



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:14 AM UTC

PDB ID : 6QIC / pdb_00006qic
Title : Crystal structure of DEAH-box ATPase Prp22-S837A with bound ssRNA
Authors : Hamann, F.; Ficner, R.; Enders, M.
Deposited on : 2019-01-18
Resolution : 2.70 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

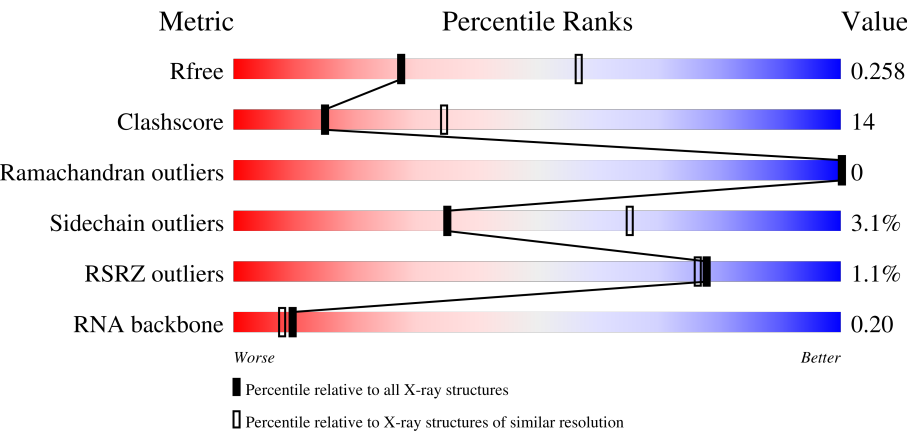
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.



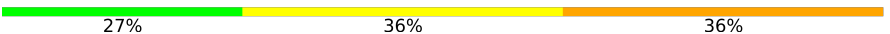
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)
RNA backbone	3983	1044 (2.90-2.50)

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	677	% 70%21%• 7%
1	B	677	2% 67%25%• 8%
1	C	677	% 73%19%• 7%
1	D	677	69%22%• 7%

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Mol	Chain	Length	Quality of chain
2	E	11	 18% 55% 27%
2	F	11	 27% 36% 36%
2	G	11	 55% 36% 9%
2	H	11	 9% 73% 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20594 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 Putative pre-mRNA splicing factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	628	Total	C	N	O	S	0	0	0
			4932	3146	831	926	29			
1	B	626	Total	C	N	O	S	0	0	0
			4934	3147	834	923	30			
1	C	627	Total	C	N	O	S	0	0	0
			4942	3154	834	925	29			
1	D	629	Total	C	N	O	S	0	1	0
			4953	3159	836	928	30			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	837	ALA	SER	engineered mutation	UNP G0S700
B	837	ALA	SER	engineered mutation	UNP G0S700
C	837	ALA	SER	engineered mutation	UNP G0S700
D	837	ALA	SER	engineered mutation	UNP G0S700

- Molecule 2 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	11	Total	C	N	O	P	0	0	0
			220	99	22	88	11			
2	F	11	Total	C	N	O	P	0	0	0
			204	90	20	83	11			
2	G	11	Total	C	N	O	P	0	0	0
			204	90	20	83	11			
2	H	11	Total	C	N	O	P	0	0	0
			204	90	20	83	11			

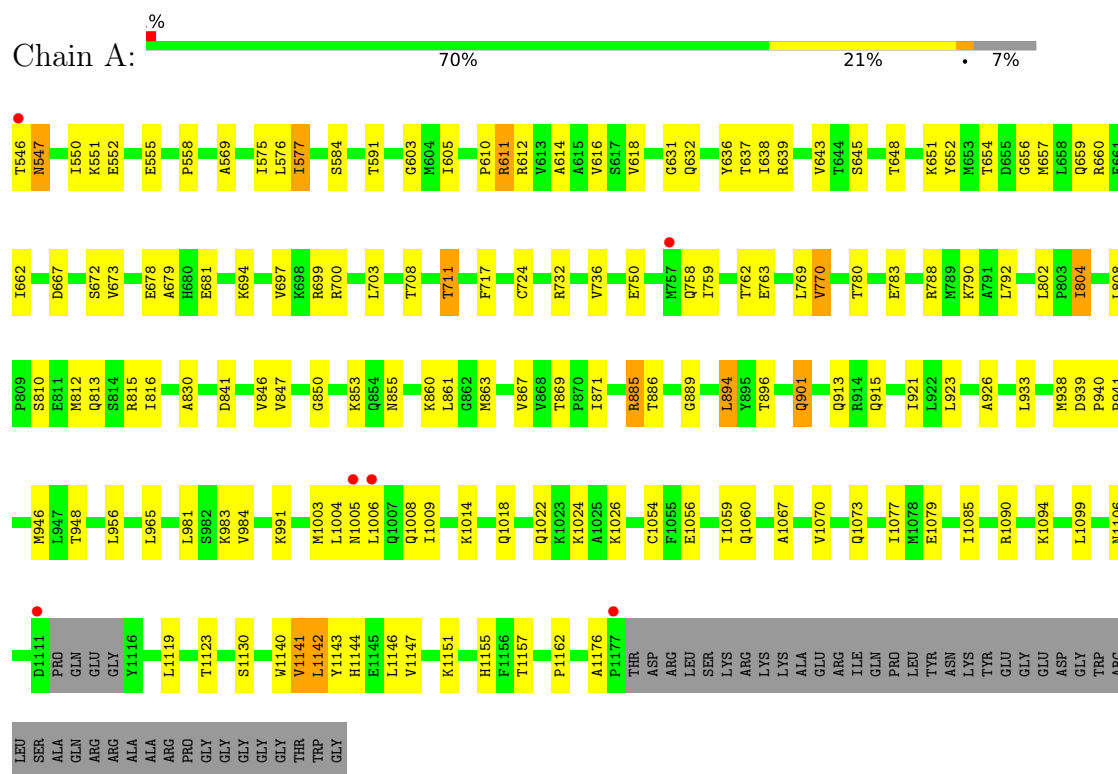
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		

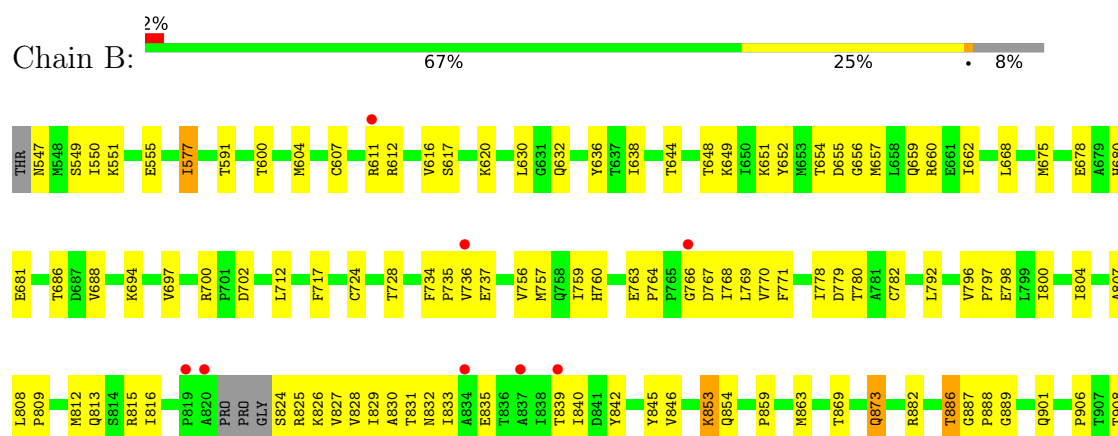
3 Residue-property plots [i](#)

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 electron 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

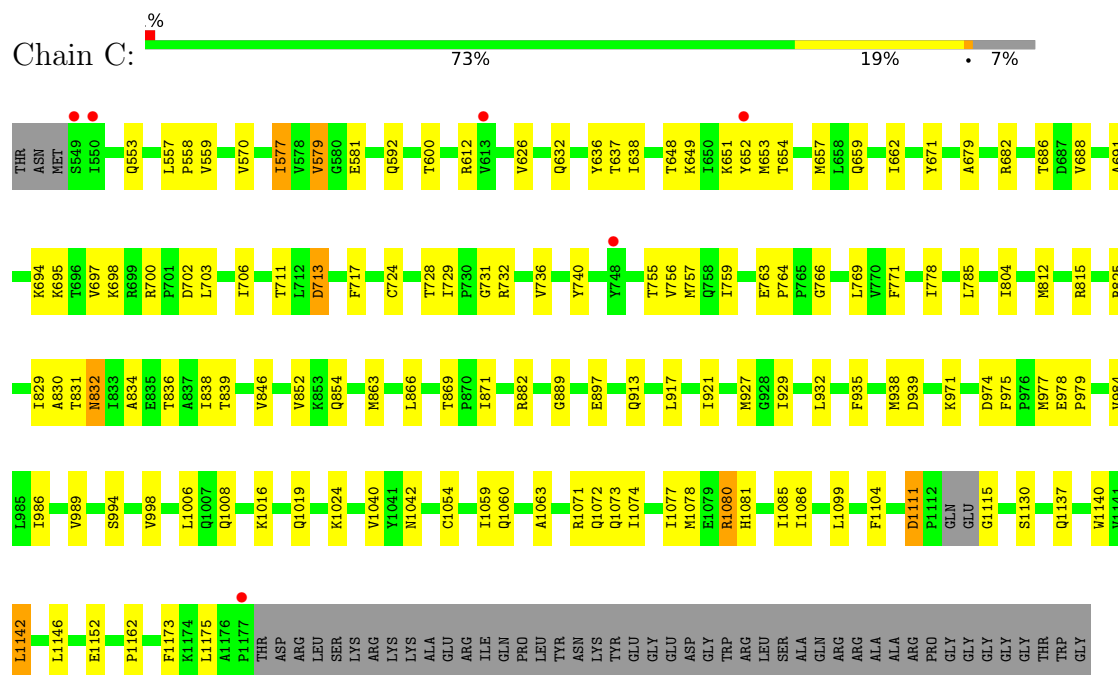
• Molecule 1: Putative pre-mRNA splicing factor



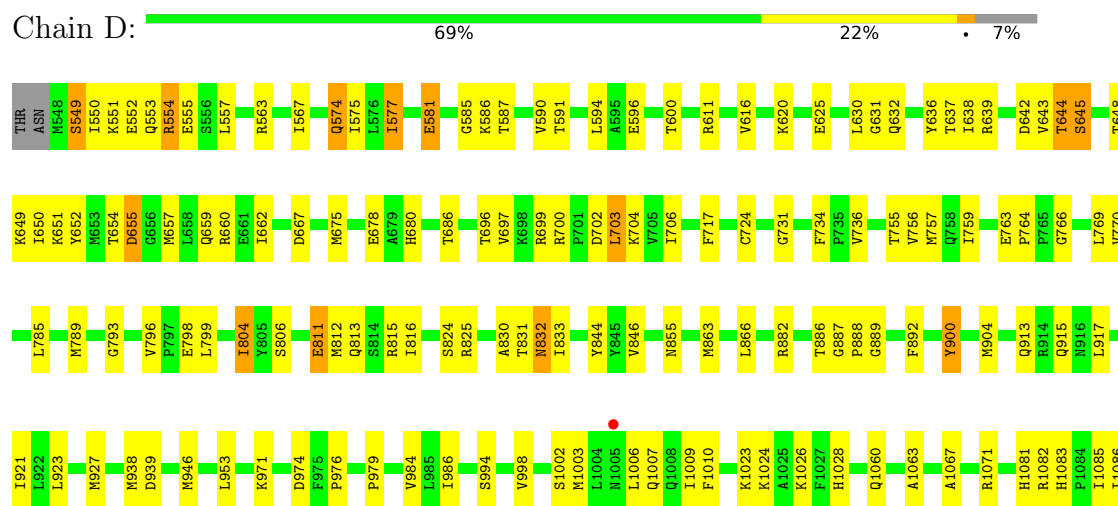
• Molecule 1: Putative pre-mRNA splicing factor

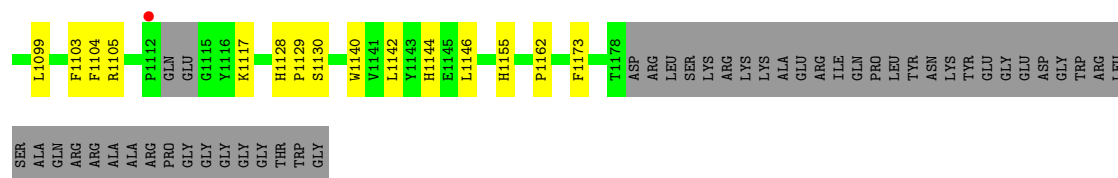


- Molecule 1: Putative pre-mRNA splicing factor



- Molecule 1: Putative pre-mRNA splicing factor





- Molecule 2: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

Chain E: 18% 55% 27%



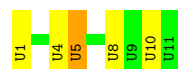
- Molecule 2: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

Chain F: 27% 36% 36%



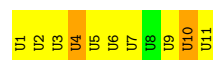
- Molecule 2: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

Chain G: 55% 36% 9%



- Molecule 2: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*UP*UP*UP*U)-3')

Chain H: 9% 73% 18%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	139.61Å 139.96Å 164.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	106.49 – 2.70 106.49 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (106.49-2.70) 99.7 (106.49-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.212 , 0.261 0.213 , 0.258	Depositor DCC
R_{free} test set	4442 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	77.2	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20594	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.67% 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](#)

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	1.04	2/5035 (0.0%)	1.42	4/6835 (0.1%)
1	B	1.05	1/5035 (0.0%)	1.41	4/6825 (0.1%)
1	C	1.04	1/5046 (0.0%)	1.41	4/6845 (0.1%)
1	D	1.05	0/5059	1.43	7/6863 (0.1%)
2	E	0.49	0/241	0.64	0/370
2	F	0.48	0/223	0.68	0/343
2	G	0.47	0/223	0.63	0/343
2	H	0.50	0/223	0.65	0/343
All	All	1.03	4/21085 (0.0%)	1.39	19/28767 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	859	PRO	C-O	-6.54	1.16	1.24
1	A	1144	HIS	CG-ND1	5.71	1.44	1.38
1	C	1086	ILE	C-O	5.48	1.29	1.24
1	A	1003	MET	C-O	5.18	1.30	1.24

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	549	SER	CA-C-O	-6.82	114.22	121.99
1	D	549	SER	CA-C-N	6.03	128.16	120.56
1	D	549	SER	C-N-CA	6.03	128.16	120.56
1	D	1144	HIS	CA-CB-CG	-5.92	107.88	113.80
1	A	1144	HIS	CA-CB-CG	-5.83	107.97	113.80
1	B	873	GLN	CB-CA-C	5.72	120.28	110.79
1	D	611	ARG	CB-CG-CD	5.62	124.23	111.30
1	C	1042	ASN	OD1-CG-ND2	-5.54	117.06	122.60
1	D	655	ASP	CA-CB-CG	5.49	118.09	112.60
1	C	711	THR	CA-C-O	-5.33	115.22	120.82
1	C	989	VAL	N-CA-C	-5.22	105.76	110.82
1	C	713	ASP	CA-CB-CG	5.20	117.80	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	574	GLN	N-CA-CB	5.18	117.80	110.13
1	A	948	THR	CA-CB-OG1	-5.12	101.93	109.60
1	A	860	LYS	CB-CA-C	5.10	120.06	110.63
1	A	711	THR	CB-CA-C	5.08	118.86	110.88
1	D	574	GLN	CA-C-O	-5.07	115.12	120.55
1	B	728	THR	CA-C-N	5.04	127.23	122.85
1	B	728	THR	C-N-CA	5.04	127.23	122.85

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4932	0	4932	135	0
1	B	4934	0	4954	168	0
1	C	4942	0	4969	116	0
1	D	4953	0	4980	141	0
2	E	220	0	111	14	0
2	F	204	0	100	15	0
2	G	204	0	100	11	0
2	H	204	0	100	14	0
3	A	1	0	0	0	0
All	All	20594	0	20246	561	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 (561) 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)
1:C:975:PHE:CE2	1:C:1078:MET:HE3	1.79	1.18
1:B:1011:TYR:HB2	1:B:1060:GLN:HE21	1.09	1.18
1:B:798:GLU:HB3	1:B:824:SER:HB3	1.21	1.15
1:C:975:PHE:CZ	1:C:1078:MET:CE	2.30	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:703:LEU:HD12	1:D:704:LYS:N	1.64	1.10
1:D:657:MET:HE1	2:H:10:U:OP1	1.50	1.10
1:A:1006:LEU:HD11	1:A:1067:ALA:HB2	1.34	1.09
1:B:757:MET:HE1	1:B:792:LEU:HD11	1.34	1.08
1:A:603:GLY:HA3	1:A:672:SER:HB3	1.35	1.07
1:B:1011:TYR:HB2	1:B:1060:GLN:NE2	1.70	1.06
1:B:1020:ALA:HA	1:B:1058:TYR:CD2	1.92	1.04
1:D:581:GLU:OE1	1:D:731:GLY:HA3	1.59	1.03
1:B:1016:LYS:HG2	1:B:1019:GLN:OE1	1.60	1.01
1:C:755:THR:O	1:C:759:ILE:HG13	1.60	1.00
1:B:736:VAL:HA	1:B:889:GLY:O	1.58	1.00
1:C:975:PHE:HZ	1:C:1078:MET:CE	1.71	1.00
1:D:734:PHE:CB	1:D:887:GLY:C	2.34	0.99
1:C:804:ILE:CG2	1:C:830:ALA:HB2	1.94	0.97
1:B:804:ILE:CG2	1:B:830:ALA:HB2	1.96	0.96
1:B:1006:LEU:HD11	1:B:1067:ALA:HB2	1.44	0.95
1:A:605:ILE:HD13	1:A:673:VAL:HB	1.46	0.95
1:A:605:ILE:CD1	1:A:673:VAL:HB	1.97	0.94
1:A:1079:GLU:HG3	1:D:1082:ARG:HH22	1.34	0.91
1:D:717:PHE:CE1	1:D:938:MET:HE1	2.06	0.91
1:D:1023:LYS:O	1:D:1026:LYS:HG2	1.70	0.91
1:C:1080:ARG:HH11	1:C:1080:ARG:HG2	1.35	0.91
1:A:632:GLN:O	1:A:648:THR:HG22	1.70	0.90
1:C:1111:ASP:HB3	1:C:1115:GLY:O	1.71	0.90
1:B:981:LEU:HD21	1:B:1004:LEU:HD12	1.51	0.89
1:B:717:PHE:CE1	1:B:938:MET:HE1	2.07	0.89
1:B:632:GLN:O	1:B:648:THR:HG22	1.74	0.88
1:C:975:PHE:CE2	1:C:1078:MET:CE	2.54	0.87
1:C:581:GLU:HB2	1:C:731:GLY:HA3	1.58	0.86
1:A:1079:GLU:HG3	1:D:1082:ARG:NH2	1.90	0.85
1:B:981:LEU:HD21	1:B:1004:LEU:CD1	2.07	0.85
1:B:873:GLN:HG2	1:B:906:PRO:HA	1.59	0.84
1:C:975:PHE:HE2	1:C:1078:MET:HE3	1.29	0.84
1:B:1011:TYR:CB	1:B:1060:GLN:HE21	1.90	0.84
1:D:844:TYR:CE1	1:D:886:THR:O	2.30	0.84
1:A:1141:VAL:HG13	1:A:1157:THR:HG23	1.59	0.84
1:C:975:PHE:HZ	1:C:1078:MET:HE2	1.44	0.83
1:C:579:VAL:HG13	1:C:728:THR:HG22	1.60	0.83
1:B:970:ARG:HD2	1:B:970:ARG:O	1.78	0.82
1:A:603:GLY:CA	1:A:672:SER:HB3	2.10	0.81
1:B:779:ASP:C	1:B:1014:LYS:NZ	2.37	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1009:ILE:H	1:A:1009:ILE:HD12	1.44	0.80
1:B:798:GLU:O	1:B:824:SER:HB2	1.81	0.79
1:C:717:PHE:CE1	1:C:938:MET:HE1	2.18	0.79
1:A:1008:GLN:O	1:A:1060:GLN:HG3	1.83	0.79
1:C:577:ILE:HD12	1:C:724:CYS:SG	2.23	0.79
1:D:921:ILE:HD11	1:D:946:MET:HE3	1.65	0.79
1:B:798:GLU:O	1:B:824:SER:CB	2.32	0.78
1:D:577:ILE:HD12	1:D:724:CYS:SG	2.25	0.77
1:D:657:MET:CE	2:H:10:U:OP1	2.30	0.77
1:A:657:MET:HE1	2:E:10:U:OP1	1.85	0.77
1:C:657:MET:HE1	2:G:10:U:OP1	1.82	0.77
1:C:804:ILE:HG23	1:C:830:ALA:CB	2.13	0.77
1:A:547:ASN:HB3	1:A:550:ILE:HB	1.66	0.77
1:B:807:ALA:HB1	1:B:1008:GLN:OE1	1.85	0.77
1:D:831:THR:OG1	1:D:833:ILE:HG22	1.85	0.77
1:D:703:LEU:HD12	1:D:704:LYS:H	1.48	0.76
1:B:657:MET:HE1	2:F:10:U:OP1	1.85	0.76
1:D:734:PHE:CB	1:D:887:GLY:O	2.34	0.76
1:B:933:LEU:HD12	1:B:933:LEU:H	1.50	0.76
1:A:591:THR:HG21	1:A:652:TYR:OH	1.85	0.75
1:C:921:ILE:HG23	1:C:932:LEU:HD12	1.66	0.75
1:B:804:ILE:HG23	1:B:830:ALA:CB	2.17	0.75
1:A:717:PHE:CE1	1:A:938:MET:HE1	2.20	0.75
1:B:757:MET:CE	1:B:792:LEU:HD11	2.15	0.75
1:D:921:ILE:CD1	1:D:946:MET:HE3	2.16	0.75
1:A:901:GLN:HA	1:A:901:GLN:NE2	2.00	0.75
1:B:779:ASP:C	1:B:1014:LYS:HZ2	1.95	0.75
1:D:844:TYR:HE1	1:D:886:THR:O	1.70	0.74
1:C:921:ILE:HG23	1:C:932:LEU:CD1	2.15	0.74
1:B:717:PHE:CD1	1:B:938:MET:HE1	2.22	0.74
1:A:614:ALA:O	1:A:618:VAL:HG23	1.88	0.74
1:B:577:ILE:HD12	1:B:724:CYS:SG	2.27	0.74
1:B:591:THR:HG21	1:B:652:TYR:OH	1.88	0.74
1:B:766:GLY:O	1:B:825:ARG:HD3	1.87	0.74
1:A:1151:LYS:NZ	2:E:1:U:O2'	2.21	0.74
1:A:616:VAL:HG12	1:A:636:TYR:CE1	2.23	0.73
1:B:734:PHE:CE2	1:B:886:THR:O	2.41	0.73
1:B:853:LYS:HE3	2:F:5:U:OP1	1.89	0.73
1:A:894:LEU:HD23	1:A:894:LEU:N	2.04	0.72
1:D:1155:HIS:NE2	2:H:4:U:H5	1.87	0.72
1:D:717:PHE:CD1	1:D:938:MET:HE1	2.24	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:793:GLY:O	1:D:796:VAL:HG12	1.90	0.71
1:C:579:VAL:CG1	1:C:728:THR:HG22	2.19	0.71
1:B:1155:HIS:CE1	2:F:4:U:H5	2.09	0.71
1:B:757:MET:HE1	1:B:792:LEU:CD1	2.18	0.71
1:B:1015:ASP:N	1:B:1015:ASP:OD1	2.23	0.70
1:B:804:ILE:CG2	1:B:830:ALA:CB	2.69	0.70
1:D:1024:LYS:HE2	1:D:1028:HIS:NE2	2.05	0.70
1:D:1024:LYS:NZ	1:D:1130:SER:O	2.25	0.70
1:A:1024:LYS:NZ	1:A:1130:SER:O	2.25	0.70
1:B:1007:GLN:H	1:B:1007:GLN:CD	1.98	0.69
1:D:971:LYS:HG3	1:D:1081:HIS:CD2	2.27	0.69
1:B:812:MET:O	1:B:815:ARG:HG2	1.92	0.69
1:B:921:ILE:HG23	1:B:932:LEU:HD12	1.73	0.69
1:B:981:LEU:CD2	1:B:1004:LEU:HD12	2.20	0.69
1:B:798:GLU:CB	1:B:824:SER:HB3	2.12	0.69
1:B:807:ALA:CB	1:B:1008:GLN:OE1	2.41	0.69
1:B:1016:LYS:CG	1:B:1019:GLN:OE1	2.39	0.69
1:D:1023:LYS:O	1:D:1026:LYS:CG	2.40	0.69
1:A:1141:VAL:CG1	1:A:1157:THR:HG23	2.23	0.69
1:B:780:THR:N	1:B:1014:LYS:HZ1	1.89	0.69
1:B:1010:PHE:CE2	1:B:1024:LYS:HG3	2.26	0.69
1:A:552:GLU:N	1:A:552:GLU:OE2	2.26	0.68
1:B:1024:LYS:NZ	1:B:1130:SER:O	2.24	0.68
1:D:1007:GLN:H	1:D:1007:GLN:CD	2.01	0.68
1:D:812:MET:O	1:D:815:ARG:HG2	1.92	0.68
1:D:766:GLY:O	1:D:825:ARG:NH1	2.26	0.68
1:B:1082:ARG:NE	1:C:1072:GLN:HE22	1.91	0.68
1:D:971:LYS:HD2	1:D:1083:HIS:NE2	2.09	0.68
1:A:616:VAL:HG12	1:A:636:TYR:CZ	2.28	0.68
1:A:1006:LEU:CD1	1:A:1067:ALA:HB2	2.18	0.68
1:B:551:LYS:HE2	1:B:555:GLU:OE1	1.95	0.67
1:C:975:PHE:CZ	1:C:1078:MET:HE1	2.26	0.67
1:D:655:ASP:OD2	1:D:686:THR:HG23	1.93	0.67
1:D:554:ARG:NH1	1:D:596:GLU:OE2	2.27	0.67
1:A:956:LEU:O	1:A:983:LYS:HE2	1.96	0.66
1:D:1006:LEU:HD22	1:D:1063:ALA:O	1.96	0.66
1:D:703:LEU:HD12	1:D:703:LEU:C	2.14	0.66
1:D:574:GLN:CD	1:D:697:VAL:HG13	2.21	0.66
1:B:971:LYS:HD3	1:B:1083:HIS:HE1	1.61	0.65
1:B:768:ILE:HB	1:B:827:VAL:HG23	1.79	0.65
1:A:1077:ILE:HG12	2:E:10:U:H1'	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1024:LYS:NZ	1:C:1130:SER:O	2.25	0.65
1:D:660:ARG:HD2	2:H:11:U:P	2.36	0.65
1:B:1073:GLN:NE2	2:F:10:U:OP2	2.29	0.65
1:A:783:GLU:HA	1:A:783:GLU:OE1	1.98	0.64
1:C:1008:GLN:O	1:C:1063:ALA:HB2	1.96	0.64
1:B:839:THR:HG22	1:B:882:ARG:HH12	1.62	0.64
1:D:974:ASP:OD2	1:D:1081:HIS:CE1	2.51	0.64
1:D:659:GLN:O	1:D:662:ILE:HG22	1.97	0.64
1:A:659:GLN:O	1:A:662:ILE:HG22	1.98	0.64
1:A:1090:ARG:HH21	1:A:1090:ARG:HB2	1.63	0.64
1:B:655:ASP:OD2	1:B:686:THR:HG23	1.97	0.64
1:C:659:GLN:O	1:C:662:ILE:HG22	1.96	0.64
1:A:850:GLY:HA2	1:A:871:ILE:HD13	1.80	0.64
1:D:574:GLN:OE1	1:D:697:VAL:HG13	1.98	0.64
1:B:839:THR:HG22	1:B:882:ARG:NH1	2.13	0.63
1:B:1142:LEU:C	1:B:1142:LEU:HD23	2.23	0.63
1:C:975:PHE:CZ	1:C:1078:MET:HE2	2.21	0.63
1:C:612:ARG:HG3	2:G:8:U:OP1	1.98	0.63
1:D:1140:TRP:CE2	1:D:1162:PRO:HG3	2.32	0.63
1:B:767:ASP:OD2	1:B:826:LYS:HE2	1.98	0.63
1:D:631:GLY:O	1:D:645:SER:HB3	1.98	0.63
1:B:659:GLN:O	1:B:662:ILE:HG22	1.98	0.63
1:B:734:PHE:CE2	1:B:886:THR:C	2.77	0.63
1:D:554:ARG:HH21	1:D:554:ARG:CG	2.11	0.63
1:C:804:ILE:HG23	1:C:830:ALA:HB2	1.67	0.63
1:A:645:SER:H	1:A:648:THR:HG23	1.64	0.62
1:B:644:THR:HG21	1:B:651:LYS:HD2	1.81	0.62
1:D:638:ILE:HG22	1:D:654:THR:HG23	1.79	0.62
1:C:592:GLN:HE21	1:C:626:VAL:HG23	1.63	0.62
1:C:978:GLU:HG3	1:C:979:PRO:HD2	1.81	0.62
1:A:611:ARG:NH1	1:A:681:GLU:OE2	2.24	0.62
1:B:1082:ARG:CZ	1:C:1072:GLN:HE22	2.11	0.62
1:B:798:GLU:HB3	1:B:824:SER:CB	2.13	0.61
1:A:981:LEU:HD21	1:A:1004:LEU:CD1	2.31	0.61
1:D:696:THR:OG1	1:D:699:ARG:NH2	2.32	0.61
1:B:636:TYR:HA	1:B:652:TYR:O	2.01	0.61
1:C:559:VAL:HG23	1:C:729:ILE:HD13	1.83	0.61
1:B:600:THR:OG1	1:B:649:LYS:HB3	2.00	0.61
1:B:921:ILE:HG23	1:B:932:LEU:CD1	2.31	0.61
1:D:971:LYS:HG3	1:D:1081:HIS:HD2	1.66	0.61
1:B:1020:ALA:HA	1:B:1058:TYR:HD2	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:766:GLY:O	1:C:825:ARG:HD2	2.00	0.61
1:A:638:ILE:HG22	1:A:654:THR:HG23	1.82	0.60
1:C:636:TYR:HA	1:C:652:TYR:O	2.02	0.60
1:C:778:ILE:HG12	1:C:829:ILE:HG22	1.82	0.60
1:B:638:ILE:HG22	1:B:654:THR:HG23	1.82	0.60
1:B:1016:LYS:HG2	1:B:1019:GLN:CD	2.26	0.60
1:A:981:LEU:HD21	1:A:1004:LEU:HD13	1.84	0.60
1:D:702:ASP:OD1	1:D:702:ASP:N	2.34	0.60
1:A:636:TYR:HA	1:A:652:TYR:O	2.01	0.60
1:B:759:ILE:O	1:B:763:GLU:HB2	2.02	0.60
1:B:779:ASP:HB3	1:B:1014:LYS:HE3	1.84	0.59
1:B:916:ASN:OD1	1:B:917:LEU:N	2.35	0.59
1:A:780:THR:OG1	1:A:1014:LYS:HE2	2.02	0.59
1:B:933:LEU:HD12	1:B:933:LEU:N	2.16	0.59
1:B:839:THR:CG2	1:B:882:ARG:HH12	2.15	0.59
1:B:1007:GLN:OE1	1:B:1007:GLN:N	2.23	0.59
1:B:970:ARG:HD2	1:B:970:ARG:C	2.25	0.59
1:D:643:VAL:HG12	1:D:643:VAL:O	2.02	0.59
1:A:1147:VAL:HG22	1:A:1155:HIS:CD2	2.38	0.59
1:D:812:MET:HE3	1:D:815:ARG:NH1	2.17	0.59
1:D:1155:HIS:NE2	2:H:4:U:C5	2.69	0.59
1:A:1106:ASN:HB3	1:A:1143:TYR:CD1	2.38	0.59
1:C:863:MET:HG3	1:C:1146:LEU:HG	1.84	0.59
1:D:591:THR:HG22	1:D:650:ILE:HD13	1.85	0.59
1:D:1006:LEU:HD11	1:D:1067:ALA:HB2	1.85	0.58
1:B:796:VAL:HG22	1:B:797:PRO:HD2	1.84	0.58
1:C:638:ILE:HG22	1:C:654:THR:HG23	1.85	0.58
1:C:700:ARG:HG2	1:C:702:ASP:OD1	2.03	0.58
1:B:941:PRO:HG2	1:B:946:MET:SD	2.44	0.58
1:C:804:ILE:HG12	1:C:838:ILE:CD1	2.34	0.58
1:D:636:TYR:HA	1:D:652:TYR:O	2.03	0.58
1:B:832:ASN:ND2	2:F:5:U:O2'	2.37	0.58
1:A:1008:GLN:O	1:A:1008:GLN:HG3	2.03	0.58
1:C:581:GLU:HB2	1:C:731:GLY:CA	2.33	0.58
1:C:592:GLN:HE21	1:C:626:VAL:CG2	2.17	0.58
1:C:998:VAL:HG12	1:C:1071:ARG:HG3	1.86	0.58
1:D:917:LEU:O	1:D:921:ILE:HG12	2.04	0.58
1:D:1140:TRP:CD2	1:D:1162:PRO:HG3	2.40	0.57
1:A:1140:TRP:CE2	1:A:1162:PRO:HG3	2.39	0.57
1:A:1147:VAL:O	1:A:1147:VAL:HG23	2.04	0.57
1:C:1152:GLU:H	1:C:1152:GLU:CD	2.13	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:5:U:H5''	2:G:5:U:H6	1.67	0.57
1:A:643:VAL:O	1:A:643:VAL:HG12	2.03	0.57
1:A:1141:VAL:HG13	1:A:1157:THR:CG2	2.33	0.57
1:B:1140:TRP:CE2	1:B:1162:PRO:HG3	2.39	0.57
1:C:974:ASP:OD2	1:C:1081:HIS:CE1	2.58	0.57
1:A:1090:ARG:HB2	1:A:1090:ARG:NH2	2.18	0.57
1:C:700:ARG:HH11	1:C:703:LEU:HB2	1.69	0.57
1:B:804:ILE:HG22	1:B:830:ALA:HB2	1.81	0.56
1:C:577:ILE:CD1	1:C:724:CYS:SG	2.92	0.56
1:C:756:VAL:HG11	1:C:785:LEU:HD21	1.86	0.56
1:B:780:THR:N	1:B:1014:LYS:NZ	2.51	0.56
1:A:694:LYS:O	1:A:697:VAL:CG2	2.53	0.56
1:A:804:ILE:CG2	1:A:830:ALA:HB2	2.36	0.56
1:B:736:VAL:CA	1:B:889:GLY:O	2.44	0.56
1:B:971:LYS:HD2	1:B:986:ILE:HD13	1.87	0.56
1:C:763:GLU:HB3	1:C:764:PRO:HD2	1.88	0.56
1:C:1080:ARG:HG2	1:C:1080:ARG:NH1	2.09	0.56
1:A:667:ASP:OD1	1:A:699:ARG:HD2	2.06	0.56
1:C:1140:TRP:CE2	1:C:1162:PRO:HG3	2.41	0.56
1:D:574:GLN:CD	1:D:697:VAL:CG1	2.79	0.56
1:A:804:ILE:CG2	1:A:830:ALA:CB	2.84	0.56
1:C:570:VAL:HG22	1:C:706:ILE:HD11	1.88	0.56
1:D:550:ILE:HD11	1:D:625:GLU:HA	1.88	0.56
1:D:551:LYS:O	1:D:555:GLU:HG3	2.06	0.56
1:A:1106:ASN:HB3	1:A:1143:TYR:CE1	2.41	0.55
1:B:734:PHE:CD2	1:B:887:GLY:HA2	2.41	0.55
1:A:1147:VAL:HG22	1:A:1155:HIS:NE2	2.21	0.55
1:C:691:ALA:CB	1:C:935:PHE:CE2	2.89	0.55
1:D:563:ARG:O	1:D:567:ILE:HD12	2.07	0.55
1:D:642:ASP:OD1	1:D:644:THR:HB	2.07	0.55
1:B:863:MET:HG3	1:B:1146:LEU:HG	1.88	0.55
1:D:1007:GLN:OE1	1:D:1007:GLN:N	2.24	0.55
1:B:700:ARG:HG2	1:B:702:ASP:OD1	2.07	0.55
1:C:600:THR:OG1	1:C:649:LYS:HB3	2.07	0.55
1:A:639:ARG:HD2	2:E:9:U:H2'	1.88	0.55
1:A:981:LEU:CD2	1:A:1004:LEU:HD12	2.37	0.55
1:C:771:PHE:HD1	1:C:831:THR:O	1.90	0.55
1:D:554:ARG:HD2	1:D:625:GLU:OE1	2.05	0.55
1:C:695:LYS:O	1:C:698:LYS:HG2	2.07	0.55
1:D:577:ILE:CD1	1:D:724:CYS:SG	2.94	0.55
1:A:1141:VAL:CG1	1:A:1157:THR:CG2	2.85	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:734:PHE:HE2	1:B:886:THR:O	1.87	0.54
1:D:832:ASN:C	1:D:832:ASN:HD22	2.14	0.54
2:F:1:U:O2	2:F:1:U:H5"	2.07	0.54
1:B:1020:ALA:CA	1:B:1058:TYR:CD2	2.81	0.54
1:A:758:GLN:O	1:A:762:THR:HG23	2.08	0.54
1:B:800:ILE:HG13	1:B:826:LYS:HA	1.89	0.54
1:A:637:THR:OG1	1:A:651:LYS:HE3	2.06	0.54
1:B:734:PHE:CD2	1:B:887:GLY:CA	2.91	0.54
1:B:759:ILE:HG23	1:B:763:GLU:HG3	1.90	0.54
1:D:756:VAL:HG11	1:D:785:LEU:HD21	1.89	0.54
1:C:971:LYS:HE3	1:C:986:ILE:HD13	1.90	0.53
1:A:1147:VAL:O	1:A:1147:VAL:CG2	2.55	0.53
1:B:998:VAL:HG12	1:B:1071:ARG:HG3	1.89	0.53
1:D:759:ILE:HD11	1:D:892:PHE:CD1	2.43	0.53
1:D:900:TYR:CD1	1:D:900:TYR:C	2.86	0.53
1:B:577:ILE:CD1	1:B:724:CYS:SG	2.96	0.53
1:D:600:THR:OG1	1:D:649:LYS:HB3	2.08	0.53
1:D:703:LEU:CD1	1:D:704:LYS:O	2.57	0.53
1:C:553:GLN:O	1:C:557:LEU:HD13	2.09	0.53
1:C:632:GLN:O	1:C:648:THR:HA	2.08	0.53
1:C:1140:TRP:CD2	1:C:1162:PRO:HG3	2.44	0.53
1:D:659:GLN:HG2	1:D:923:LEU:HD21	1.90	0.53
1:D:755:THR:O	1:D:759:ILE:HG13	2.07	0.53
1:A:660:ARG:HD2	2:E:11:U:P	2.49	0.53
1:A:1140:TRP:CD2	1:A:1162:PRO:HG3	2.44	0.53
1:B:933:LEU:H	1:B:933:LEU:CD1	2.22	0.53
1:B:1140:TRP:CD2	1:B:1162:PRO:HG3	2.43	0.53
1:C:653:MET:HE1	1:C:671:TYR:CE2	2.44	0.53
1:A:736:VAL:HA	1:A:889:GLY:O	2.09	0.53
1:A:1009:ILE:HD12	1:A:1009:ILE:N	2.19	0.53
1:C:736:VAL:HA	1:C:889:GLY:O	2.09	0.53
1:A:610:PRO:HD3	1:A:678:GLU:HB2	1.91	0.52
1:B:796:VAL:CG2	1:B:797:PRO:HD2	2.39	0.52
1:A:659:GLN:HG2	1:A:923:LEU:HD21	1.91	0.52
1:C:1080:ARG:HH11	1:C:1080:ARG:CG	2.13	0.52
1:A:558:PRO:HG2	1:A:584:SER:HA	1.92	0.52
1:A:853:LYS:NZ	2:E:5:U:OP1	2.43	0.52
1:B:984:VAL:CG1	1:B:997:MET:SD	2.97	0.52
1:B:1016:LYS:HG2	1:B:1019:GLN:HB2	1.92	0.52
1:C:637:THR:OG1	1:C:651:LYS:HE3	2.09	0.52
1:D:1010:PHE:CE1	1:D:1024:LYS:HG3	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:832:ASN:ND2	2:G:5:U:O2'	2.41	0.52
1:C:1137:GLN:HA	1:C:1137:GLN:NE2	2.24	0.52
1:A:605:ILE:HD12	1:A:673:VAL:HB	1.88	0.52
1:D:554:ARG:CG	1:D:554:ARG:NH2	2.73	0.52
1:D:632:GLN:O	1:D:648:THR:HA	2.09	0.52
1:A:1073:GLN:NE2	2:E:10:U:OP2	2.42	0.51
1:D:590:VAL:HG11	1:D:675:MET:SD	2.50	0.51
1:B:760:HIS:ND1	1:B:825:ARG:NE	2.52	0.51
1:D:974:ASP:OD2	1:D:1081:HIS:NE2	2.43	0.51
1:C:971:LYS:HG3	1:C:1081:HIS:CD2	2.46	0.51
1:D:798:GLU:OE2	1:D:824:SER:CB	2.58	0.51
2:F:7:U:H2'	2:F:8:U:C6	2.45	0.51
1:C:592:GLN:NE2	1:C:626:VAL:HG23	2.25	0.51
1:D:832:ASN:HD22	1:D:833:ILE:N	2.09	0.51
1:A:632:GLN:O	1:A:648:THR:CG2	2.51	0.51
1:A:850:GLY:C	1:A:871:ILE:HG23	2.36	0.51
1:C:978:GLU:CG	1:C:979:PRO:HD2	2.41	0.51
1:A:750:GLU:OE1	1:A:788:ARG:NH2	2.32	0.50
1:A:861:LEU:HD22	1:A:1123:THR:HG21	1.93	0.50
2:E:1:U:H5''	2:E:1:U:O2	2.11	0.50
1:D:763:GLU:OE1	1:D:764:PRO:HD2	2.10	0.50
1:B:1012:ARG:HG2	1:B:1017:GLN:HG2	1.92	0.50
1:B:1077:ILE:HG12	2:F:10:U:H1'	1.93	0.50
1:A:790:LYS:O	1:A:792:LEU:O	2.29	0.50
1:D:590:VAL:CG1	1:D:675:MET:SD	3.00	0.50
1:D:998:VAL:O	1:D:1002:SER:OG	2.28	0.50
1:A:612:ARG:HG3	2:E:8:U:OP1	2.12	0.50
1:D:586:LYS:O	1:D:590:VAL:HG12	2.12	0.49
1:B:764:PRO:HG2	1:B:845:TYR:OH	2.13	0.49
1:C:804:ILE:CG2	1:C:830:ALA:CB	2.71	0.49
1:B:1011:TYR:CD1	1:B:1011:TYR:C	2.90	0.49
1:D:1104:PHE:HB2	1:D:1173:PHE:HE2	1.77	0.49
1:A:679:ALA:HB3	1:A:708:THR:O	2.13	0.49
1:A:759:ILE:HA	1:A:763:GLU:HG3	1.94	0.49
1:C:1040:VAL:HG13	1:C:1059:ILE:CD1	2.42	0.49
1:B:632:GLN:O	1:B:648:THR:CG2	2.53	0.49
1:A:547:ASN:N	1:A:547:ASN:ND2	2.61	0.49
1:A:694:LYS:O	1:A:697:VAL:HG22	2.12	0.49
1:B:804:ILE:HG23	1:B:830:ALA:HB1	1.94	0.49
1:B:1155:HIS:NE2	2:F:4:U:H5	2.11	0.49
1:B:1082:ARG:NE	1:C:1072:GLN:NE2	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:852:VAL:CG1	1:C:871:ILE:HG22	2.43	0.49
1:D:757:MET:HE1	1:D:789:MET:SD	2.53	0.49
1:A:547:ASN:N	1:A:547:ASN:HD22	2.09	0.49
1:A:1018:GLN:O	1:A:1022:GLN:CG	2.60	0.49
1:C:691:ALA:HB1	1:C:935:PHE:CE2	2.48	0.49
1:D:632:GLN:O	1:D:648:THR:OG1	2.24	0.49
1:A:700:ARG:HH11	1:A:703:LEU:HB2	1.77	0.48
1:C:866:LEU:HD11	2:G:5:U:C5	2.48	0.48
1:D:998:VAL:HG12	1:D:1071:ARG:HG3	1.94	0.48
1:D:799:LEU:HD12	1:D:825:ARG:O	2.14	0.48
1:D:1085:ILE:O	1:D:1086:ILE:HD13	2.14	0.48
1:B:974:ASP:OD2	1:B:1081:HIS:CE1	2.66	0.48
1:A:1005:ASN:HD22	1:A:1070:VAL:HG11	1.79	0.48
1:A:1006:LEU:HD11	1:A:1067:ALA:CB	2.24	0.48
1:B:917:LEU:O	1:B:921:ILE:HG12	2.14	0.48
1:A:610:PRO:CD	1:A:678:GLU:HB2	2.44	0.48
1:B:591:THR:CG2	1:B:652:TYR:OH	2.60	0.48
1:D:575:ILE:HD11	1:D:697:VAL:HG21	1.96	0.48
2:G:1:U:H3'	2:G:1:U:O2	2.13	0.48
1:D:1128:HIS:NE2	2:H:4:U:OP1	2.41	0.48
1:A:758:GLN:O	1:A:762:THR:N	2.40	0.48
1:A:863:MET:HG3	1:A:1146:LEU:HG	1.96	0.48
1:D:637:THR:OG1	1:D:651:LYS:HE3	2.14	0.48
1:A:700:ARG:HH11	1:A:703:LEU:HD12	1.79	0.47
1:C:1073:GLN:NE2	2:G:10:U:OP2	2.38	0.47
1:D:1155:HIS:CE1	2:H:4:U:H5	2.32	0.47
1:B:779:ASP:C	1:B:1014:LYS:HZ1	2.10	0.47
1:D:574:GLN:NE2	1:D:697:VAL:CG1	2.78	0.47
1:D:984:VAL:HG11	1:D:1099:LEU:HD23	1.96	0.47
1:B:984:VAL:HG13	1:B:997:MET:HE3	1.96	0.47
1:C:917:LEU:O	1:C:921:ILE:HG12	2.15	0.47
1:A:921:ILE:HD11	1:A:946:MET:HE3	1.97	0.47
1:A:1079:GLU:CG	1:D:1082:ARG:HH22	2.17	0.47
1:B:735:PRO:HG2	1:B:888:PRO:HB3	1.97	0.47
1:C:929:ILE:HD13	1:C:935:PHE:HE1	1.80	0.47
1:D:549:SER:O	1:D:553:GLN:HG3	2.15	0.47
1:D:591:THR:HG21	1:D:650:ILE:HG23	1.96	0.47
1:A:804:ILE:HG23	1:A:804:ILE:O	2.15	0.47
1:B:1012:ARG:NH2	2:F:3:U:O4	2.48	0.47
1:C:657:MET:CE	2:G:10:U:OP1	2.58	0.47
1:A:551:LYS:HE3	1:A:555:GLU:OE2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:631:GLY:O	1:A:645:SER:OG	2.30	0.47
1:D:994:SER:O	1:D:998:VAL:HG23	2.15	0.47
1:D:734:PHE:CB	1:D:888:PRO:N	2.76	0.46
1:D:1117:LYS:HB3	1:D:1117:LYS:HE2	1.68	0.46
2:H:1:U:O2	2:H:1:U:H3'	2.15	0.46
1:A:616:VAL:CG1	1:A:636:TYR:CE1	2.97	0.46
1:A:770:VAL:HG12	1:A:847:VAL:HB	1.98	0.46
1:A:577:ILE:HD12	1:A:724:CYS:SG	2.55	0.46
1:B:813:GLN:O	1:B:816:ILE:HG12	2.16	0.46
1:C:682:ARG:NH1	1:C:938:MET:HE2	2.30	0.46
1:A:1009:ILE:H	1:A:1009:ILE:CD1	2.21	0.46
1:D:804:ILE:CG2	1:D:830:ALA:HB2	2.45	0.46
1:A:926:ALA:HA	1:A:965:LEU:HD22	1.96	0.46
1:A:1018:GLN:O	1:A:1022:GLN:HG3	2.16	0.46
1:B:913:GLN:NE2	1:B:939:ASP:HB3	2.30	0.46
1:C:913:GLN:NE2	1:C:939:ASP:HB3	2.31	0.46
1:C:1077:ILE:HG12	2:G:10:U:H1'	1.98	0.46
1:A:981:LEU:HD22	1:A:1004:LEU:HD12	1.98	0.46
1:B:981:LEU:HD21	1:B:1004:LEU:HD13	1.92	0.46
1:C:994:SER:O	1:C:998:VAL:HG23	2.15	0.46
1:D:703:LEU:HD12	1:D:704:LYS:O	2.16	0.46
1:A:558:PRO:HG3	1:A:732:ARG:HG2	1.97	0.46
1:B:612:ARG:HG3	2:F:8:U:OP1	2.16	0.46
1:B:734:PHE:CE2	1:B:887:GLY:HA2	2.51	0.46
1:D:804:ILE:CG2	1:D:830:ALA:CB	2.94	0.46
1:A:576:LEU:HD23	1:A:577:ILE:C	2.41	0.46
1:A:1147:VAL:CG2	1:A:1155:HIS:NE2	2.79	0.46
1:B:756:VAL:HG21	1:B:770:VAL:HG21	1.97	0.46
1:C:1077:ILE:HD11	2:G:10:U:H1'	1.98	0.46
1:D:804:ILE:HG23	1:D:804:ILE:O	2.16	0.46
1:D:806:SER:OG	2:H:7:U:OP2	2.28	0.46
1:D:855:ASN:HB3	1:D:915:GLN:HE22	1.81	0.46
1:D:1006:LEU:HD21	1:D:1067:ALA:N	2.30	0.46
1:B:1011:TYR:CB	1:B:1060:GLN:NE2	2.60	0.46
1:D:813:GLN:O	1:D:816:ILE:HG12	2.15	0.46
1:A:591:THR:CG2	1:A:652:TYR:OH	2.59	0.45
1:A:981:LEU:CD2	1:A:1004:LEU:CD1	2.94	0.45
1:B:760:HIS:HA	1:B:825:ARG:NH1	2.31	0.45
1:D:1103:PHE:HB2	1:D:1142:LEU:HD11	1.97	0.45
1:C:977:MET:HE1	1:C:1074:ILE:CD1	2.46	0.45
1:A:1176:ALA:O	1:B:1174:LYS:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:581:GLU:OE1	1:D:731:GLY:CA	2.46	0.45
1:D:812:MET:HE3	1:D:815:ARG:HH12	1.82	0.45
1:A:991:LYS:O	1:A:1094:LYS:HD3	2.16	0.45
1:A:1077:ILE:CG1	2:E:10:U:H1'	2.44	0.45
1:B:778:ILE:HG12	1:B:829:ILE:HG22	1.98	0.45
1:D:863:MET:HG3	1:D:1146:LEU:HG	1.98	0.45
1:A:855:ASN:HB3	1:A:915:GLN:HE22	1.82	0.45
1:B:798:GLU:O	1:B:824:SER:CA	2.64	0.45
1:C:1111:ASP:N	1:C:1115:GLY:O	2.46	0.45
1:D:699:ARG:HE	1:D:699:ARG:HB2	1.59	0.45
1:B:620:LYS:HG2	1:B:630:LEU:HD22	1.98	0.45
1:B:1008:GLN:O	1:B:1063:ALA:HB2	2.16	0.45
1:D:696:THR:HA	1:D:699:ARG:HH21	1.81	0.45
1:B:1007:GLN:CD	1:B:1007:GLN:N	2.72	0.45
1:D:667:ASP:O	1:D:700:ARG:HD3	2.18	0.44
1:D:900:TYR:HA	1:D:904:MET:HG3	1.99	0.44
1:A:804:ILE:CG2	1:A:830:ALA:HA	2.47	0.44
1:A:808:LEU:HD11	1:A:812:MET:HE2	1.99	0.44
1:C:984:VAL:HG11	1:C:1099:LEU:HD23	1.99	0.44
1:D:1007:GLN:CD	1:D:1007:GLN:N	2.73	0.44
1:A:841:ASP:OD1	1:A:885:ARG:NH2	2.46	0.44
1:A:1054:CYS:HB3	1:A:1059:ILE:O	2.18	0.44
1:B:760:HIS:O	1:B:825:ARG:NH1	2.49	0.44
1:B:766:GLY:O	1:B:825:ARG:CD	2.62	0.44
1:B:1104:PHE:HB2	1:B:1173:PHE:HE2	1.82	0.44
1:B:827:VAL:HG13	1:B:827:VAL:O	2.16	0.44
1:B:1155:HIS:CE1	2:F:4:U:C5	2.97	0.44
1:B:1010:PHE:CD2	1:B:1024:LYS:HG3	2.52	0.44
1:C:1080:ARG:NH1	1:C:1080:ARG:CG	2.73	0.44
1:D:660:ARG:HD2	2:H:11:U:OP1	2.17	0.44
1:B:771:PHE:HD1	1:B:831:THR:O	2.01	0.44
1:C:579:VAL:HG13	1:C:728:THR:CG2	2.38	0.44
1:D:703:LEU:HD11	1:D:704:LYS:O	2.18	0.44
1:D:811:GLU:H	1:D:811:GLU:HG2	1.60	0.44
1:B:1006:LEU:N	1:B:1006:LEU:HD23	2.33	0.44
1:B:769:LEU:O	1:B:846:VAL:HA	2.18	0.43
1:B:1006:LEU:HD13	1:B:1063:ALA:O	2.18	0.43
1:B:678:GLU:OE1	1:B:680:HIS:HE1	2.01	0.43
1:B:737:GLU:N	1:B:889:GLY:O	2.51	0.43
1:A:656:GLY:HA3	2:E:8:U:O2'	2.19	0.43
1:B:757:MET:HE2	1:B:757:MET:HB3	1.95	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:763:GLU:HB3	1:D:764:PRO:HD2	1.99	0.43
1:B:644:THR:HA	1:B:648:THR:HG21	2.00	0.43
1:B:1105:ARG:HD3	1:B:1105:ARG:HA	1.94	0.43
1:C:804:ILE:HG12	1:C:838:ILE:HD13	1.98	0.43
1:D:1129:PRO:HG2	2:H:3:U:C6	2.54	0.43
1:A:1155:HIS:NE2	2:E:4:U:H5	2.16	0.43
1:B:668:LEU:O	1:B:700:ARG:NH1	2.52	0.43
1:D:913:GLN:NE2	1:D:939:ASP:HB3	2.34	0.43
1:A:1119:LEU:HD21	1:A:1140:TRP:HZ3	1.84	0.43
1:D:736:VAL:HA	1:D:889:GLY:O	2.19	0.43
1:D:1142:LEU:C	1:D:1142:LEU:HD23	2.43	0.43
1:B:659:GLN:HG2	1:B:923:LEU:HD21	2.01	0.43
1:B:767:ASP:HB2	1:B:842:TYR:O	2.19	0.43
1:A:678:GLU:HB3	1:A:681:GLU:HG3	2.01	0.43
1:A:984:VAL:HG11	1:A:1099:LEU:HD23	2.00	0.43
1:A:1018:GLN:O	1:A:1022:GLN:HG2	2.19	0.43
1:B:611:ARG:HH12	1:B:681:GLU:CD	2.27	0.43
1:B:985:LEU:O	1:B:988:SER:OG	2.29	0.43
1:C:978:GLU:HG3	1:C:979:PRO:CD	2.47	0.43
1:B:994:SER:O	1:B:998:VAL:HG23	2.20	0.42
1:B:1071:ARG:O	1:B:1075:VAL:HG23	2.19	0.42
1:C:717:PHE:HE1	1:C:938:MET:HE1	1.81	0.42
1:C:771:PHE:CE1	1:C:834:ALA:HB3	2.54	0.42
1:D:554:ARG:HH21	1:D:554:ARG:HG3	1.81	0.42
1:A:605:ILE:CD1	1:A:673:VAL:CB	2.84	0.42
1:B:833:ILE:HB	2:F:6:U:H5'	2.01	0.42
1:B:1053:TRP:CH2	1:B:1059:ILE:HD12	2.54	0.42
1:B:1103:PHE:HB2	1:B:1142:LEU:HD11	2.01	0.42
1:A:913:GLN:NE2	1:A:939:ASP:HB3	2.34	0.42
1:C:804:ILE:HG12	1:C:838:ILE:HD11	2.01	0.42
1:C:1104:PHE:HB2	1:C:1173:PHE:HE2	1.85	0.42
1:D:620:LYS:HG2	1:D:630:LEU:HD22	2.01	0.42
1:D:678:GLU:OE1	1:D:680:HIS:HE1	2.02	0.42
1:B:656:GLY:HA3	2:F:8:U:O2'	2.19	0.42
1:B:760:HIS:HA	1:B:825:ARG:CZ	2.49	0.42
1:C:679:ALA:HA	1:C:686:THR:HG21	2.01	0.42
1:D:557:LEU:HD13	1:D:585:GLY:HA3	2.01	0.42
1:A:645:SER:H	1:A:648:THR:CG2	2.31	0.42
1:A:759:ILE:O	1:A:763:GLU:HB2	2.19	0.42
1:B:694:LYS:O	1:B:697:VAL:HG22	2.19	0.42
1:C:1152:GLU:OE1	1:C:1152:GLU:N	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1119:LEU:HD21	1:A:1140:TRP:CZ3	2.55	0.42
1:A:1142:LEU:HD23	1:A:1142:LEU:C	2.45	0.42
1:B:547:ASN:N	1:B:547:ASN:HD22	2.16	0.42
1:C:659:GLN:HE21	1:C:688:VAL:HG11	1.85	0.42
1:C:866:LEU:HD23	1:C:866:LEU:HA	1.87	0.42
1:D:639:ARG:HG3	2:H:10:U:OP2	2.19	0.42
1:A:575:ILE:HD11	1:A:697:VAL:HG21	2.01	0.42
1:A:1155:HIS:CE1	2:E:4:U:H5	2.37	0.42
1:B:808:LEU:HA	1:B:809:PRO:HD3	1.94	0.42
1:D:971:LYS:HE3	1:D:986:ILE:HD13	2.01	0.42
1:A:612:ARG:NH2	1:A:639:ARG:O	2.52	0.42
1:A:802:LEU:HD22	1:A:815:ARG:HG3	2.02	0.42
1:C:1016:LYS:HD3	1:C:1019:GLN:NE2	2.35	0.42
1:D:549:SER:HB3	1:D:552:GLU:CD	2.45	0.42
1:A:769:LEU:O	1:A:846:VAL:HA	2.20	0.41
1:A:1005:ASN:ND2	1:A:1070:VAL:HG11	2.35	0.41
1:B:926:ALA:HA	1:B:965:LEU:HD22	2.02	0.41
1:B:660:ARG:HG2	2:F:10:U:H5"	2.02	0.41
1:C:927:MET:HE2	1:C:927:MET:HB3	1.91	0.41
1:A:940:PRO:HA	1:A:941:PRO:HD3	1.85	0.41
1:C:558:PRO:HG3	1:C:732:ARG:CD	2.50	0.41
1:C:769:LEU:O	1:C:846:VAL:HA	2.20	0.41
1:D:769:LEU:O	1:D:846:VAL:HA	2.20	0.41
1:A:813:GLN:O	1:A:816:ILE:HG12	2.20	0.41
1:C:839:THR:HG22	1:C:882:ARG:HH12	1.84	0.41
1:D:594:LEU:HD11	1:D:706:ILE:HD11	2.02	0.41
1:D:804:ILE:HG23	1:D:830:ALA:HA	2.01	0.41
1:B:659:GLN:HE21	1:B:688:VAL:HG11	1.85	0.41
2:E:11:U:H6	2:E:11:U:O5'	2.03	0.41
1:A:885:ARG:CZ	1:A:885:ARG:HB3	2.50	0.41
1:C:740:TYR:CE2	1:C:897:GLU:HA	2.56	0.41
1:C:1111:ASP:CB	1:C:1115:GLY:O	2.57	0.41
1:C:1137:GLN:HA	1:C:1137:GLN:HE21	1.85	0.41
1:B:863:MET:HE2	1:B:1146:LEU:HD21	2.01	0.41
1:C:804:ILE:HG22	1:C:830:ALA:HB2	1.93	0.41
1:C:812:MET:HE2	1:C:815:ARG:NH2	2.36	0.41
1:C:1008:GLN:O	1:C:1063:ALA:CB	2.64	0.41
1:C:1142:LEU:HD23	1:C:1142:LEU:C	2.46	0.41
1:D:586:LYS:HG3	1:D:587:THR:N	2.35	0.41
1:D:804:ILE:CG2	1:D:830:ALA:HA	2.50	0.41
1:D:882:ARG:HD3	1:D:882:ARG:HA	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:861:LEU:CD2	1:A:1123:THR:HG21	2.51	0.41
1:B:712:LEU:C	1:B:712:LEU:HD23	2.46	0.41
1:B:734:PHE:HE2	1:B:886:THR:C	2.23	0.41
1:B:779:ASP:HB3	1:B:1014:LYS:CE	2.48	0.41
1:B:835:GLU:HA	1:B:882:ARG:HD2	2.03	0.41
1:B:1020:ALA:HA	1:B:1058:TYR:CE2	2.51	0.41
1:B:607:CYS:HA	1:B:675:MET:O	2.21	0.41
1:C:1054:CYS:HB3	1:C:1059:ILE:O	2.20	0.41
1:A:569:ALA:HB1	1:A:576:LEU:HD11	2.02	0.40
1:B:782:CYS:SG	1:B:829:ILE:HD12	2.61	0.40
1:C:757:MET:HE2	1:C:757:MET:HB2	1.93	0.40
1:C:832:ASN:ND2	1:C:832:ASN:H	2.18	0.40
1:D:866:LEU:HD23	1:D:866:LEU:HA	1.91	0.40
1:B:816:ILE:HD12	1:B:840:ILE:HD11	2.03	0.40
1:C:694:LYS:O	1:C:697:VAL:HG22	2.20	0.40
1:C:929:ILE:HD13	1:C:935:PHE:CE1	2.55	0.40
1:C:1006:LEU:HD23	1:C:1006:LEU:HA	1.96	0.40
1:D:657:MET:SD	2:H:10:U:OP1	2.79	0.40
1:D:927:MET:HE2	1:D:927:MET:HB3	1.89	0.40
2:G:1:U:O2	2:G:1:U:C3'	2.70	0.40
1:D:953:LEU:HD21	1:D:979:PRO:HG2	2.04	0.40
1:D:1003:MET:HE3	1:D:1009:ILE:HD13	2.02	0.40
1:C:771:PHE:CD1	1:C:831:THR:O	2.73	0.40
1:D:976:PRO:CB	2:H:9:U:O2'	2.70	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 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	624/677 (92%)	607 (97%)	17 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	620/677 (92%)	596 (96%)	24 (4%)	0	100	100
1	C	623/677 (92%)	598 (96%)	25 (4%)	0	100	100
1	D	626/677 (92%)	606 (97%)	20 (3%)	0	100	100
All	All	2493/2708 (92%)	2407 (97%)	86 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	535/580 (92%)	514 (96%)	21 (4%)	28	57
1	B	538/580 (93%)	519 (96%)	19 (4%)	32	61
1	C	539/580 (93%)	526 (98%)	13 (2%)	43	72
1	D	540/580 (93%)	526 (97%)	14 (3%)	40	70
All	All	2152/2320 (93%)	2085 (97%)	67 (3%)	35	65

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	546	THR
1	A	547	ASN
1	A	577	ILE
1	A	611	ARG
1	A	711	THR
1	A	770	VAL
1	A	804	ILE
1	A	810	SER
1	A	867	VAL
1	A	869	THR
1	A	885	ARG
1	A	886	THR
1	A	894	LEU

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Mol	Chain	Res	Type
1	A	896	THR
1	A	901	GLN
1	A	933	LEU
1	A	1026	LYS
1	A	1056	GLU
1	A	1085	ILE
1	A	1141	VAL
1	A	1142	LEU
1	B	550	ILE
1	B	577	ILE
1	B	604	MET
1	B	616	VAL
1	B	617	SER
1	B	828	VAL
1	B	853	LYS
1	B	854	GLN
1	B	869	THR
1	B	886	THR
1	B	901	GLN
1	B	908	THR
1	B	914	ARG
1	B	980	SER
1	B	1015	ASP
1	B	1085	ILE
1	B	1105	ARG
1	B	1142	LEU
1	B	1161	GLU
1	C	577	ILE
1	C	579	VAL
1	C	713	ASP
1	C	832	ASN
1	C	836	THR
1	C	854	GLN
1	C	869	THR
1	C	1060	GLN
1	C	1080	ARG
1	C	1085	ILE
1	C	1111	ASP
1	C	1142	LEU
1	C	1175	LEU
1	D	554	ARG
1	D	577	ILE

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Mol	Chain	Res	Type
1	D	581	GLU
1	D	616	VAL
1	D	644	THR
1	D	645	SER
1	D	703	LEU
1	D	770	VAL
1	D	804	ILE
1	D	811	GLU
1	D	832	ASN
1	D	900	TYR
1	D	1060	GLN
1	D	1105	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	547	ASN
1	A	568	GLN
1	A	609	GLN
1	A	813	GLN
1	A	901	GLN
1	A	915	GLN
1	A	1005	ASN
1	A	1144	HIS
1	B	573	ASN
1	B	592	GLN
1	B	609	GLN
1	B	632	GLN
1	B	680	HIS
1	B	775	GLN
1	B	832	ASN
1	B	855	ASN
1	B	878	GLN
1	B	913	GLN
1	B	915	GLN
1	B	1046	ASN
1	B	1060	GLN
1	B	1081	HIS
1	B	1083	HIS
1	B	1097	GLN
1	B	1137	GLN
1	B	1155	HIS

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Mol	Chain	Res	Type
1	C	553	GLN
1	C	573	ASN
1	C	592	GLN
1	C	609	GLN
1	C	680	HIS
1	C	832	ASN
1	C	915	GLN
1	C	1051	ASN
1	C	1060	GLN
1	C	1072	GLN
1	C	1081	HIS
1	C	1097	GLN
1	C	1137	GLN
1	D	592	GLN
1	D	680	HIS
1	D	878	GLN
1	D	915	GLN
1	D	944	ASN
1	D	1022	GLN
1	D	1060	GLN
1	D	1081	HIS
1	D	1097	GLN
1	D	1106	ASN
1	D	1137	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	E	10/11 (90%)	5 (50%)	1 (10%)
2	F	9/11 (81%)	3 (33%)	1 (11%)
2	G	9/11 (81%)	2 (22%)	0
2	H	9/11 (81%)	5 (55%)	0
All	All	37/44 (84%)	15 (40%)	2 (5%)

All (15) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	E	3	U
2	E	4	U
2	E	5	U
2	E	7	U

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Mol	Chain	Res	Type
2	E	10	U
2	F	4	U
2	F	5	U
2	F	10	U
2	G	4	U
2	G	5	U
2	H	2	U
2	H	4	U
2	H	5	U
2	H	6	U
2	H	10	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	E	3	U
2	F	3	U

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 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	628/677 (92%)	-0.07	6 (0%) 79 78	58, 86, 128, 159	0
1	B	626/677 (92%)	-0.02	13 (2%) 63 61	60, 91, 131, 159	0
1	C	627/677 (92%)	-0.05	6 (0%) 79 78	63, 94, 139, 169	0
1	D	629/677 (92%)	-0.06	2 (0%) 90 89	52, 85, 121, 165	1 (0%)
2	E	11/11 (100%)	-0.11	0 100 100	77, 87, 129, 178	0
2	F	11/11 (100%)	-0.33	0 100 100	78, 92, 142, 154	0
2	G	11/11 (100%)	-0.41	0 100 100	81, 94, 156, 166	0
2	H	11/11 (100%)	-0.28	0 100 100	70, 85, 134, 145	0
All	All	2554/2752 (92%)	-0.06	27 (1%) 78 76	52, 89, 131, 178	1 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1177	PRO	4.2
1	C	1177	PRO	3.9
1	B	820	ALA	3.6
1	A	546	THR	3.5
1	B	837	ALA	3.5
1	A	1006	LEU	3.1
1	B	1009	ILE	2.8
1	C	652	TYR	2.7
1	D	1005	ASN	2.7
1	B	1005	ASN	2.7
1	C	549	SER	2.6
1	B	1177	PRO	2.5
1	A	1005	ASN	2.5
1	B	1112	PRO	2.5
1	B	819	PRO	2.4
1	D	1112	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	736	VAL	2.3
1	B	839	THR	2.3
1	B	611	ARG	2.3
1	A	1111	ASP	2.2
1	C	748	TYR	2.2
1	C	613	VAL	2.2
1	A	757	MET	2.2
1	B	766	GLY	2.2
1	B	978	GLU	2.2
1	C	550	ILE	2.2
1	B	834	ALA	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.