



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 08:55 AM EDT

PDB ID : 3SNP / pdb_00003snp
Title : Crystal structure analysis of iron regulatory protein 1 in complex with ferritin
H IRE RNA
Authors : Volz, K.; Selezneva, A.I.; Walden, W.E.
Deposited on : 2011-06-29
Resolution : 2.80 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

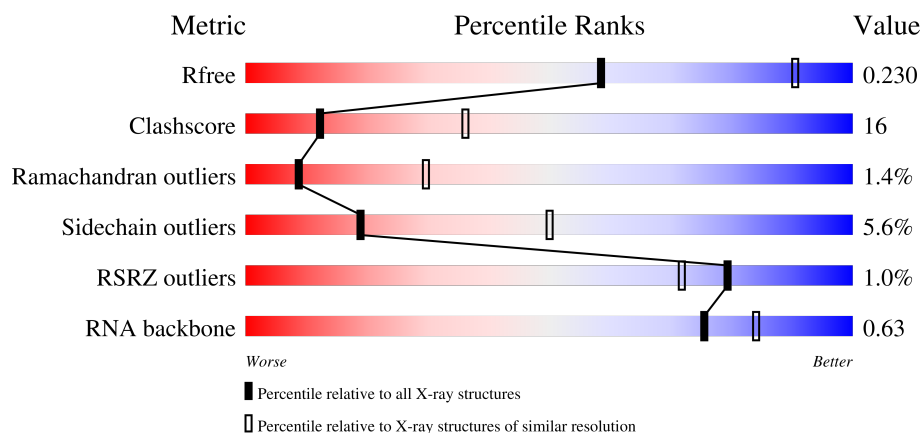
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)
RNA backbone	3983	1114 (3.00-2.60)

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	908	61% 29% 6%
1	B	908	59% 30% 6%
2	C	30	63% 27% 10%
2	D	30	57% 30% 10%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 15113 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 Cytoplasmic aconitate hydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	850	Total	C	N	O	S	0	0	0
			6657	4259	1143	1231	24			
1	B	850	Total	C	N	O	S	0	0	0
			6657	4259	1143	1231	24			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	expression tag	UNP Q01059
A	-17	GLY	-	expression tag	UNP Q01059
A	-16	HIS	-	expression tag	UNP Q01059
A	-15	HIS	-	expression tag	UNP Q01059
A	-14	HIS	-	expression tag	UNP Q01059
A	-13	HIS	-	expression tag	UNP Q01059
A	-12	HIS	-	expression tag	UNP Q01059
A	-11	HIS	-	expression tag	UNP Q01059
A	-10	ALA	-	expression tag	UNP Q01059
A	-9	ASP	-	expression tag	UNP Q01059
A	-8	ASP	-	expression tag	UNP Q01059
A	-7	ASP	-	expression tag	UNP Q01059
A	-6	ASP	-	expression tag	UNP Q01059
A	-5	LYS	-	expression tag	UNP Q01059
A	-4	ASP	-	expression tag	UNP Q01059
A	-3	GLY	-	expression tag	UNP Q01059
A	-2	VAL	-	expression tag	UNP Q01059
A	-1	ASP	-	expression tag	UNP Q01059
A	0	LYS	-	expression tag	UNP Q01059
A	1	LEU	-	expression tag	UNP Q01059
A	283	PRO	LEU	SEE REMARK 999	UNP Q01059
A	437	SER	CYS	engineered mutation	UNP Q01059
A	503	SER	CYS	engineered mutation	UNP Q01059
A	874	PHE	LEU	SEE REMARK 999	UNP Q01059
B	-18	MET	-	expression tag	UNP Q01059

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	GLY	-	expression tag	UNP Q01059
B	-16	HIS	-	expression tag	UNP Q01059
B	-15	HIS	-	expression tag	UNP Q01059
B	-14	HIS	-	expression tag	UNP Q01059
B	-13	HIS	-	expression tag	UNP Q01059
B	-12	HIS	-	expression tag	UNP Q01059
B	-11	HIS	-	expression tag	UNP Q01059
B	-10	ALA	-	expression tag	UNP Q01059
B	-9	ASP	-	expression tag	UNP Q01059
B	-8	ASP	-	expression tag	UNP Q01059
B	-7	ASP	-	expression tag	UNP Q01059
B	-6	ASP	-	expression tag	UNP Q01059
B	-5	LYS	-	expression tag	UNP Q01059
B	-4	ASP	-	expression tag	UNP Q01059
B	-3	GLY	-	expression tag	UNP Q01059
B	-2	VAL	-	expression tag	UNP Q01059
B	-1	ASP	-	expression tag	UNP Q01059
B	0	LYS	-	expression tag	UNP Q01059
B	1	LEU	-	expression tag	UNP Q01059
B	283	PRO	LEU	SEE REMARK 999	UNP Q01059
B	437	SER	CYS	engineered mutation	UNP Q01059
B	503	SER	CYS	engineered mutation	UNP Q01059
B	874	PHE	LEU	SEE REMARK 999	UNP Q01059

- Molecule 2 is a RNA chain called ferritin H IRE RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	30	Total	C	N	O	P	19	0	0
			621	275	106	210	30			
2	D	30	Total	C	N	O	P	19	0	0
			621	275	106	210	30			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	259	Total	O	0	0
			259	259		
3	B	226	Total	O	0	0
			226	226		
3	C	40	Total	O	0	0
			40	40		
3	D	32	Total	O	0	0
			32	32		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	109.56Å 80.85Å 142.87Å 90.00° 92.03° 90.00°	Depositor
Resolution (Å)	99.00 – 2.80 99.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	91.1 (99.00-2.80) 98.7 (99.00-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.69Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.198 , 0.218 0.217 , 0.230	Depositor DCC
R_{free} test set	3442 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	55.4	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 66.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.169 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15113	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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

5.1 Standard geometry

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	0.48	1/6813 (0.0%)	1.00	34/9249 (0.4%)
1	B	0.46	1/6813 (0.0%)	1.00	37/9249 (0.4%)
2	C	0.45	1/691 (0.1%)	0.94	1/1072 (0.1%)
2	D	0.43	1/691 (0.1%)	0.94	3/1072 (0.3%)
All	All	0.47	4/15008 (0.0%)	0.99	75/20642 (0.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	1
2	D	0	3
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	6	U	O3'-P	-5.72	1.52	1.61
2	D	6	U	O3'-P	-5.57	1.52	1.61
1	B	707	ASN	CG-OD1	5.39	1.33	1.23
1	A	707	ASN	CG-OD1	5.34	1.33	1.23

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	19	U	C2'-C3'-O3'	10.68	125.51	109.50
2	C	19	U	C2'-C3'-O3'	9.45	123.67	109.50
1	B	864	ARG	N-CA-C	7.86	122.89	110.70
1	A	464	LYS	CA-C-N	7.64	128.03	119.32
1	A	464	LYS	C-N-CA	7.64	128.03	119.32
1	A	864	ARG	N-CA-C	7.61	122.50	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	865	PHE	N-CA-C	-7.60	96.14	108.52
1	B	464	LYS	CA-C-N	7.57	127.94	119.32
1	B	464	LYS	C-N-CA	7.57	127.94	119.32
1	B	865	PHE	N-CA-C	-7.35	96.55	108.52
1	A	235	ILE	N-CA-C	-7.08	99.54	108.89
1	A	472	SER	CA-C-N	6.62	126.25	119.56
1	A	472	SER	C-N-CA	6.62	126.25	119.56
1	A	204	THR	N-CA-C	-6.60	105.54	112.93
1	A	529	VAL	N-CA-C	6.46	117.90	108.53
1	A	604	VAL	N-CA-C	6.38	116.53	110.53
1	B	235	ILE	N-CA-C	-6.38	100.47	108.89
1	B	529	VAL	N-CA-C	6.36	117.75	108.53
1	B	204	THR	N-CA-C	-6.29	105.88	112.93
1	B	252	HIS	CA-C-N	6.29	125.78	119.24
1	B	252	HIS	C-N-CA	6.29	125.78	119.24
1	B	472	SER	CA-C-N	6.28	126.40	119.87
1	B	472	SER	C-N-CA	6.28	126.40	119.87
1	A	252	HIS	CA-C-N	6.23	125.72	119.24
1	A	252	HIS	C-N-CA	6.23	125.72	119.24
1	B	604	VAL	N-CA-C	6.17	116.33	110.53
1	B	241	GLN	N-CA-C	-6.17	102.22	110.55
1	A	241	GLN	N-CA-C	-6.10	102.32	110.55
1	B	740	GLN	N-CA-C	6.09	117.15	108.74
1	B	814	ILE	N-CA-C	5.94	113.76	107.76
1	A	740	GLN	N-CA-C	5.93	116.93	108.74
1	B	752	VAL	N-CA-C	5.88	116.54	110.36
1	B	798	GLU	N-CA-C	-5.82	104.90	112.23
1	B	12	LEU	N-CA-C	-5.78	105.06	111.36
1	A	798	GLU	N-CA-C	-5.77	104.96	112.23
1	A	752	VAL	N-CA-C	5.70	116.35	110.36
1	A	649	PHE	N-CA-C	5.61	119.44	112.59
1	B	663	VAL	N-CA-C	5.61	115.96	108.11
1	B	747	GLY	N-CA-C	-5.58	107.62	115.32
1	A	240	PRO	N-CA-C	5.55	120.00	111.57
1	B	359	VAL	N-CA-C	-5.54	100.47	108.89
1	A	663	VAL	N-CA-C	5.52	115.83	108.11
1	B	649	PHE	N-CA-C	5.46	119.25	112.59
1	A	814	ILE	N-CA-C	5.46	113.27	107.76
1	A	12	LEU	N-CA-C	-5.44	105.43	111.36
1	B	113	LYS	N-CA-C	-5.38	105.97	112.54
1	A	869	VAL	N-CA-C	-5.38	105.48	110.53
1	B	869	VAL	N-CA-C	-5.38	105.48	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	169	ILE	N-CA-C	5.36	116.00	108.17
1	A	575	LEU	N-CA-C	-5.34	105.63	111.82
1	A	359	VAL	N-CA-C	-5.32	100.81	108.89
1	B	202	VAL	N-CA-C	-5.30	101.35	108.35
1	B	58	GLU	N-CA-C	-5.30	105.67	111.82
1	A	58	GLU	N-CA-C	-5.29	105.69	111.82
1	B	575	LEU	N-CA-C	-5.27	105.71	111.82
1	A	747	GLY	N-CA-C	-5.26	108.07	115.32
1	B	171	PRO	N-CA-C	5.26	117.11	110.70
1	A	202	VAL	N-CA-C	-5.22	101.46	108.35
1	A	169	ILE	N-CA-C	5.22	115.79	108.17
1	B	240	PRO	N-CA-C	5.21	119.49	111.57
1	B	239	LEU	N-CA-C	-5.18	102.35	109.71
1	A	113	LYS	N-CA-C	-5.18	106.22	112.54
1	A	171	PRO	N-CA-C	5.16	116.99	110.70
1	A	839	ILE	N-CA-C	5.12	113.03	108.63
1	B	125	ASP	N-CA-C	5.10	118.99	112.26
1	B	801	GLU	N-CA-C	-5.08	101.67	109.76
1	A	125	ASP	N-CA-C	5.08	118.97	112.26
1	A	866	ASP	N-CA-C	5.06	119.51	113.23
2	D	19	U	C4'-C3'-O3'	5.06	116.99	109.40
1	B	306	THR	N-CA-C	-5.06	105.68	111.14
2	D	7	G	N9-C1'-C2'	5.05	121.58	114.00
1	B	633	LEU	N-CA-C	-5.03	101.20	109.40
1	A	633	LEU	N-CA-C	-5.03	101.21	109.40
1	B	405	ALA	CA-C-N	5.01	126.11	119.84
1	B	405	ALA	C-N-CA	5.01	126.11	119.84

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	16	G	Sidechain
2	D	16	G	Sidechain
2	D	18	G	Sidechain
2	D	20	U	Sidechain

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	6657	0	6662	213	0
1	B	6657	0	6662	223	0
2	C	621	0	311	8	0
2	D	621	0	311	9	0
3	A	259	0	0	11	0
3	B	226	0	0	18	0
3	C	40	0	0	2	0
3	D	32	0	0	0	0
All	All	15113	0	13946	442	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 16.

All (442) 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:A:450:GLY:HA3	1:A:487:VAL:HG11	1.44	1.00
1:B:450:GLY:HA3	1:B:487:VAL:HG11	1.44	0.99
1:B:537:ASN:HD21	1:B:540:GLY:HA3	1.39	0.88
1:A:537:ASN:HD21	1:A:540:GLY:HA3	1.40	0.87
1:A:158:LYS:HD2	1:A:623:TRP:CE2	2.12	0.84
1:B:158:LYS:HD2	1:B:623:TRP:CE2	2.12	0.83
1:B:704:ARG:NH2	2:D:29:A:H4'	1.92	0.83
1:B:120:VAL:HG22	1:B:167:MET:HG3	1.64	0.80
1:A:120:VAL:HG22	1:A:167:MET:HG3	1.64	0.80
1:A:773:GLU:H	1:A:799:SER:HB3	1.46	0.80
1:B:704:ARG:HH21	2:D:29:A:H4'	1.50	0.77
1:A:809:VAL:HG11	1:A:870:GLU:HG2	1.64	0.77
1:B:773:GLU:H	1:B:799:SER:HB3	1.48	0.76
1:B:809:VAL:HG11	1:B:870:GLU:HG2	1.64	0.76
1:A:513:LEU:HD12	1:A:513:LEU:H	1.51	0.76
1:B:176:ILE:HD13	1:B:176:ILE:H	1.52	0.75
1:B:513:LEU:H	1:B:513:LEU:HD12	1.52	0.75
1:A:176:ILE:H	1:A:176:ILE:HD13	1.53	0.74
1:B:2:SER:HB3	3:B:1204:HOH:O	1.87	0.73
1:A:531:VAL:HG12	1:A:557:VAL:HG22	1.71	0.71
1:A:364:LYS:NZ	1:A:364:LYS:HB3	2.06	0.71
1:B:450:GLY:CA	1:B:487:VAL:HG11	2.19	0.71
1:A:450:GLY:CA	1:A:487:VAL:HG11	2.20	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:LYS:NZ	1:B:364:LYS:HB3	2.07	0.69
1:B:536:ARG:HH22	2:D:20:U:H4'	1.57	0.69
1:B:347:SER:O	1:B:349:PRO:HD3	1.93	0.69
1:B:531:VAL:HG12	1:B:557:VAL:HG22	1.74	0.68
1:B:863:ILE:HD11	1:B:865:PHE:CE2	2.29	0.68
2:C:29:A:O2'	2:C:30:C:H5'	1.94	0.68
1:A:721:ARG:HG3	1:A:721:ARG:HH11	1.59	0.68
1:A:803:ILE:HD12	1:A:803:ILE:H	1.60	0.67
1:B:803:ILE:HD12	1:B:803:ILE:H	1.60	0.67
1:A:863:ILE:HD11	1:A:865:PHE:CE2	2.30	0.67
1:A:347:SER:O	1:A:349:PRO:HD3	1.95	0.67
1:B:48:ARG:NH2	3:B:1163:HOH:O	2.27	0.66
1:B:721:ARG:HG3	1:B:721:ARG:HH11	1.61	0.66
1:A:707:ASN:HD22	1:A:708:SER:H	1.42	0.66
1:A:168:ARG:HD2	3:A:1241:HOH:O	1.96	0.65
1:B:533:SER:HB2	1:B:557:VAL:HG21	1.78	0.65
1:A:233:GLN:OE1	1:A:234:PRO:HD2	1.96	0.65
1:B:233:GLN:OE1	1:B:234:PRO:HD2	1.97	0.65
1:B:168:ARG:HD2	3:B:1201:HOH:O	1.97	0.64
1:A:533:SER:HB2	1:A:557:VAL:HG21	1.80	0.64
1:A:93:SER:HB2	3:A:1163:HOH:O	1.97	0.62
1:B:536:ARG:NH2	2:D:20:U:H4'	2.14	0.62
1:B:375:ARG:HD3	1:B:399:PHE:CD2	2.33	0.62
1:A:375:ARG:HD3	1:A:399:PHE:CD2	2.34	0.62
1:B:707:ASN:HD22	1:B:708:SER:H	1.46	0.62
1:A:518:VAL:HG13	1:A:545:ASN:ND2	2.16	0.61
1:A:866:ASP:HB3	3:A:1230:HOH:O	2.00	0.61
1:A:696:LEU:HD22	1:A:701:LEU:HD12	1.81	0.61
1:B:609:PHE:O	1:B:613:TYR:HD2	1.84	0.60
1:B:745:PRO:HG3	3:B:1223:HOH:O	2.01	0.60
1:B:454:LYS:HA	1:B:490:TYR:CD2	2.37	0.60
1:A:454:LYS:HA	1:A:490:TYR:CD2	2.37	0.60
1:A:743:HIS:CD2	1:A:745:PRO:HD2	2.37	0.60
1:B:180:VAL:HG11	1:B:612:VAL:HG21	1.84	0.60
1:A:180:VAL:HG11	1:A:612:VAL:HG21	1.84	0.59
1:B:93:SER:HB2	3:B:1018:HOH:O	2.01	0.59
1:A:776:SER:HA	1:A:801:GLU:HG3	1.82	0.59
1:B:814:ILE:HG23	1:B:863:ILE:HD13	1.83	0.59
1:A:765:PRO:HA	3:A:1159:HOH:O	2.01	0.59
1:A:814:ILE:HG23	1:A:863:ILE:HD13	1.83	0.59
1:B:518:VAL:HG13	1:B:545:ASN:ND2	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:609:PHE:O	1:A:613:TYR:HD2	1.86	0.59
1:B:776:SER:HA	1:B:801:GLU:HG3	1.82	0.59
1:B:696:LEU:HD22	1:B:701:LEU:HD12	1.84	0.59
1:A:803:ILE:HD12	1:A:803:ILE:N	2.18	0.58
1:A:411:ASP:O	1:A:412:HIS:HB3	2.03	0.58
1:A:707:ASN:ND2	1:A:708:SER:H	2.01	0.58
1:B:803:ILE:HD12	1:B:803:ILE:N	2.19	0.58
1:A:146:GLU:HG3	1:A:149:ARG:HD2	1.83	0.58
1:A:850:GLN:NE2	1:A:852:LYS:HB3	2.18	0.58
1:A:78:PHE:HA	3:A:1137:HOH:O	2.04	0.58
1:A:727:ILE:HA	1:A:739:PRO:HD3	1.86	0.58
1:B:415:PHE:HE2	1:B:424:LEU:HD13	1.68	0.58
1:A:758:ARG:HG2	1:A:762:GLU:OE1	2.04	0.58
1:B:300:CYS:HA	1:B:303:TYR:CZ	2.39	0.58
2:D:3:U:O2'	2:D:4:C:H5'	2.03	0.58
1:B:411:ASP:O	1:B:412:HIS:HB3	2.04	0.58
1:B:758:ARG:HG2	1:B:762:GLU:OE1	2.04	0.58
1:A:850:GLN:HE21	1:A:852:LYS:HB3	1.69	0.58
1:B:45:ALA:HA	3:B:1163:HOH:O	2.02	0.58
1:B:146:GLU:HG3	1:B:149:ARG:HD2	1.83	0.58
1:B:850:GLN:NE2	1:B:852:LYS:HB3	2.19	0.58
1:A:415:PHE:HE2	1:A:424:LEU:HD13	1.69	0.57
1:B:816:LEU:HD11	1:B:837:ILE:HD13	1.86	0.57
1:B:850:GLN:HE21	1:B:852:LYS:HB3	1.69	0.57
1:B:552:ALA:HB1	1:B:556:LEU:HD23	1.86	0.57
1:B:743:HIS:CD2	1:B:745:PRO:HD2	2.40	0.57
1:A:621:ALA:C	1:A:623:TRP:H	2.12	0.57
1:B:537:ASN:ND2	1:B:540:GLY:HA3	2.14	0.57
1:A:552:ALA:HB1	1:A:556:LEU:HD23	1.86	0.57
1:B:592:THR:OG1	1:B:595:GLU:HG3	2.04	0.57
1:A:574:PRO:HB3	1:A:583:GLN:HB2	1.86	0.56
1:A:816:LEU:HD11	1:A:837:ILE:HD13	1.87	0.56
1:A:537:ASN:ND2	1:A:540:GLY:HA3	2.15	0.56
1:A:592:THR:OG1	1:A:595:GLU:HG3	2.05	0.56
1:B:574:PRO:HB3	1:B:583:GLN:HB2	1.87	0.56
1:B:621:ALA:C	1:B:623:TRP:H	2.13	0.56
1:B:727:ILE:HA	1:B:739:PRO:HD3	1.88	0.56
1:B:707:ASN:ND2	1:B:708:SER:H	2.04	0.56
1:A:319:TYR:O	1:A:323:THR:HG23	2.07	0.55
1:B:542:VAL:HG12	1:B:542:VAL:O	2.06	0.55
1:A:300:CYS:HA	1:A:303:TYR:CZ	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:TYR:O	1:B:323:THR:HG23	2.08	0.54
1:A:542:VAL:HG12	1:A:542:VAL:O	2.07	0.54
1:B:6:ALA:HB1	3:B:1068:HOH:O	2.08	0.54
1:B:180:VAL:CG1	1:B:612:VAL:HG21	2.38	0.54
1:A:180:VAL:CG1	1:A:612:VAL:HG21	2.38	0.54
1:A:536:ARG:HH22	2:C:20:U:H4'	1.72	0.53
1:A:569:ASP:OD2	1:A:572:LYS:HG2	2.09	0.53
1:B:189:VAL:HG21	1:B:333:ILE:HG23	1.91	0.53
1:B:362:ASP:O	1:B:365:THR:HB	2.09	0.53
1:A:746:SER:OG	1:A:748:GLU:HB2	2.08	0.53
1:A:819:LEU:HD11	1:A:860:GLN:HB2	1.90	0.53
1:B:325:ARG:HD3	3:B:1142:HOH:O	2.09	0.53
1:B:863:ILE:HD11	1:B:865:PHE:CD2	2.43	0.53
1:A:578:ASN:C	1:A:580:LYS:N	2.65	0.53
1:B:569:ASP:OD2	1:B:572:LYS:HG2	2.09	0.53
1:A:847:MET:O	1:A:863:ILE:HG22	2.09	0.52
1:B:513:LEU:HD13	1:B:518:VAL:HG22	1.91	0.52
1:A:100:MET:CE	1:A:883:MET:HG2	2.39	0.52
1:B:80:PRO:HB3	1:B:201:LEU:HA	1.91	0.52
1:B:578:ASN:C	1:B:580:LYS:N	2.65	0.52
1:B:390:GLU:HB3	3:B:1212:HOH:O	2.08	0.52
1:A:362:ASP:O	1:A:365:THR:HB	2.10	0.52
1:A:594:GLU:N	1:A:594:GLU:OE2	2.43	0.52
1:B:484:GLU:HG2	1:B:603:TYR:O	2.10	0.52
1:B:594:GLU:OE2	1:B:594:GLU:N	2.43	0.52
1:A:80:PRO:HB3	1:A:201:LEU:HA	1.92	0.52
1:A:513:LEU:HD13	1:A:518:VAL:HG22	1.92	0.52
1:A:570:PHE:HA	3:A:1168:HOH:O	2.09	0.52
1:B:578:ASN:C	1:B:580:LYS:H	2.16	0.52
1:B:746:SER:OG	1:B:748:GLU:HB2	2.09	0.52
1:B:100:MET:HE3	1:B:883:MET:CG	2.40	0.52
1:A:578:ASN:C	1:A:580:LYS:H	2.16	0.51
1:B:96:ASP:OD2	1:B:227:GLU:OE1	2.27	0.51
1:A:385:MET:HG3	1:A:560:TYR:OH	2.10	0.51
1:A:300:CYS:N	1:A:301:PRO:CD	2.73	0.51
1:A:863:ILE:HD11	1:A:865:PHE:CD2	2.45	0.51
1:B:189:VAL:CG2	1:B:333:ILE:HG23	2.41	0.51
1:A:764:HIS:ND1	3:A:1227:HOH:O	2.34	0.51
1:B:23:ASN:OD1	1:B:25:ASN:HB2	2.11	0.51
1:B:100:MET:CE	1:B:883:MET:HG2	2.41	0.51
1:B:819:LEU:HD11	1:B:860:GLN:HB2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:VAL:HG21	1:A:333:ILE:HG23	1.93	0.51
1:A:814:ILE:HG23	1:A:863:ILE:CD1	2.41	0.51
1:A:875:HIS:HE1	3:A:1224:HOH:O	1.94	0.51
1:B:300:CYS:N	1:B:301:PRO:CD	2.73	0.51
1:B:404:VAL:O	1:B:405:ALA:C	2.53	0.51
1:B:616:ILE:HG22	1:B:617:GLU:N	2.25	0.51
1:B:847:MET:O	1:B:863:ILE:HG22	2.11	0.51
1:A:616:ILE:HG22	1:A:617:GLU:N	2.25	0.51
1:A:412:HIS:HD1	1:A:412:HIS:C	2.19	0.51
1:B:412:HIS:C	1:B:412:HIS:HD1	2.19	0.51
1:B:517:VAL:O	1:B:521:ILE:HG13	2.11	0.51
1:B:385:MET:HG3	1:B:560:TYR:OH	2.11	0.50
1:A:86:GLN:HB2	1:A:205:ASP:HB2	1.93	0.50
1:A:371:SER:O	1:A:550:TYR:HA	2.10	0.50
1:B:371:SER:O	1:B:550:TYR:HA	2.10	0.50
1:B:78:PHE:HA	3:B:1160:HOH:O	2.11	0.50
1:B:313:ASP:HB2	3:B:1087:HOH:O	2.11	0.50
1:B:814:ILE:CG2	1:B:863:ILE:HD13	2.42	0.50
1:A:484:GLU:HG2	1:A:603:TYR:O	2.11	0.50
1:B:86:GLN:HB2	1:B:205:ASP:HB2	1.93	0.50
1:B:375:ARG:HD3	1:B:399:PHE:CE2	2.47	0.50
1:B:814:ILE:HG23	1:B:863:ILE:CD1	2.41	0.50
1:A:404:VAL:O	1:A:405:ALA:C	2.54	0.50
1:A:803:ILE:H	1:A:803:ILE:CD1	2.24	0.50
1:A:708:SER:HB3	2:C:7:G:H4'	1.93	0.50
1:A:814:ILE:CG2	1:A:863:ILE:HD13	2.42	0.50
1:B:52:LYS:NZ	3:B:1204:HOH:O	2.44	0.50
1:B:803:ILE:H	1:B:803:ILE:CD1	2.24	0.50
1:A:701:LEU:HA	1:A:705:GLU:OE2	2.11	0.50
1:B:91:VAL:N	1:B:92:PRO:HD2	2.27	0.50
1:B:721:ARG:HG3	1:B:721:ARG:NH1	2.27	0.50
1:A:96:ASP:OD2	1:A:227:GLU:OE1	2.29	0.49
1:B:454:LYS:HG3	1:B:490:TYR:CE2	2.47	0.49
1:A:364:LYS:HB3	1:A:364:LYS:HZ3	1.76	0.49
1:A:375:ARG:HD3	1:A:399:PHE:CE2	2.47	0.49
1:A:707:ASN:HD22	1:A:708:SER:N	2.09	0.49
1:A:733:PHE:HZ	1:A:831:GLY:O	1.94	0.49
1:A:100:MET:HE3	1:A:883:MET:CG	2.42	0.49
1:A:769:LEU:HD23	1:A:796:LEU:HB3	1.95	0.49
1:A:189:VAL:CG2	1:A:333:ILE:HG23	2.43	0.49
1:A:517:VAL:O	1:A:521:ILE:HG13	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ASN:OD1	1:A:25:ASN:HB2	2.13	0.49
1:B:785:LYS:HE3	3:B:1166:HOH:O	2.12	0.49
1:B:364:LYS:HB3	1:B:364:LYS:HZ3	1.76	0.48
1:B:375:ARG:HB2	1:B:378:ASP:OD2	2.13	0.48
1:A:364:LYS:HB3	1:A:364:LYS:HZ2	1.76	0.48
1:A:518:VAL:HG13	1:A:545:ASN:HD21	1.78	0.48
1:B:252:HIS:HB3	3:B:1119:HOH:O	2.11	0.48
1:B:703:PRO:HA	1:B:706:PHE:CE2	2.49	0.48
1:A:616:ILE:O	1:A:619:VAL:HG12	2.14	0.48
1:B:616:ILE:O	1:B:619:VAL:HG12	2.14	0.48
1:A:91:VAL:N	1:A:92:PRO:HD2	2.29	0.48
1:B:769:LEU:HD23	1:B:796:LEU:HB3	1.96	0.48
1:A:536:ARG:NH2	2:C:20:U:H4'	2.29	0.48
1:A:703:PRO:HA	1:A:706:PHE:CE2	2.49	0.48
1:A:115:ASN:ND2	1:A:116:PRO:HD2	2.28	0.47
1:B:463:VAL:CG2	1:B:562:ILE:HD13	2.44	0.47
1:B:314:GLU:O	1:B:318:LYS:HG3	2.14	0.47
1:B:518:VAL:HG13	1:B:545:ASN:HD21	1.80	0.47
1:B:729:LEU:O	1:B:737:GLN:HA	2.15	0.47
1:B:396:LYS:HD3	1:B:403:GLN:OE1	2.14	0.47
1:B:701:LEU:HA	1:B:705:GLU:OE2	2.13	0.47
1:A:850:GLN:HG2	1:A:851:VAL:N	2.29	0.47
1:B:733:PHE:HZ	1:B:831:GLY:O	1.96	0.47
1:A:454:LYS:HG3	1:A:490:TYR:CE2	2.50	0.47
1:A:727:ILE:HG23	3:A:1182:HOH:O	2.15	0.47
1:B:691:PRO:O	1:B:721:ARG:HD3	2.14	0.47
1:A:375:ARG:HB2	1:A:378:ASP:OD2	2.15	0.47
1:A:729:LEU:O	1:A:737:GLN:HA	2.15	0.47
1:B:850:GLN:HG2	1:B:851:VAL:N	2.30	0.47
1:A:463:VAL:CG2	1:A:562:ILE:HD13	2.44	0.47
1:B:364:LYS:HB3	1:B:364:LYS:HZ2	1.78	0.47
1:A:187:ARG:O	1:A:188:VAL:HB	2.15	0.47
1:A:314:GLU:O	1:A:318:LYS:HG3	2.15	0.47
1:B:531:VAL:CG1	1:B:557:VAL:HG22	2.44	0.47
1:A:13:ASP:HB2	1:A:335:LYS:NZ	2.30	0.46
1:A:344:ARG:HG2	1:A:346:TYR:CE2	2.51	0.46
1:A:408:HIS:HA	1:A:411:ASP:OD1	2.16	0.46
1:B:28:ASP:HB3	1:B:64:LEU:HD21	1.96	0.46
1:B:56:LYS:O	1:B:59:ASP:HB2	2.16	0.46
1:B:450:GLY:CA	1:B:487:VAL:CG1	2.93	0.46
1:B:707:ASN:HD22	1:B:708:SER:N	2.12	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:LYS:O	1:A:59:ASP:HB2	2.16	0.46
1:A:702:THR:OG1	1:A:705:GLU:HG2	2.15	0.46
1:B:100:MET:HE3	1:B:883:MET:HG2	1.98	0.46
1:B:344:ARG:HG2	1:B:346:TYR:CE2	2.51	0.46
1:B:408:HIS:HA	1:B:411:ASP:OD1	2.16	0.46
1:A:464:LYS:HA	1:A:465:PRO:HD3	1.81	0.46
1:B:45:ALA:CB	3:B:1163:HOH:O	2.64	0.46
1:B:464:LYS:HA	1:B:465:PRO:HD3	1.81	0.46
1:A:450:GLY:CA	1:A:487:VAL:CG1	2.93	0.46
1:B:212:ASP:HA	1:B:216:VAL:O	2.16	0.46
1:A:472:SER:HB3	1:A:499:VAL:HG23	1.97	0.46
1:B:62:ASN:HB3	1:B:74:ILE:HD11	1.97	0.46
1:B:115:ASN:ND2	1:B:116:PRO:HD2	2.31	0.46
1:B:472:SER:HB3	1:B:499:VAL:HG23	1.98	0.46
1:A:212:ASP:HA	1:A:216:VAL:O	2.16	0.46
1:A:393:LEU:HD21	1:A:527:VAL:HG11	1.98	0.46
1:A:405:ALA:HB3	1:A:408:HIS:HD2	1.81	0.46
1:B:393:LEU:HD21	1:B:527:VAL:HG11	1.98	0.46
1:B:702:THR:OG1	1:B:705:GLU:HG2	2.15	0.46
1:A:535:ASN:ND2	3:A:1186:HOH:O	2.48	0.46
1:A:396:LYS:HD3	1:A:403:GLN:OE1	2.16	0.45
1:A:605:ILE:HB	1:A:606:PRO:HD3	1.98	0.45
1:B:405:ALA:HB3	1:B:408:HIS:HD2	1.81	0.45
1:A:531:VAL:HG11	1:A:557:VAL:HA	1.99	0.45
2:C:15:A:H4'	2:C:16:G:O5'	2.15	0.45
1:B:13:ASP:HB2	1:B:335:LYS:NZ	2.31	0.45
1:A:531:VAL:CG1	1:A:557:VAL:HG22	2.42	0.45
1:B:605:ILE:HB	1:B:606:PRO:HD3	1.99	0.45
2:C:22:G:N7	3:C:1007:HOH:O	2.36	0.45
1:A:780:ARG:C	3:C:1026:HOH:O	2.60	0.45
1:A:847:MET:O	1:A:863:ILE:CG2	2.64	0.45
1:B:245:TYR:HD1	1:B:268:LEU:HD21	1.82	0.45
1:A:515:GLU:N	1:A:516:PRO:CD	2.80	0.45
1:B:188:VAL:HG21	1:B:320:LEU:HD21	1.99	0.45
1:B:463:VAL:HG23	1:B:562:ILE:HD13	1.99	0.45
1:B:515:GLU:N	1:B:516:PRO:CD	2.80	0.45
1:B:847:MET:O	1:B:863:ILE:CG2	2.65	0.45
1:A:100:MET:HE3	1:A:883:MET:HG2	1.99	0.44
1:A:575:LEU:CD1	1:A:589:ILE:HD11	2.47	0.44
1:B:187:ARG:O	1:B:188:VAL:HB	2.17	0.44
1:B:246:ARG:HA	1:B:281:PHE:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:593:ARG:O	1:B:597:GLN:HG3	2.18	0.44
1:A:392:CYS:O	1:A:401:GLY:HA2	2.17	0.44
1:A:621:ALA:O	1:A:623:TRP:N	2.51	0.44
1:A:62:ASN:HB3	1:A:74:ILE:HD11	1.98	0.44
1:A:594:GLU:CD	1:A:594:GLU:H	2.26	0.44
1:A:642:TYR:C	1:A:643:ILE:HD12	2.43	0.44
1:A:643:ILE:HD12	1:A:643:ILE:N	2.32	0.44
1:A:691:PRO:O	1:A:721:ARG:HD3	2.17	0.44
1:B:146:GLU:HG3	1:B:149:ARG:HB2	1.99	0.44
1:B:392:CYS:O	1:B:401:GLY:HA2	2.17	0.44
1:B:643:ILE:HD12	1:B:643:ILE:N	2.33	0.44
1:A:538:PHE:C	1:A:540:GLY:H	2.26	0.44
1:A:327:GLU:O	1:A:331:LYS:HG3	2.17	0.44
1:A:404:VAL:HG23	1:A:409:HIS:CD2	2.53	0.44
1:A:621:ALA:C	1:A:623:TRP:N	2.76	0.44
1:A:721:ARG:HG3	1:A:721:ARG:NH1	2.26	0.44
1:B:250:LYS:HG3	1:B:251:PRO:HD2	2.00	0.44
1:B:531:VAL:HG11	1:B:557:VAL:HA	2.00	0.44
1:B:538:PHE:C	1:B:540:GLY:H	2.26	0.44
1:A:463:VAL:HG23	1:A:562:ILE:HD13	2.00	0.44
1:B:415:PHE:CE2	1:B:424:LEU:HD13	2.51	0.44
1:B:594:GLU:H	1:B:594:GLU:CD	2.26	0.44
1:A:146:GLU:HG3	1:A:149:ARG:HB2	1.99	0.44
1:A:424:LEU:O	1:A:526:LEU:HD13	2.18	0.43
1:B:568:ILE:HG12	1:B:569:ASP:N	2.33	0.43
1:B:642:TYR:C	1:B:643:ILE:HD12	2.43	0.43
1:A:568:ILE:HG12	1:A:569:ASP:N	2.33	0.43
1:A:593:ARG:O	1:A:597:GLN:HG3	2.18	0.43
1:B:266:LYS:HE2	1:B:270:GLN:OE1	2.18	0.43
1:B:424:LEU:O	1:B:526:LEU:HD13	2.18	0.43
1:A:708:SER:CB	2:C:7:G:H4'	2.49	0.43
1:B:62:ASN:HB3	1:B:74:ILE:CD1	2.48	0.43
1:B:110:ASP:HB3	1:B:113:LYS:HG2	1.99	0.43
1:B:190:PHE:O	1:B:196:TYR:HA	2.18	0.43
1:B:575:LEU:CD1	1:B:589:ILE:HD11	2.48	0.43
1:A:266:LYS:HE2	1:A:270:GLN:OE1	2.19	0.43
1:A:415:PHE:CE2	1:A:424:LEU:HD13	2.51	0.43
1:B:367:VAL:O	1:B:367:VAL:HG23	2.18	0.43
1:B:621:ALA:O	1:B:623:TRP:N	2.52	0.43
1:A:62:ASN:HB3	1:A:74:ILE:CD1	2.49	0.43
1:A:245:TYR:HD1	1:A:268:LEU:HD21	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:621:ALA:C	1:B:623:TRP:N	2.77	0.43
1:A:835:TYR:CD1	1:A:835:TYR:N	2.87	0.43
1:B:835:TYR:CD1	1:B:835:TYR:N	2.86	0.43
1:A:512:PRO:N	1:A:543:HIS:HE2	2.17	0.43
1:A:652:LEU:C	1:A:652:LEU:HD23	2.44	0.43
1:B:404:VAL:HG23	1:B:409:HIS:CD2	2.54	0.43
1:A:188:VAL:HG21	1:A:320:LEU:HD21	2.01	0.43
1:A:78:PHE:O	1:A:80:PRO:HD3	2.18	0.43
1:A:426:HIS:CD2	1:A:527:VAL:HG12	2.53	0.43
1:B:243:ILE:HG13	1:B:273:VAL:HG12	2.00	0.43
1:B:247:LEU:HB3	1:B:284:GLY:HA3	2.01	0.43
1:B:367:VAL:O	1:B:367:VAL:CG2	2.67	0.43
1:B:396:LYS:O	1:B:397:GLN:C	2.62	0.43
1:B:704:ARG:HH21	2:D:29:A:C4'	2.24	0.43
1:A:404:VAL:HG23	1:A:409:HIS:HD2	1.83	0.42
1:A:416:ILE:HD12	1:A:416:ILE:N	2.34	0.42
1:B:100:MET:HE3	1:B:883:MET:HG3	2.01	0.42
1:A:352:ASP:HA	1:A:353:PRO:HD3	1.77	0.42
1:A:364:LYS:NZ	1:A:364:LYS:CB	2.78	0.42
1:A:743:HIS:NE2	1:A:745:PRO:HD2	2.34	0.42
1:B:48:ARG:HD2	3:B:1102:HOH:O	2.19	0.42
1:A:110:ASP:HB3	1:A:113:LYS:HG2	2.00	0.42
1:A:28:ASP:HB3	1:A:64:LEU:HD21	2.00	0.42
1:B:229:VAL:CG2	1:B:235:ILE:CD1	2.97	0.42
1:B:512:PRO:N	1:B:543:HIS:HE2	2.17	0.42
1:B:578:ASN:O	1:B:580:LYS:N	2.52	0.42
1:A:405:ALA:O	1:A:408:HIS:N	2.50	0.42
1:A:652:LEU:HD23	1:A:653:THR:N	2.34	0.42
1:B:381:ALA:HB1	3:B:1224:HOH:O	2.20	0.42
1:A:246:ARG:HA	1:A:281:PHE:O	2.19	0.42
1:A:396:LYS:O	1:A:397:GLN:C	2.62	0.42
1:A:578:ASN:O	1:A:580:LYS:N	2.52	0.42
1:B:352:ASP:HA	1:B:353:PRO:HD3	1.78	0.42
1:B:416:ILE:N	1:B:416:ILE:HD12	2.34	0.42
1:A:158:LYS:HD2	1:A:623:TRP:NE1	2.32	0.42
1:A:250:LYS:HG3	1:A:251:PRO:HD2	2.02	0.42
1:B:81:ALA:HA	3:B:1025:HOH:O	2.20	0.42
1:B:327:GLU:O	1:B:331:LYS:HG3	2.19	0.42
1:B:652:LEU:HD23	1:B:652:LEU:C	2.45	0.42
1:A:100:MET:HE1	1:A:883:MET:HG2	2.02	0.42
1:A:735:ASN:O	1:A:736:LYS:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:LYS:HD2	1:B:623:TRP:NE1	2.32	0.42
1:B:652:LEU:HD23	1:B:653:THR:N	2.34	0.42
1:A:769:LEU:CD2	1:A:796:LEU:HD23	2.50	0.42
1:A:782:TRP:CE3	1:A:785:LYS:HB2	2.55	0.42
1:B:426:HIS:CD2	1:B:527:VAL:HG12	2.54	0.42
1:B:733:PHE:C	1:B:734:LEU:HD23	2.45	0.42
1:A:385:MET:HE1	1:A:550:TYR:CD2	2.54	0.42
1:A:287:GLN:HG3	3:A:1162:HOH:O	2.19	0.41
1:A:697:THR:C	1:A:699:ARG:H	2.27	0.41
1:B:94:VAL:HG12	1:B:156:PHE:CZ	2.55	0.41
1:B:404:VAL:HG23	1:B:409:HIS:HD2	1.84	0.41
1:A:247:LEU:HB3	1:A:284:GLY:HA3	2.02	0.41
1:A:367:VAL:O	1:A:367:VAL:HG23	2.20	0.41
1:B:203:GLY:C	1:B:205:ASP:H	2.28	0.41
1:B:282:GLY:O	1:B:285:VAL:HG23	2.20	0.41
1:B:10:GLU:C	1:B:20:LYS:HG2	2.44	0.41
1:B:405:ALA:O	1:B:408:HIS:N	2.51	0.41
1:B:542:VAL:HG13	1:B:549:ASN:ND2	2.35	0.41
2:C:28:A:O2'	2:C:29:A:H5'	2.21	0.41
1:A:151:ARG:O	1:A:155:GLU:HG2	2.20	0.41
1:A:203:GLY:C	1:A:205:ASP:H	2.28	0.41
1:A:243:ILE:HG13	1:A:273:VAL:HG12	2.01	0.41
1:A:828:GLY:O	1:A:829:LEU:C	2.63	0.41
1:B:472:SER:HB3	1:B:499:VAL:CG2	2.51	0.41
1:B:828:GLY:O	1:B:829:LEU:C	2.63	0.41
1:A:367:VAL:O	1:A:367:VAL:CG2	2.69	0.41
1:A:413:LYS:HE3	1:A:523:GLN:HB3	2.02	0.41
1:A:733:PHE:C	1:A:734:LEU:HD23	2.45	0.41
1:B:295:THR:HG23	1:B:440:THR:OG1	2.20	0.41
1:B:371:SER:HA	1:B:378:ASP:O	2.20	0.41
1:B:536:ARG:HE	1:B:536:ARG:HB3	1.70	0.41
1:B:645:SER:HA	1:B:646:PRO:HD3	1.88	0.41
1:A:472:SER:HB3	1:A:499:VAL:CG2	2.51	0.41
1:A:542:VAL:HG13	1:A:549:ASN:ND2	2.35	0.41
1:A:578:ASN:N	1:A:582:GLN:O	2.53	0.41
1:B:733:PHE:O	1:B:734:LEU:HD23	2.21	0.41
1:A:170:ILE:HG23	1:A:616:ILE:HD11	2.03	0.41
1:A:577:THR:O	1:A:578:ASN:C	2.63	0.41
1:B:78:PHE:O	1:B:80:PRO:HD3	2.20	0.41
1:B:364:LYS:NZ	1:B:364:LYS:CB	2.79	0.41
1:A:295:THR:HG23	1:A:440:THR:OG1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:658:PRO:HA	1:A:659:PRO:HD3	1.97	0.41
1:B:697:THR:C	1:B:699:ARG:H	2.28	0.41
1:A:158:LYS:HD2	1:A:623:TRP:CZ2	2.55	0.41
1:A:371:SER:HA	1:A:378:ASP:O	2.21	0.41
1:A:412:HIS:C	1:A:412:HIS:ND1	2.78	0.41
1:A:536:ARG:HE	1:A:536:ARG:HB3	1.70	0.41
1:A:835:TYR:CE2	1:A:853:LEU:HD21	2.56	0.41
1:B:385:MET:HE1	1:B:550:TYR:CD2	2.55	0.41
1:B:402:PHE:O	1:B:403:GLN:O	2.39	0.41
1:B:577:THR:O	1:B:578:ASN:C	2.63	0.41
1:B:578:ASN:HD22	1:B:582:GLN:NE2	2.19	0.41
2:D:10:U:H2'	2:D:11:C:C6	2.55	0.41
1:A:82:ARG:NH2	1:A:186:ALA:HA	2.36	0.41
1:A:146:GLU:O	1:A:146:GLU:HG2	2.21	0.41
1:A:229:VAL:CG2	1:A:235:ILE:CD1	2.99	0.41
1:A:514:PRO:HB2	1:A:517:VAL:HG23	2.02	0.41
1:A:578:ASN:HD22	1:A:582:GLN:NE2	2.19	0.41
1:A:785:LYS:HG2	1:A:811:MET:HE2	2.03	0.41
1:B:170:ILE:HG23	1:B:616:ILE:HD11	2.03	0.41
1:B:412:HIS:C	1:B:412:HIS:ND1	2.78	0.41
1:B:514:PRO:HB2	1:B:517:VAL:HG23	2.02	0.41
1:A:370:CYS:HA	1:A:551:LEU:O	2.22	0.40
1:A:645:SER:HA	1:A:646:PRO:HD3	1.88	0.40
1:B:598:ALA:HA	1:B:601:ARG:NH1	2.36	0.40
1:B:608:MET:O	1:B:612:VAL:HG23	2.21	0.40
1:A:10:GLU:C	1:A:20:LYS:HG2	2.45	0.40
1:A:35:LEU:CD2	1:A:66:TRP:HB3	2.51	0.40
1:B:21:PHE:CD1	1:B:21:PHE:C	2.99	0.40
1:B:158:LYS:HD2	1:B:623:TRP:CZ2	2.55	0.40
1:B:171:PRO:HA	1:B:172:PRO:HD3	1.89	0.40
1:B:413:LYS:HE3	1:B:523:GLN:HB3	2.03	0.40
1:B:602:GLN:HG3	1:B:603:TYR:CE1	2.56	0.40
1:B:708:SER:HB3	2:D:7:G:H4'	2.03	0.40
1:B:34:ARG:HA	1:B:34:ARG:NE	2.37	0.40
1:B:683:ALA:HA	2:D:8:C:C5	2.56	0.40
1:A:402:PHE:O	1:A:403:GLN:O	2.39	0.40
1:A:464:LYS:N	1:A:464:LYS:HD2	2.36	0.40
1:B:447:LEU:HD22	1:B:451:LEU:HG	2.02	0.40
1:B:480:TYR:O	1:B:484:GLU:HB2	2.20	0.40
1:B:658:PRO:HA	1:B:659:PRO:HD3	1.99	0.40
1:B:735:ASN:O	1:B:736:LYS:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:769:LEU:CD2	1:B:796:LEU:HD23	2.51	0.40
1:A:602:GLN:HG3	1:A:603:TYR:CE1	2.56	0.40
1:A:743:HIS:CE1	1:A:745:PRO:HD2	2.56	0.40
1:A:773:GLU:N	1:A:799:SER:HB3	2.26	0.40
1:A:784:ALA:C	1:A:787:PRO:HD2	2.46	0.40
1:B:82:ARG:NH2	1:B:186:ALA:HA	2.37	0.40
1:B:488:MET:N	1:B:489:PRO:CD	2.85	0.40
1:B:782:TRP:CE3	1:B:785:LYS:HB2	2.57	0.40
1:B:805:ARG:HG2	1:B:805:ARG:HH11	1.86	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	842/908 (93%)	781 (93%)	49 (6%)	12 (1%)	9	30
1	B	842/908 (93%)	782 (93%)	48 (6%)	12 (1%)	9	30
All	All	1684/1816 (93%)	1563 (93%)	97 (6%)	24 (1%)	9	30

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	395	ALA
1	A	403	GLN
1	A	536	ARG
1	B	395	ALA
1	B	403	GLN
1	B	536	ARG
1	A	28	ASP
1	A	397	GLN
1	A	412	HIS

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Mol	Chain	Res	Type
1	A	622	SER
1	B	28	ASP
1	B	397	GLN
1	B	412	HIS
1	B	622	SER
1	A	537	ASN
1	A	539	GLU
1	A	648	PHE
1	A	698	ASN
1	B	537	ASN
1	B	539	GLU
1	B	648	PHE
1	B	698	ASN
1	A	406	PRO
1	B	406	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	725/772 (94%)	685 (94%)	40 (6%)	19	51
1	B	725/772 (94%)	684 (94%)	41 (6%)	18	49
All	All	1450/1544 (94%)	1369 (94%)	81 (6%)	19	50

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	67	ASN
1	A	72	MET
1	A	100	MET
1	A	120	VAL
1	A	126	HIS
1	A	176	ILE
1	A	214	LEU

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Mol	Chain	Res	Type
1	A	235	ILE
1	A	240	PRO
1	A	247	LEU
1	A	248	MET
1	A	354	ASP
1	A	364	LYS
1	A	365	THR
1	A	367	VAL
1	A	377	GLN
1	A	380	VAL
1	A	387	LYS
1	A	406	PRO
1	A	412	HIS
1	A	437	SER
1	A	447	LEU
1	A	452	LEU
1	A	484	GLU
1	A	513	LEU
1	A	515	GLU
1	A	527	VAL
1	A	543	HIS
1	A	582	GLN
1	A	610	THR
1	A	616	ILE
1	A	617	GLU
1	A	633	LEU
1	A	750	LEU
1	A	761	GLN
1	A	826	SER
1	A	852	LYS
1	A	858	THR
1	A	863	ILE
1	B	24	LEU
1	B	67	ASN
1	B	72	MET
1	B	100	MET
1	B	120	VAL
1	B	126	HIS
1	B	176	ILE
1	B	214	LEU
1	B	235	ILE
1	B	240	PRO

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Mol	Chain	Res	Type
1	B	247	LEU
1	B	248	MET
1	B	354	ASP
1	B	364	LYS
1	B	365	THR
1	B	367	VAL
1	B	377	GLN
1	B	380	VAL
1	B	387	LYS
1	B	406	PRO
1	B	412	HIS
1	B	437	SER
1	B	447	LEU
1	B	452	LEU
1	B	484	GLU
1	B	513	LEU
1	B	515	GLU
1	B	527	VAL
1	B	543	HIS
1	B	571	GLU
1	B	582	GLN
1	B	610	THR
1	B	616	ILE
1	B	617	GLU
1	B	633	LEU
1	B	750	LEU
1	B	761	GLN
1	B	826	SER
1	B	852	LYS
1	B	858	THR
1	B	863	ILE

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

Mol	Chain	Res	Type
1	A	67	ASN
1	A	71	HIS
1	A	179	GLN
1	A	252	HIS
1	A	338	GLN
1	A	377	GLN
1	A	397	GLN

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Mol	Chain	Res	Type
1	A	409	HIS
1	A	426	HIS
1	A	493	GLN
1	A	537	ASN
1	A	582	GLN
1	A	620	ASN
1	A	651	ASN
1	A	698	ASN
1	A	707	ASN
1	A	764	HIS
1	A	807	ASN
1	A	850	GLN
1	A	875	HIS
1	B	16	GLN
1	B	67	ASN
1	B	71	HIS
1	B	252	HIS
1	B	338	GLN
1	B	377	GLN
1	B	397	GLN
1	B	409	HIS
1	B	426	HIS
1	B	493	GLN
1	B	537	ASN
1	B	582	GLN
1	B	620	ASN
1	B	651	ASN
1	B	698	ASN
1	B	707	ASN
1	B	737	GLN
1	B	764	HIS
1	B	807	ASN
1	B	850	GLN
1	B	875	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	27/30 (90%)	3 (11%)	1 (3%)
2	D	27/30 (90%)	4 (14%)	1 (3%)
All	All	54/60 (90%)	7 (12%)	2 (3%)

All (7) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	9	U
2	C	19	U
2	C	20	U
2	D	8	C
2	D	9	U
2	D	19	U
2	D	20	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	19	U
2	D	19	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	850/908 (93%)	-0.10	9 (1%) 78 70	22, 50, 90, 124	0
1	B	850/908 (93%)	-0.09	9 (1%) 78 70	23, 50, 91, 124	0
2	C	30/30 (100%)	-0.34	0 100 100	16, 56, 85, 93	2 (6%)
2	D	30/30 (100%)	-0.32	0 100 100	16, 58, 81, 94	2 (6%)
All	All	1760/1876 (93%)	-0.10	18 (1%) 79 72	16, 50, 91, 124	4 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	542	VAL	4.1
1	A	538	PHE	3.4
1	B	14	PRO	3.1
1	A	542	VAL	3.0
1	B	538	PHE	3.0
1	B	512	PRO	2.9
1	B	18	GLY	2.9
1	A	512	PRO	2.8
1	A	618	THR	2.6
1	A	146	GLU	2.6
1	B	126	HIS	2.6
1	A	541	ARG	2.4
1	B	616	ILE	2.4
1	B	499	VAL	2.4
1	B	146	GLU	2.3
1	A	499	VAL	2.2
1	A	126	HIS	2.1
1	A	127	SER	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.