:H:4:DA:H8	1.72	0.55
3:D:459:ARG:HA	3:D:462:ASP:HB2	1.88	0.55
3:D:749:TYR:OH	3:D:784:GLU:OE1	2.20	0.54
1:B:170:PRO:HA	1:B:199:LYS:HD2	1.89	0.54
3:D:1274:PRO:HB3	4:E:82:LEU:HD11	1.89	0.54
3:D:24:SER:OG	3:D:26:GLY:O	2.26	0.54
3:D:238:GLU:OE1	5:F:237:LYS:NZ	2.41	0.54
2:C:483:MET:HE3	2:C:498:GLY:HA3	1.90	0.54
2:C:628:THR:N	2:C:631:GLU:OE2	2.37	0.54
3:D:363:PRO:HD2	3:D:366:ILE:HD12	1.89	0.54
5:F:240:LEU:HD21	5:F:301:ARG:HD2	1.90	0.54
2:C:43:LYS:NZ	2:C:544:ALA:O	2.40	0.54
3:D:1092:GLU:HG3	3:D:1094:GLY:H	1.73	0.53
3:D:460:LEU:HD11	3:D:483:VAL:HG12	1.89	0.53
3:D:1051:GLY:HA2	3:D:1069:ASP:HB2	1.91	0.53
3:D:107:PHE:CZ	3:D:126:GLU:HG2	2.43	0.53
1:A:56:ILE:HG12	1:A:136:VAL:HG13	1.90	0.53
2:C:38:ARG:HH11	2:C:712:GLU:CD	2.17	0.53
4:E:38:PRO:HD2	4:E:96:LEU:HD23	1.90	0.53
2:C:824:ILE:HG12	5:F:514:LEU:HD13	1.90	0.53
2:C:1125:LEU:HD22	2:C:1135:VAL:HG11	1.90	0.53
3:D:530:GLU:HB2	3:D:578:ARG:HD2	1.91	0.53
2:C:704:ASP:HB2	2:C:708:THR:HB	1.91	0.52
2:C:1086:GLN:O	2:C:1090:THR:OG1	2.17	0.52
3:D:820:MET:HE3	3:D:837:LYS:HA	1.91	0.52
5:F:256:GLY:HA3	5:F:288:GLY:HA3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1265:ASN:OD1	3:D:1268:ARG:NH2	2.38	0.52
5:F:231:TYR:CE2	5:F:235:ILE:HD11	2.45	0.52
2:C:1112:ILE:HG13	3:D:548:SER:HA	1.90	0.52
5:F:470:ARG:HH11	5:F:506:ILE:HD11	1.75	0.52
6:H:22:DT:H1'	6:H:23:DG:H5'	1.91	0.52
2:C:757:ILE:HD12	2:C:837:LEU:HB2	1.91	0.52
2:C:1087:GLU:HG3	2:C:1091:ILE:HD11	1.92	0.52
3:D:238:GLU:OE2	3:D:242:ARG:NH1	2.41	0.52
5:F:515:ARG:O	5:F:519:ARG:N	2.43	0.52
3:D:905:ALA:HB2	3:D:911:ILE:HG13	1.93	0.51
5:F:489:LEU:H	5:F:489:LEU:HD23	1.75	0.51
2:C:317:ASN:O	2:C:321:GLY:N	2.37	0.51
5:F:401:LYS:HA	5:F:405:ILE:HA	1.92	0.51
1:A:129:ASN:ND2	2:C:652:GLU:HG3	2.26	0.51
3:D:365:ILE:HG21	5:F:297:GLU:HG2	1.91	0.51
1:B:92:PRO:HB3	1:B:141:GLU:HG2	1.92	0.50
2:C:721:VAL:HG23	2:C:915:ILE:HG23	1.92	0.50
3:D:778:TRP:CE2	3:D:835:PRO:HG3	2.46	0.50
3:D:834:ARG:NH2	3:D:847:LEU:O	2.40	0.50
5:F:470:ARG:HB3	5:F:506:ILE:HG21	1.94	0.50
5:F:505:GLN:HG3	5:F:509:LYS:HE3	1.92	0.50
1:A:64:THR:OG1	1:A:65:THR:N	2.42	0.50
2:C:809:LYS:HD2	2:C:833:ARG:HD3	1.91	0.50
1:A:21:PHE:HZ	1:A:205:ARG:HA	1.76	0.50
2:C:88:GLU:CD	2:C:310:ARG:HH12	2.19	0.50
2:C:451:HIS:O	2:C:451:HIS:ND1	2.42	0.50
5:F:474:VAL:HA	5:F:477:LEU:HD12	1.93	0.50
2:C:944:TRP:NE1	2:C:963:LEU:O	2.40	0.50
2:C:206:PRO:HA	2:C:308:LEU:HD23	1.93	0.49
6:H:20:DG:H1	7:G:8:DC:H42	1.59	0.49
2:C:641:VAL:HG12	2:C:705:GLY:H	1.77	0.49
1:A:175:THR:OG1	1:A:176:TYR:N	2.46	0.49
3:D:895:ARG:NH1	3:D:967:THR:O	2.46	0.49
1:A:87:SER:O	1:A:142:ARG:NH1	2.39	0.49
2:C:1088:LEU:HD23	2:C:1092:LYS:HD2	1.94	0.49
5:F:392:ARG:NH1	5:F:398:GLU:OE2	2.45	0.49
5:F:313:ARG:HB2	5:F:359:MET:HE2	1.94	0.49
3:D:915:TYR:HA	3:D:1143:ARG:HH12	1.78	0.49
3:D:756:VAL:HG21	3:D:777:ILE:HD11	1.95	0.49
2:C:441:ASP:HB2	2:C:451:HIS:HD2	1.78	0.49
2:C:107:PHE:HE1	2:C:418:ILE:HD11	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:896:GLY:HA2	3:D:431:VAL:HG13	1.95	0.49
2:C:919:THR:HG23	3:D:731:VAL:HG23	1.94	0.49
2:C:840:PRO:HB2	2:C:843:GLU:HG3	1.94	0.48
3:D:642:PRO:HG2	3:D:647:GLU:HB2	1.94	0.48
3:D:1030:ARG:HH21	3:D:1137:GLU:HG2	1.78	0.48
1:A:225:LEU:HD13	1:A:225:LEU:H	1.78	0.48
2:C:46:GLU:N	2:C:47:PRO:HD3	2.28	0.48
2:C:641:VAL:HG11	2:C:708:THR:HG21	1.96	0.48
3:D:127:LYS:O	3:D:133:ALA:N	2.46	0.48
3:D:230:ALA:N	3:D:233:GLN:OE1	2.46	0.48
6:H:19:DG:H2''	6:H:20:DG:C8	2.48	0.48
3:D:45:GLY:H	3:D:48:CYS:HB2	1.77	0.48
5:F:303:VAL:HG22	5:F:351:ILE:HD13	1.94	0.48
2:C:926:MET:HE1	3:D:817:LEU:HA	1.95	0.48
2:C:178:GLN:HG2	2:C:180:VAL:HG13	1.94	0.48
2:C:475:VAL:O	3:D:854:HIS:ND1	2.37	0.48
3:D:67:ARG:HD2	3:D:69:ARG:NE	2.29	0.48
2:C:604:ARG:NH1	2:C:925:ARG:HD2	2.28	0.48
5:F:477:LEU:HB3	5:F:492:ILE:HD13	1.96	0.48
1:A:120:ASN:OD1	1:A:120:ASN:N	2.46	0.48
2:C:371:ASP:OD1	6:H:14:DG:N1	2.47	0.48
4:E:33:LEU:HD23	4:E:33:LEU:H	1.79	0.48
2:C:232:GLN:OE1	2:C:280:LYS:HG3	2.14	0.48
2:C:313:ARG:HH22	2:C:337:ASP:CG	2.21	0.47
2:C:604:ARG:HG3	2:C:925:ARG:HB3	1.94	0.47
3:D:745:ILE:HD13	3:D:784:GLU:HG2	1.95	0.47
3:D:595:ASP:HB3	3:D:631:ALA:HB2	1.97	0.47
3:D:922:ALA:HB3	3:D:1150:HIS:CE1	2.49	0.47
2:C:185:VAL:HG12	2:C:316:VAL:HG22	1.97	0.47
3:D:1166:THR:HB	3:D:1206:VAL:HG21	1.95	0.47
3:D:103:HIS:CE1	3:D:105:TRP:HB2	2.49	0.47
2:C:344:TYR:OH	2:C:365:VAL:HA	2.15	0.47
2:C:373:PHE:HA	2:C:376:ARG:HB3	1.96	0.47
3:D:821:LYS:HB3	3:D:836:VAL:HB	1.95	0.47
2:C:257:ILE:HD11	2:C:346:VAL:HG23	1.96	0.47
2:C:446:LEU:HB2	2:C:713:MET:HE2	1.95	0.47
2:C:557:PRO:O	2:C:573:SER:N	2.48	0.47
2:C:599:HIS:HB3	2:C:928:ILE:HD12	1.97	0.47
3:D:49:GLU:OE2	3:D:55:THR:N	2.37	0.47
5:F:460:LEU:HA	5:F:463:VAL:HB	1.97	0.47
3:D:473:LYS:HD2	5:F:448:VAL:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:320:ILE:HG23	3:D:325:ARG:HD3	1.96	0.47
6:H:18:DC:H2"	6:H:19:DG:C8	2.49	0.47
2:C:599:HIS:ND1	3:D:840:PHE:O	2.48	0.47
2:C:805:LYS:HE3	2:C:835:THR:HB	1.97	0.47
3:D:924:THR:HG22	3:D:943:ASP:HA	1.96	0.47
2:C:677:ARG:HE	2:C:753:GLU:HA	1.80	0.46
3:D:899:VAL:HG11	3:D:920:ALA:HB2	1.97	0.46
5:F:256:GLY:HA2	5:F:284:ILE:HG22	1.97	0.46
2:C:1044:ARG:NH1	2:C:1048:PRO:HD2	2.31	0.46
3:D:127:LYS:HA	3:D:132:ALA:HB3	1.97	0.46
3:D:139:VAL:HG12	3:D:231:PRO:HD3	1.96	0.46
2:C:597:LEU:HB3	2:C:976:VAL:HG13	1.98	0.46
3:D:113:ARG:H	3:D:113:ARG:HD2	1.80	0.46
1:B:84:VAL:HG12	1:B:199:LYS:HD3	1.98	0.46
3:D:823:LEU:HD13	3:D:831:PHE:HB3	1.98	0.46
3:D:828:LYS:HE3	3:D:830:GLU:HB2	1.96	0.46
5:F:299:ASN:OD1	6:H:6:DT:N3	2.39	0.46
6:H:11:DG:H5"	6:H:12:DC:C4	2.51	0.46
2:C:476:HIS:CG	2:C:477:PRO:HD2	2.50	0.46
2:C:549:ASP:HB3	2:C:553:ARG:H	1.79	0.46
3:D:565:ILE:HG23	3:D:575:ALA:HB3	1.98	0.46
6:H:15:DT:H2"	6:H:16:DC:H6	1.79	0.46
2:C:815:THR:HG21	5:F:453:PHE:HE1	1.80	0.46
2:C:230:ARG:HH12	7:G:8:DC:H5"	1.81	0.46
2:C:281:LEU:HD22	2:C:282:ARG:HG3	1.98	0.46
2:C:53:LEU:HD13	2:C:449:LEU:HD21	1.98	0.45
2:C:821:LEU:HD22	5:F:456:LEU:HD11	1.98	0.45
2:C:1054:GLN:NE2	5:F:441:ASP:OD2	2.49	0.45
3:D:905:ALA:HB3	3:D:908:GLY:O	2.16	0.45
3:D:459:ARG:NH2	4:E:88:GLN:HG2	2.31	0.45
3:D:1065:THR:HG23	3:D:1076:VAL:HB	1.97	0.45
3:D:1139:GLN:O	3:D:1143:ARG:HG2	2.16	0.45
2:C:173:ARG:NH1	2:C:437:SER:O	2.50	0.45
2:C:516:TYR:HD2	2:C:531:LEU:HD13	1.82	0.45
3:D:365:ILE:HD11	5:F:234:GLN:HB3	1.98	0.45
3:D:384:ASN:HB2	3:D:401:SER:HB3	1.97	0.45
3:D:907:ASP:OD1	3:D:907:ASP:N	2.48	0.45
2:C:396:MET:HE3	2:C:396:MET:HB2	1.77	0.45
3:D:101:VAL:HG23	3:D:375:GLN:CD	2.41	0.45
5:F:432:ASP:OD1	5:F:433:SER:N	2.50	0.45
2:C:568:VAL:HG21	3:D:847:LEU:HD23	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:503:ILE:HA	5:F:506:ILE:HD12	1.99	0.45
1:B:44:SER:O	1:B:144:ARG:HB3	2.17	0.45
2:C:1007:LYS:HA	2:C:1024:THR:HA	1.99	0.45
5:F:403:MET:HG3	5:F:404:ASP:N	2.32	0.45
2:C:347:ARG:HH11	2:C:352:GLN:HE22	1.65	0.45
1:A:65:THR:HG22	1:A:72:ASP:HB3	1.99	0.45
3:D:392:THR:HB	3:D:396:ASN:HA	1.99	0.45
3:D:416:ASN:O	3:D:420:LYS:HG2	2.17	0.45
3:D:717:LYS:HB3	3:D:717:LYS:HE2	1.67	0.45
5:F:502:ARG:O	5:F:506:ILE:HG13	2.17	0.45
2:C:251:ARG:NH2	2:C:343:GLU:OE1	2.42	0.44
2:C:815:THR:HG21	5:F:453:PHE:CE1	2.52	0.44
5:F:253:ILE:O	5:F:257:LEU:N	2.37	0.44
2:C:737:LEU:HG	2:C:895:ILE:HD12	2.00	0.44
3:D:707:ILE:HD11	4:E:32:PRO:HB3	1.99	0.44
1:B:208:LEU:O	1:B:212:GLY:N	2.49	0.44
3:D:876:ARG:HG2	3:D:1226:PHE:HZ	1.82	0.44
2:C:315:LYS:HD2	2:C:315:LYS:HA	1.81	0.44
2:C:762:THR:HG23	2:C:764:LEU:H	1.83	0.44
3:D:589:THR:HB	3:D:687:GLN:HA	2.00	0.44
3:D:891:CYS:O	3:D:892:GLN:HB2	2.18	0.44
4:E:95:ALA:O	4:E:98:GLU:HG2	2.17	0.44
1:B:110:ILE:O	1:B:112:PRO:HD3	2.17	0.44
2:C:197:LYS:HD2	2:C:218:LYS:HA	2.00	0.44
2:C:305:ARG:HH12	6:H:10:DA:N6	2.16	0.44
2:C:546:SER:HA	2:C:547:PRO:HD3	1.88	0.44
2:C:1020:PRO:HB2	2:C:1021:TYR:CD2	2.53	0.44
3:D:445:LYS:NZ	3:D:518:GLU:OE2	2.44	0.44
5:F:344:SER:OG	6:H:5:DA:OP2	2.30	0.44
5:F:423:LEU:HA	5:F:435:LEU:HD23	1.99	0.44
1:A:185:GLN:HG2	1:A:186:ARG:H	1.82	0.44
3:D:116:TYR:O	3:D:295:ARG:NE	2.44	0.44
3:D:1218:ASP:OD1	3:D:1218:ASP:N	2.51	0.44
1:A:100:GLN:HG2	1:A:101:GLY:H	1.83	0.43
2:C:377:ARG:NH2	2:C:383:GLU:OE1	2.51	0.43
2:C:549:ASP:OD2	2:C:550:ALA:N	2.50	0.43
2:C:818:GLU:OE2	2:C:822:ARG:NH1	2.49	0.43
3:D:308:SER:HA	3:D:309:PRO:HD3	1.80	0.43
3:D:558:LEU:HB2	3:D:559:MET:HG3	2.00	0.43
2:C:111:ARG:NH2	2:C:151:THR:OG1	2.51	0.43
3:D:445:LYS:HB3	3:D:484:TRP:CZ3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:740:PRO:HD3	3:D:792:HIS:ND1	2.33	0.43
5:F:345:THR:HA	6:H:5:DA:N7	2.33	0.43
2:C:253:GLY:HA2	2:C:259:ARG:HE	1.83	0.43
3:D:888:GLU:HG2	3:D:889:HIS:H	1.83	0.43
5:F:403:MET:HE3	5:F:403:MET:HB2	1.92	0.43
2:C:408:ASP:O	2:C:412:ILE:HG13	2.18	0.43
2:C:206:PRO:HB3	2:C:306:TYR:CZ	2.54	0.43
2:C:522:GLY:O	2:C:553:ARG:HA	2.19	0.43
3:D:369:ASN:O	3:D:373:MET:HG3	2.18	0.43
3:D:409:LYS:O	3:D:415:GLN:HB2	2.18	0.43
3:D:1088:VAL:HG11	3:D:1109:GLN:CD	2.44	0.43
2:C:157:PHE:HA	2:C:158:PRO:HD3	1.92	0.43
2:C:377:ARG:HH22	2:C:383:GLU:CD	2.27	0.43
2:C:615:ALA:HB3	2:C:715:LEU:HD22	2.01	0.43
2:C:809:LYS:N	2:C:831:GLU:O	2.42	0.43
3:D:36:TYR:CZ	3:D:37:ARG:HG3	2.53	0.43
3:D:350:ARG:NH1	3:D:373:MET:HB3	2.33	0.43
1:A:40:ARG:HB2	1:B:33:THR:HG23	2.01	0.43
3:D:832:ILE:HG22	3:D:834:ARG:H	1.83	0.43
2:C:220:ASP:HB3	2:C:257:ILE:HG22	2.00	0.43
3:D:778:TRP:CD2	3:D:835:PRO:HG3	2.53	0.43
5:F:342:LYS:HE2	5:F:342:LYS:HB3	1.86	0.43
3:D:580:ASP:OD1	3:D:580:ASP:N	2.52	0.43
5:F:383:GLN:O	5:F:387:LEU:HB2	2.19	0.43
1:B:95:MET:HE3	1:B:140:VAL:HG21	2.01	0.42
2:C:152:VAL:HG21	2:C:418:ILE:HD12	2.01	0.42
2:C:200:HIS:CE1	2:C:348:LEU:HG	2.53	0.42
2:C:485:PRO:HG3	2:C:586:MET:HE3	2.01	0.42
2:C:1052:ILE:O	3:D:89:ARG:NH1	2.52	0.42
5:F:329:ILE:O	5:F:333:GLU:HG2	2.19	0.42
5:F:372:MET:O	5:F:376:ILE:HG13	2.19	0.42
3:D:1273:GLN:O	4:E:105:GLU:N	2.39	0.42
5:F:485:GLN:HA	5:F:486:PRO:HD3	1.82	0.42
5:F:502:ARG:HA	5:F:502:ARG:HD3	1.82	0.42
1:B:84:VAL:HG23	1:B:119:HIS:HB2	2.00	0.42
2:C:472:VAL:HG22	6:H:14:DG:N2	2.34	0.42
2:C:759:ALA:HB1	2:C:767:GLU:O	2.19	0.42
3:D:1127:PRO:O	3:D:1130:VAL:HG12	2.19	0.42
5:F:467:LEU:HG	5:F:510:THR:HG23	2.00	0.42
1:B:182:ARG:HA	1:B:187:THR:HA	2.01	0.42
1:B:77:ILE:HG22	1:B:81:LYS:HE3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:321:PRO:HA	3:D:322:PRO:HD3	1.77	0.42
3:D:676:LEU:HD23	3:D:716:LEU:HD23	2.02	0.42
2:C:356:THR:HG22	2:C:362:GLU:HA	2.00	0.42
2:C:958:ARG:N	2:C:958:ARG:HD2	2.34	0.42
2:C:600:ASP:O	3:D:849:TYR:OH	2.34	0.42
3:D:789:LEU:HD22	3:D:793:TYR:CE2	2.54	0.42
1:A:181:THR:O	1:A:188:ASP:HA	2.20	0.42
1:B:157:ALA:C	1:B:159:ILE:H	2.28	0.42
2:C:653:VAL:HG23	2:C:658:ILE:HG12	2.02	0.42
2:C:1044:ARG:HH12	2:C:1048:PRO:HD2	1.85	0.42
2:C:252:PHE:HB3	2:C:258:MET:HG3	2.01	0.42
2:C:687:CYS:HA	2:C:688:PRO:HD3	1.83	0.42
3:D:113:ARG:HD2	3:D:113:ARG:N	2.34	0.42
2:C:38:ARG:NH2	2:C:971:ILE:HB	2.35	0.42
3:D:35:ASN:OD1	3:D:36:TYR:N	2.52	0.42
3:D:87:VAL:C	3:D:89:ARG:H	2.28	0.42
2:C:371:ASP:OD1	6:H:14:DG:N2	2.51	0.41
2:C:1141:ASP:OD1	2:C:1142:GLY:N	2.47	0.41
3:D:69:ARG:HD2	3:D:70:PHE:CZ	2.55	0.41
3:D:409:LYS:NZ	7:G:15:DT:OP2	2.48	0.41
3:D:454:PRO:HA	3:D:457:MET:HE2	2.02	0.41
5:F:262:LEU:O	5:F:266:LEU:HG	2.20	0.41
1:B:83:LEU:HB2	1:B:123:MET:HE1	2.00	0.41
2:C:419:ASN:O	2:C:422:PRO:HD2	2.20	0.41
2:C:848:ILE:HD13	2:C:874:ALA:HB2	2.02	0.41
3:D:125:LEU:HD12	3:D:125:LEU:HA	1.93	0.41
3:D:923:ARG:NH2	3:D:1155:GLU:OE2	2.35	0.41
3:D:1003:ILE:HA	3:D:1149:ILE:HD13	2.01	0.41
4:E:84:GLU:CD	4:E:84:GLU:H	2.27	0.41
2:C:182:SER:HB2	2:C:377:ARG:HB2	2.02	0.41
2:C:1067:ARG:NH2	3:D:415:GLN:O	2.46	0.41
3:D:1176:LEU:H	3:D:1176:LEU:HD12	1.85	0.41
5:F:384:ARG:HA	5:F:387:LEU:HB2	2.02	0.41
2:C:1078:ALA:HB1	3:D:998:VAL:HG22	2.01	0.41
3:D:48:CYS:SG	3:D:50:LYS:HB3	2.60	0.41
3:D:153:ALA:O	3:D:157:VAL:HG23	2.21	0.41
3:D:293:LEU:HD23	3:D:293:LEU:HA	1.96	0.41
3:D:14:LEU:HD13	3:D:106:TYR:OH	2.20	0.41
3:D:453:LYS:O	3:D:457:MET:HG3	2.20	0.41
3:D:760:PHE:CG	3:D:770:ARG:HD2	2.54	0.41
3:D:1036:GLU:OE2	3:D:1211:THR:OG1	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:LYS:HE2	1:A:131:LYS:HB3	1.95	0.41
3:D:34:ILE:HG22	3:D:41:PRO:HA	2.02	0.41
5:F:515:ARG:HA	5:F:518:SER:HB3	2.02	0.41
2:C:215:ASP:HB2	2:C:231:ARG:NH2	2.35	0.41
2:C:544:ALA:HB2	2:C:580:ASP:HB2	2.03	0.41
2:C:546:SER:O	2:C:548:ILE:HG13	2.21	0.41
2:C:894:VAL:HG22	3:D:536:PHE:O	2.20	0.41
2:C:1131:LEU:HD23	3:D:382:PHE:CD1	2.56	0.41
3:D:600:GLN:HB2	3:D:609:THR:HB	2.02	0.41
3:D:1030:ARG:NH2	3:D:1034:LEU:HD21	2.36	0.41
3:D:1220:TRP:CD1	3:D:1243:ASP:HB2	2.55	0.41
1:B:97:LEU:HD22	1:B:110:ILE:HG12	2.03	0.41
2:C:150:GLN:HG2	2:C:414:PRO:HG2	2.03	0.41
3:D:329:GLN:HB2	3:D:335:PHE:CE1	2.55	0.41
3:D:497:LEU:O	3:D:543:VAL:HA	2.21	0.41
3:D:666:THR:HG21	3:D:683:PHE:CE2	2.55	0.41
3:D:1248:LEU:HD23	3:D:1259:PRO:HD2	2.03	0.41
1:B:27:GLU:O	1:B:30:PHE:HB2	2.20	0.41
2:C:544:ALA:HA	2:C:579:MET:HE2	2.03	0.41
2:C:1114:GLU:HA	2:C:1115:PRO:HD3	1.92	0.41
2:C:1126:LYS:HA	2:C:1126:LYS:HD3	1.86	0.41
3:D:106:TYR:CD2	3:D:312:MET:HG2	2.56	0.41
3:D:337:THR:HG1	3:D:341:ASN:HD22	1.67	0.41
3:D:389:ARG:HA	3:D:390:PRO:HD3	1.86	0.41
3:D:500:ARG:HB2	3:D:541:MET:HE2	2.03	0.41
3:D:556:ARG:HG3	4:E:35:ILE:HG12	2.02	0.41
3:D:953:LEU:O	3:D:953:LEU:HD12	2.20	0.41
3:D:1168:ILE:HB	3:D:1202:ALA:HB3	2.03	0.41
1:A:18:ARG:HH21	1:A:195:ASP:CG	2.28	0.41
1:A:129:ASN:OD1	1:A:130:ASP:N	2.49	0.41
2:C:421:ARG:HB3	2:C:422:PRO:HD3	2.03	0.41
1:B:148:PRO:O	1:B:152:ASN:ND2	2.53	0.40
2:C:211:TRP:HH2	6:H:14:DG:OP1	2.04	0.40
2:C:658:ILE:HG21	2:C:702:ILE:HD12	2.03	0.40
3:D:67:ARG:CZ	5:F:486:PRO:HD3	2.51	0.40
3:D:76:GLU:CD	3:D:76:GLU:H	2.29	0.40
3:D:114:LEU:HA	3:D:114:LEU:HD23	1.79	0.40
5:F:242:ASN:HB3	5:F:245:GLU:HG3	2.02	0.40
2:C:815:THR:O	2:C:819:ARG:N	2.40	0.40
5:F:387:LEU:HD12	5:F:387:LEU:HA	1.94	0.40
3:D:400:LYS:HA	3:D:400:LYS:HD3	1.90	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1125:GLN:HB3	3:D:1129:GLU:HG2	2.02	0.40
7:G:15:DT:H2'	7:G:16:DC:C6	2.57	0.40
1:B:90:ASP:HA	1:B:142:ARG:HD3	2.02	0.40
2:C:519:VAL:HG22	2:C:524:VAL:HA	2.03	0.40
2:C:1057:LEU:HD11	5:F:440:GLU:HA	2.04	0.40
3:D:339:ASP:OD1	5:F:422:SER:OG	2.25	0.40
3:D:584:GLY:O	3:D:587:TYR:HB3	2.22	0.40
3:D:1208:MET:HE3	3:D:1208:MET:HB2	1.93	0.40
4:E:40:ILE:O	4:E:44:LEU:HG	2.21	0.40
5:F:470:ARG:NH1	5:F:506:ILE:HD11	2.35	0.40
5:F:474:VAL:HA	5:F:477:LEU:HB2	2.03	0.40
1:B:95:MET:HB3	1:B:113:PRO:HD3	2.03	0.40
2:C:314:TYR:CZ	2:C:318:LYS:HD3	2.57	0.40
2:C:720:LEU:HD23	2:C:913:VAL:HA	2.02	0.40
3:D:724:ALA:O	3:D:727:SER:OG	2.26	0.40
3:D:793:TYR:HB3	3:D:800:ILE:HG13	2.02	0.40
5:F:302:LEU:O	5:F:305:SER:OG	2.30	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	222/347 (64%)	209 (94%)	12 (5%)	1 (0%)	24	62
1	B	225/347 (65%)	207 (92%)	15 (7%)	3 (1%)	9	41
2	C	1124/1178 (95%)	1043 (93%)	75 (7%)	6 (0%)	24	62
3	D	1261/1316 (96%)	1202 (95%)	53 (4%)	6 (0%)	24	62
4	E	79/110 (72%)	76 (96%)	3 (4%)	0	100	100
5	F	320/528 (61%)	306 (96%)	12 (4%)	2 (1%)	21	58
All	All	3231/3826 (84%)	3043 (94%)	170 (5%)	18 (1%)	21	58

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	678	PRO
1	A	184	GLU
3	D	971	SER
2	C	47	PRO
2	C	564	LYS
2	C	1134	ASN
1	B	35	GLY
1	B	227	VAL
2	C	33	PRO
2	C	922	VAL
3	D	658	PRO
3	D	703	ARG
5	F	486	PRO
3	D	607	PRO
5	F	405	ILE
1	B	67	PRO
2	C	32	VAL
3	D	593	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	192/297 (65%)	188 (98%)	4 (2%)	47	65
1	B	192/297 (65%)	192 (100%)	0	100	100
2	C	948/998 (95%)	920 (97%)	28 (3%)	36	57
3	D	1048/1095 (96%)	1023 (98%)	25 (2%)	43	63
4	E	68/90 (76%)	66 (97%)	2 (3%)	37	58
5	F	271/427 (64%)	262 (97%)	9 (3%)	33	55
All	All	2719/3204 (85%)	2651 (98%)	68 (2%)	42	62

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	136	VAL
1	A	159	ILE
1	A	225	LEU
2	C	44	LEU
2	C	48	LEU
2	C	80	VAL
2	C	185	VAL
2	C	224	VAL
2	C	257	ILE
2	C	311	VAL
2	C	322	LEU
2	C	332	THR
2	C	373	PHE
2	C	378	LEU
2	C	506	VAL
2	C	512	ILE
2	C	588	SER
2	C	621	SER
2	C	647	SER
2	C	649	VAL
2	C	691	ASP
2	C	710	ASP
2	C	727	GLU
2	C	835	THR
2	C	850	ILE
2	C	854	SER
2	C	858	GLU
2	C	916	ILE
2	C	928	ILE
2	C	1002	VAL
2	C	1079	TYR
3	D	60	CYS
3	D	75	CYS
3	D	82	VAL
3	D	83	THR
3	D	101	VAL
3	D	114	LEU
3	D	261	ILE
3	D	447	MET
3	D	582	VAL
3	D	583	THR
3	D	589	THR

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Mol	Chain	Res	Type
3	D	629	VAL
3	D	684	VAL
3	D	867	THR
3	D	901	LEU
3	D	910	LEU
3	D	930	VAL
3	D	968	CYS
3	D	978	CYS
3	D	1054	ARG
3	D	1098	VAL
3	D	1099	LEU
3	D	1194	VAL
3	D	1240	CYS
3	D	1249	LYS
4	E	48	SER
4	E	75	ILE
5	F	269	ARG
5	F	271	GLU
5	F	367	ARG
5	F	434	GLN
5	F	448	VAL
5	F	463	VAL
5	F	485	GLN
5	F	503	ILE
5	F	527	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	HIS
1	B	200	ASN
2	C	37	ASN
2	C	232	GLN
2	C	349	HIS
2	C	969	ASN
2	C	981	GLN
2	C	1066	GLN
3	D	813	GLN
3	D	1246	ASN
4	E	37	ASN
4	E	69	ASN
4	E	70	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	I	3/4 (75%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 3 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	224/347 (64%)	0.02	0 100 100	21, 62, 124, 153	0
1	B	227/347 (65%)	0.25	1 (0%) 88 77	54, 100, 146, 167	0
2	C	1126/1178 (95%)	-0.05	4 (0%) 88 77	5, 43, 126, 147	0
3	D	1265/1316 (96%)	-0.16	4 (0%) 90 80	4, 40, 109, 148	0
4	E	81/110 (73%)	-0.02	1 (1%) 76 61	20, 48, 85, 101	0
5	F	322/528 (60%)	0.16	5 (1%) 70 54	11, 69, 170, 191	0
6	H	23/23 (100%)	0.25	0 100 100	34, 93, 168, 189	0
7	G	16/16 (100%)	0.29	0 100 100	30, 59, 186, 194	0
8	I	4/4 (100%)	-0.57	0 100 100	30, 44, 46, 51	0
All	All	3288/3869 (84%)	-0.04	15 (0%) 87 75	4, 50, 130, 194	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	F	481	LEU	3.5
3	D	1194	VAL	2.9
2	C	820	LEU	2.9
5	F	503	ILE	2.6
3	D	152	GLU	2.5
3	D	1281	ALA	2.5
1	B	225	LEU	2.4
3	D	5	ASN	2.4
5	F	405	ILE	2.2
4	E	75	ILE	2.2
5	F	489	LEU	2.2
2	C	979	GLY	2.1
2	C	1127	GLU	2.1
5	F	309	ARG	2.0
2	C	288	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	ZN	D	2001	1/1	0.95	0.04	32,32,32,32	0
9	ZN	D	2002	1/1	1.00	0.01	35,35,35,35	0
10	MG	D	2003	1/1	1.00	0.02	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.