



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:01 PM UTC

PDB ID : 6JE9 / pdb_00006je9
Title : Crystal structure of Nme1Cas9-sgRNA dimer mediated by double protein inhibitor AcrIIC3 monomers
Authors : Sun, W.; Yang, J.; Cheng, Z.; Liu, C.; Wang, K.; Huang, X.; Wang, Y.
Deposited on : 2019-02-04
Resolution : 3.46 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:

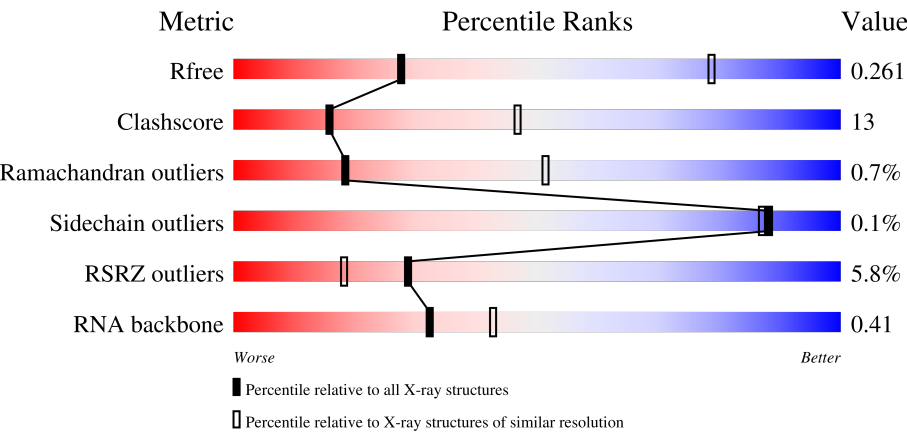
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 3.46 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	1070 (3.50-3.42)
Clashscore	190562	1128 (3.50-3.42)
Ramachandran outliers	187476	1101 (3.50-3.42)
Sidechain outliers	187428	1102 (3.50-3.42)
RSRZ outliers	180081	1070 (3.50-3.42)
RNA backbone	3983	1186 (3.92-3.00)

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	1092	4% 73% 23% • •
1	C	1092	8% 76% 18% • 5%
2	B	135	7% 28% 43% 14% 15%
2	D	135	7% 25% 44% 15% 16%

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Mol	Chain	Length	Quality of chain
3	E	117	% 71% 29%
3	F	117	3% 62% 38%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22158 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 CRISPR-associated endonuclease Cas9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1056	Total	C	N	O	S	0	0	0
			8035	5081	1462	1468	24			
1	C	1036	Total	C	N	O	S	0	0	0
			7535	4735	1390	1389	21			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1083	SER	-	expression tag	UNP C9X1G5
A	1084	GLU	-	expression tag	UNP C9X1G5
A	1085	HIS	-	expression tag	UNP C9X1G5
A	1086	HIS	-	expression tag	UNP C9X1G5
A	1087	HIS	-	expression tag	UNP C9X1G5
A	1088	HIS	-	expression tag	UNP C9X1G5
A	1089	HIS	-	expression tag	UNP C9X1G5
A	1090	HIS	-	expression tag	UNP C9X1G5
A	1091	HIS	-	expression tag	UNP C9X1G5
A	1092	HIS	-	expression tag	UNP C9X1G5
C	1083	SER	-	expression tag	UNP C9X1G5
C	1084	GLU	-	expression tag	UNP C9X1G5
C	1085	HIS	-	expression tag	UNP C9X1G5
C	1086	HIS	-	expression tag	UNP C9X1G5
C	1087	HIS	-	expression tag	UNP C9X1G5
C	1088	HIS	-	expression tag	UNP C9X1G5
C	1089	HIS	-	expression tag	UNP C9X1G5
C	1090	HIS	-	expression tag	UNP C9X1G5
C	1091	HIS	-	expression tag	UNP C9X1G5
C	1092	HIS	-	expression tag	UNP C9X1G5

- Molecule 2 is a RNA chain called sgRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	115	Total	C	N	O	P	0	0	0
			2433	1089	421	808	115			
2	D	113	Total	C	N	O	P	0	0	0
			2393	1071	416	793	113			

- Molecule 3 is a protein called AcrIIC3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	117	Total	C	N	O	S	0	0	0
			890	572	156	159	3			
3	F	117	Total	C	N	O	S	0	0	0
			872	560	152	158	2			

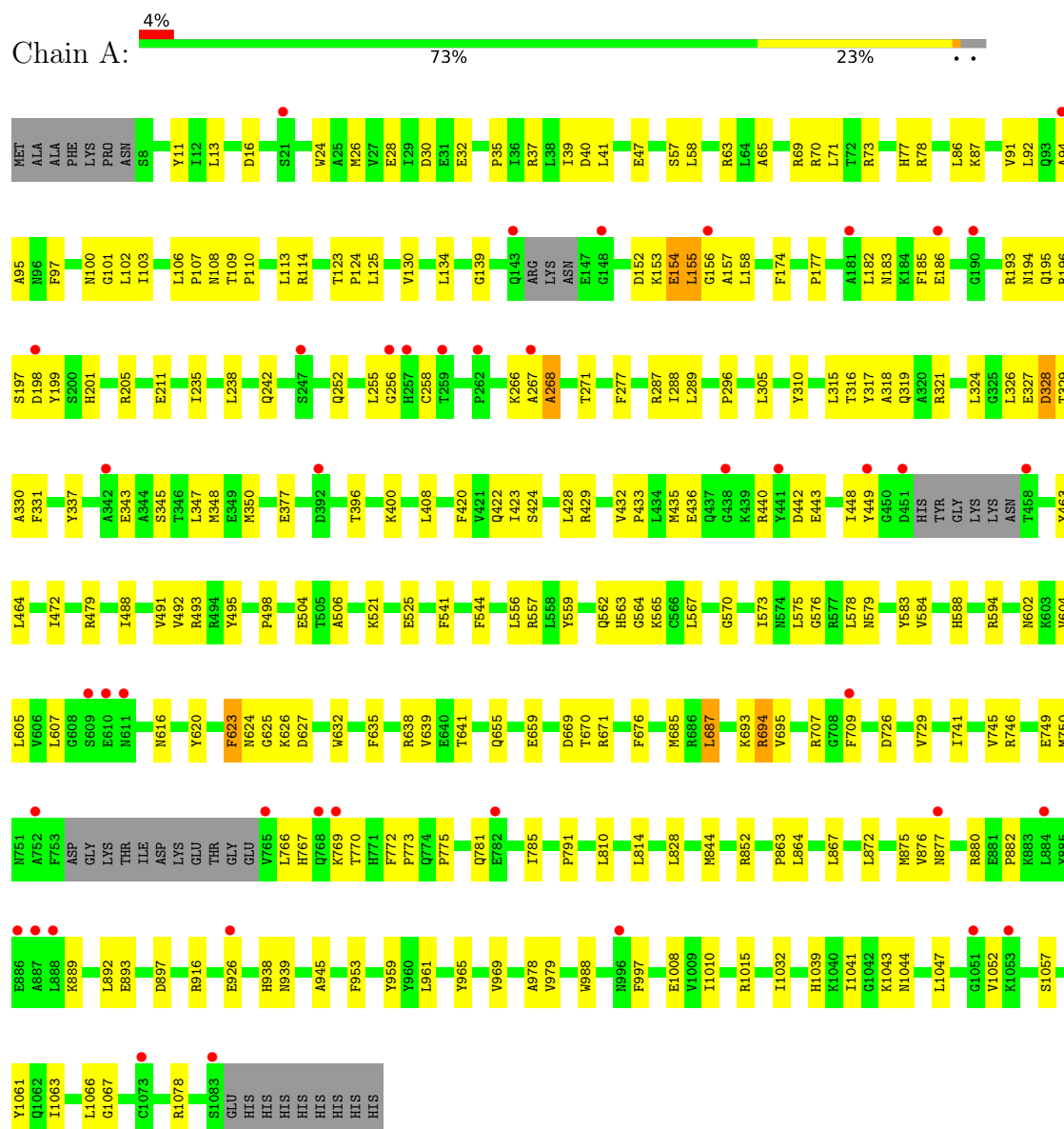
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	0	SER	-	expression tag	UNP A0A3E2QDI5
F	0	SER	-	expression tag	UNP A0A3E2QDI5

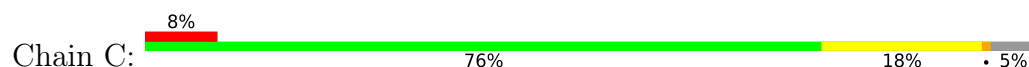
3 Residue-property plots

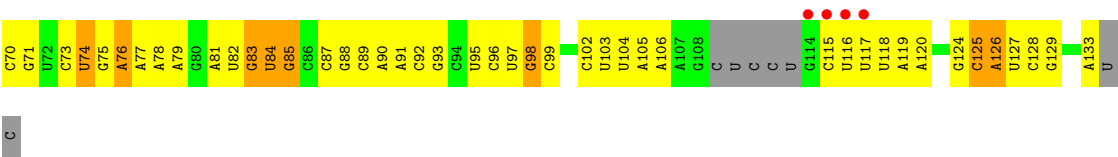
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: CRISPR-associated endonuclease Cas9

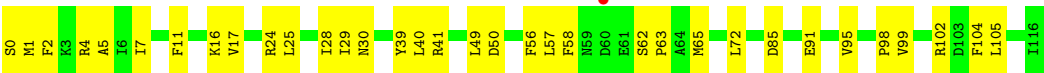


• Molecule 1: CRISPR-associated endonuclease Cas9

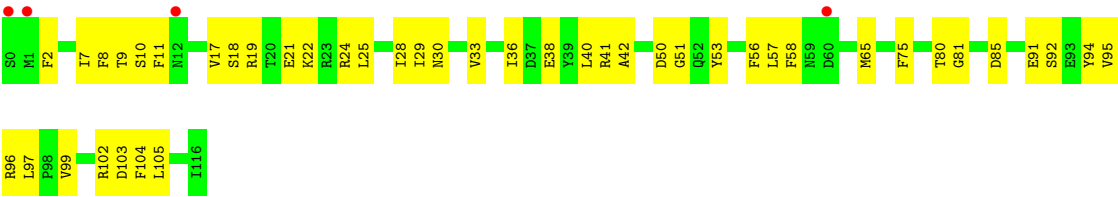




• Molecule 3: AcrIIC3



• Molecule 3: AcrIIC3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	112.84Å 156.20Å 151.60Å 90.00° 108.58° 90.00°	Depositor
Resolution (Å)	45.60 – 3.46 45.60 – 3.46	Depositor EDS
% Data completeness (in resolution range)	77.4 (45.60-3.46) 77.3 (45.60-3.46)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 3.48Å)	Xtriage
Refinement program	PHENIX (1.14_3247: ???)	Depositor
R, R_{free}	0.221 , 0.261 0.221 , 0.261	Depositor DCC
R_{free} test set	2463 reflections (3.76%)	wwPDB-VP
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	22158	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.19% 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	0.22	0/8191	0.57	6/11103 (0.1%)
1	C	0.21	2/7675 (0.0%)	0.51	2/10440 (0.0%)
2	B	0.19	0/2715	0.50	0/4222
2	D	0.22	0/2671	0.50	0/4154
3	E	0.24	0/906	0.54	0/1224
3	F	0.27	0/888	0.53	0/1205
All	All	0.22	2/23046 (0.0%)	0.53	8/32348 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	13	LEU	C-N	-6.95	1.28	1.33
1	C	14	GLY	CA-C	-6.60	1.47	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	ASP	N-CA-C	9.10	126.48	108.18
1	A	268	ALA	N-CA-C	7.56	119.76	110.91
1	A	195	GLN	CA-C-N	5.55	132.13	121.54
1	A	195	GLN	C-N-CA	5.55	132.13	121.54
1	A	623	PHE	CA-C-N	5.54	132.13	121.54
1	A	623	PHE	C-N-CA	5.54	132.13	121.54
1	C	195	GLN	CA-C-N	5.50	132.04	121.54
1	C	195	GLN	C-N-CA	5.50	132.04	121.54

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	583	TYR	Peptide
1	C	791	PRO	Peptide
1	C	92	LEU	Peptide

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	8035	0	7638	197	0
1	C	7535	0	6823	151	0
2	B	2433	0	1235	71	0
2	D	2393	0	1214	75	0
3	E	890	0	857	35	0
3	F	872	0	813	50	0
All	All	22158	0	18580	533	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 13.

All (533) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:7:ILE:HG12	3:F:56:PHE:HE2	1.20	1.03
1:A:154:GLU:O	1:A:156:GLY:N	1.96	0.99
1:A:154:GLU:O	1:A:157:ALA:N	1.95	0.99
3:E:30:ASN:HD21	3:E:41:ARG:HG3	1.31	0.94
3:F:33:VAL:HG12	3:F:40:LEU:HD12	1.55	0.89
3:E:99:VAL:HA	3:E:102:ARG:HG3	1.55	0.88
1:A:154:GLU:O	1:A:155:LEU:C	2.13	0.86
1:A:916:ARG:H	1:A:916:ARG:HD3	1.40	0.85
3:F:7:ILE:HG12	3:F:56:PHE:CE2	2.11	0.80
1:C:687:LEU:O	1:C:694:ARG:NH2	2.15	0.79
3:F:30:ASN:HD21	3:F:42:ALA:H	1.28	0.78
3:F:33:VAL:HG12	3:F:40:LEU:CD1	2.14	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:563:HIS:O	3:F:102:ARG:NH1	2.17	0.76
1:A:321:ARG:NH2	1:A:343:GLU:OE2	2.18	0.76
3:F:17:VAL:HG11	3:F:22:LYS:HE2	1.66	0.76
3:F:99:VAL:HA	3:F:102:ARG:HG3	1.69	0.74
1:C:90:GLY:O	1:C:93:GLN:HA	1.88	0.73
1:A:1010:ILE:HB	1:A:1057:SER:HB3	1.70	0.73
1:C:559:TYR:CE2	1:C:575:LEU:HD22	2.25	0.72
2:D:73:C:H2'	2:D:74:U:H5'	1.71	0.72
1:A:73:ARG:NH2	2:B:87:C:OP2	2.22	0.72
1:A:575:LEU:HD22	3:F:2:PHE:CE2	2.25	0.72
2:B:118:U:H2'	2:B:119:A:C8	2.24	0.72
1:C:1010:ILE:HB	1:C:1057:SER:HB3	1.72	0.71
1:A:707:ARG:HG2	1:A:729:VAL:HG21	1.70	0.71
2:B:73:C:H2'	2:B:74:U:H5'	1.73	0.71
1:C:28:GLU:HB2	1:C:39:ILE:HD11	1.71	0.71
1:A:624:ASN:O	1:A:626:LYS:HG2	1.91	0.71
1:A:266:LYS:HD3	1:A:422:GLN:HA	1.71	0.70
3:F:58:PHE:HB2	3:F:65:MET:HE3	1.73	0.70
1:A:40:ASP:OD2	1:A:495:TYR:OH	2.11	0.69
2:D:16:U:O2'	2:D:17:A:O5'	2.09	0.69
1:A:154:GLU:C	1:A:156:GLY:N	2.49	0.69
1:A:669:ASP:OD2	1:A:671:ARG:NH1	2.27	0.69
2:D:90:A:H2'	2:D:91:A:C8	2.27	0.68
1:A:205:ARG:NH1	1:A:242:GLN:OE1	2.23	0.68
1:C:947:MET:HE1	1:C:1023:CYS:HB3	1.74	0.68
3:F:7:ILE:CG1	3:F:56:PHE:HE2	2.03	0.68
1:C:122:LEU:O	1:C:123:THR:OG1	2.11	0.68
1:A:488:ILE:HG21	1:A:685:MET:HE1	1.75	0.67
2:B:48:A:H2'	2:B:49:G:H8	1.60	0.67
1:C:30:ASP:HB2	1:C:34:ASN:H	1.57	0.67
1:C:602:ASN:HB2	1:C:651:ARG:HH12	1.59	0.67
1:C:618:THR:HG22	1:C:619:PRO:HD2	1.75	0.67
1:A:321:ARG:HG3	1:A:326:LEU:HB3	1.76	0.66
1:A:103:ILE:HD13	1:A:106:LEU:HD13	1.76	0.66
2:B:90:A:H2'	2:B:91:A:C8	2.30	0.66
1:C:650:GLN:O	1:C:654:LEU:HD13	1.96	0.66
2:D:118:U:H2'	2:D:119:A:C8	2.30	0.66
1:A:428:LEU:O	1:A:432:VAL:HG22	1.96	0.66
2:B:70:C:H2'	2:B:71:G:H8	1.61	0.65
1:A:567:LEU:O	1:A:638:ARG:NH2	2.29	0.65
2:D:119:A:H2'	2:D:120:A:C8	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:ARG:NH2	2:D:87:C:OP2	2.30	0.65
1:A:70:ARG:NH1	2:B:87:C:OP1	2.29	0.65
2:B:119:A:H2'	2:B:120:A:C8	2.32	0.65
1:A:559:TYR:CE2	1:A:564:GLY:HA2	2.33	0.64
2:B:118:U:H2'	2:B:119:A:H8	1.62	0.64
1:A:288:ILE:HD12	1:A:326:LEU:HD21	1.80	0.63
1:A:575:LEU:HD22	3:F:2:PHE:HE2	1.62	0.63
1:A:938:HIS:O	1:A:938:HIS:ND1	2.32	0.62
2:D:70:C:H2'	2:D:71:G:H8	1.64	0.62
1:A:92:LEU:HD11	1:A:103:ILE:HD12	1.81	0.61
1:A:1039:HIS:ND1	1:A:1044:ASN:HA	2.14	0.61
1:C:138:ARG:HH22	1:C:242:GLN:CG	2.13	0.61
1:A:110:PRO:HG2	2:B:60:A:H4'	1.82	0.61
2:B:48:A:H2'	2:B:49:G:C8	2.35	0.61
2:B:60:A:H2'	2:B:61:C:C6	2.35	0.61
1:C:318:ALA:N	1:C:343:GLU:O	2.34	0.61
3:F:104:PHE:HD1	3:F:105:LEU:HD12	1.66	0.61
1:A:277:PHE:CD1	1:A:435:MET:HE1	2.35	0.60
1:A:623:PHE:O	1:A:632:TRP:HB2	2.00	0.60
2:B:90:A:H2'	2:B:91:A:H8	1.66	0.60
1:C:366:LYS:HE3	1:C:367:ASP:O	2.01	0.60
3:F:8:PHE:CE2	3:F:10:SER:HB2	2.37	0.60
1:A:687:LEU:HD21	1:A:694:ARG:HH21	1.65	0.60
1:C:254:MET:HG2	2:D:18:A:N3	2.16	0.60
1:A:772:PHE:HD1	1:A:773:PRO:HD2	1.66	0.60
2:B:73:C:C2'	2:B:74:U:H5'	2.30	0.60
1:C:707:ARG:HG3	1:C:712:LEU:HB2	1.82	0.60
1:A:114:ARG:NH2	1:A:211:GLU:OE1	2.35	0.60
2:B:54:G:H2'	2:B:55:U:O4'	2.02	0.60
1:C:569:SER:HB2	1:C:610:GLU:HG2	1.84	0.60
2:D:90:A:H2'	2:D:91:A:H8	1.67	0.60
1:A:28:GLU:HB2	1:A:39:ILE:HD11	1.84	0.59
1:A:506:ALA:O	1:A:671:ARG:NH2	2.35	0.59
1:C:707:ARG:HD2	1:C:725:LEU:HD22	1.85	0.59
1:A:693:LYS:O	1:A:695:VAL:N	2.36	0.59
1:A:556:LEU:HG	3:F:36:ILE:HD13	1.84	0.59
1:A:852:ARG:NH2	1:A:926:GLU:OE2	2.34	0.59
1:A:288:ILE:HD11	1:A:330:ALA:HB1	1.83	0.59
1:A:544:PHE:HE1	1:A:579:ASN:ND2	2.01	0.59
2:B:30:A:H2'	2:B:31:G:C8	2.39	0.58
1:A:889:LYS:O	1:A:893:GLU:HG2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:110:PRO:HG2	2:D:60:A:H4'	1.85	0.58
1:A:92:LEU:HD13	1:A:97:PHE:HE1	1.68	0.58
1:C:620:TYR:CG	1:C:655:GLN:HB3	2.38	0.58
1:A:177:PRO:HD2	1:A:211:GLU:HB2	1.86	0.58
1:A:316:THR:HB	1:A:319:GLN:HB2	1.85	0.58
1:C:938:HIS:O	1:C:938:HIS:ND1	2.36	0.57
1:A:867:LEU:HD11	1:A:872:LEU:HB2	1.86	0.57
1:A:13:LEU:HD13	1:A:491:VAL:HG11	1.86	0.57
2:D:75:G:H2'	2:D:77:A:N7	2.19	0.57
2:D:75:G:N2	2:D:77:A:H3'	2.19	0.57
2:D:70:C:H2'	2:D:71:G:C8	2.40	0.57
1:A:892:LEU:HD23	1:A:897:ASP:HA	1.86	0.57
2:D:98:G:H2'	2:D:99:C:C6	2.39	0.57
2:D:119:A:H2'	2:D:120:A:H8	1.70	0.57
1:A:327:GLU:O	1:A:327:GLU:HG2	2.04	0.56
1:A:767:HIS:O	1:A:770:THR:OG1	2.23	0.56
3:F:19:ARG:HH11	3:F:19:ARG:HB2	1.70	0.56
3:F:30:ASN:HD21	3:F:41:ARG:HD3	1.70	0.56
1:C:30:ASP:N	1:C:34:ASN:O	2.34	0.56
1:C:86:LEU:HD11	1:C:131:LEU:HD11	1.87	0.56
2:D:118:U:H2'	2:D:119:A:H8	1.69	0.56
2:B:47:G:H2'	2:B:48:A:C8	2.40	0.56
1:C:141:LEU:HB2	1:C:205:ARG:NH1	2.20	0.56
2:B:116:U:H2'	2:B:117:U:C6	2.40	0.56
2:B:119:A:H2'	2:B:120:A:H8	1.69	0.56
2:D:75:G:H5''	2:D:75:G:H8	1.70	0.56
1:A:766:LEU:HD11	1:A:769:LYS:H	1.71	0.56
2:B:116:U:H2'	2:B:117:U:H6	1.71	0.55
1:A:305:LEU:HD22	1:A:315:LEU:HD21	1.87	0.55
2:D:73:C:C2'	2:D:74:U:H5'	2.36	0.55
1:C:877:ASN:HB3	1:C:881:GLU:HG3	1.89	0.55
2:D:60:A:H2'	2:D:61:C:C6	2.41	0.55
2:B:60:A:H2'	2:B:61:C:H6	1.71	0.55
1:C:133:HIS:CD2	2:D:60:A:H5''	2.41	0.55
1:C:577:ARG:O	1:C:579:ASN:N	2.34	0.55
1:C:785:ILE:HD11	1:C:791:PRO:HA	1.89	0.55
2:B:76:A:H2'	2:B:77:A:C8	2.41	0.55
1:C:138:ARG:O	1:C:205:ARG:NE	2.40	0.55
3:E:4:ARG:HG3	3:E:91:GLU:HG2	1.89	0.55
3:F:94:TYR:HA	3:F:97:LEU:HD12	1.89	0.55
1:A:255:LEU:HD13	1:A:266:LYS:HG2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ARG:HH21	1:A:134:LEU:HD21	1.71	0.54
1:A:594:ARG:NH2	1:A:659:GLU:OE2	2.39	0.54
2:B:117:U:H2'	2:B:118:U:C6	2.42	0.54
3:F:9:THR:O	3:F:51:GLY:N	2.40	0.54
2:B:104:U:H2'	2:B:105:A:C8	2.43	0.54
2:B:117:U:H2'	2:B:118:U:H6	1.71	0.54
1:C:139:GLY:O	1:C:205:ARG:NH2	2.41	0.54
2:B:17:A:H3'	2:B:18:A:H5''	1.90	0.54
1:C:123:THR:HG22	1:C:124:PRO:HD2	1.89	0.54
2:D:24:C:C2	2:D:25:G:C8	2.96	0.54
1:C:41:LEU:HD13	1:C:828:LEU:HD12	1.90	0.54
3:E:25:LEU:HD23	3:E:49:LEU:HD13	1.90	0.54
1:A:196:ARG:NH1	1:A:863:PRO:HG3	2.22	0.54
1:A:268:ALA:HB1	1:A:271:THR:OG1	2.08	0.54
1:A:123:THR:HG22	1:A:124:PRO:HD2	1.90	0.54
1:A:429:ARG:O	1:A:433:PRO:HD2	2.08	0.54
1:A:624:ASN:O	1:A:626:LYS:N	2.41	0.54
2:D:27:U:H2'	2:D:28:G:C8	2.43	0.54
2:D:48:A:H2'	2:D:49:G:C8	2.42	0.53
1:A:47:GLU:O	1:A:479:ARG:NH2	2.33	0.53
1:A:616:ASN:HB3	1:A:750:MET:HE2	1.90	0.53
1:A:557:ARG:HA	3:F:36:ILE:HG23	1.88	0.53
1:C:308:GLU:OE1	1:C:312:LYS:NZ	2.33	0.53
1:A:565:LYS:HE3	1:A:638:ARG:HD3	1.89	0.53
2:B:98:G:H2'	2:B:99:C:C6	2.44	0.53
1:A:277:PHE:HD1	1:A:435:MET:HE1	1.74	0.53
1:C:141:LEU:HD23	1:C:143:GLN:H	1.73	0.53
1:C:1039:HIS:ND1	1:C:1044:ASN:HA	2.23	0.53
2:D:116:U:H2'	2:D:117:U:C6	2.44	0.53
2:B:128:C:H2'	2:B:129:G:H8	1.74	0.53
1:C:396:THR:O	1:C:400:LYS:CB	2.57	0.53
2:D:47:G:H2'	2:D:48:A:C8	2.43	0.53
3:F:25:LEU:HB2	3:F:75:PHE:CE2	2.44	0.53
2:D:27:U:H2'	2:D:28:G:H8	1.74	0.53
2:D:104:U:H2'	2:D:105:A:H8	1.74	0.53
3:F:38:GLU:OE2	3:F:41:ARG:NE	2.37	0.53
1:A:106:LEU:HD12	1:A:107:PRO:HD2	1.91	0.52
1:A:464:LEU:HD13	1:A:685:MET:HE3	1.91	0.52
2:D:104:U:H2'	2:D:105:A:C8	2.43	0.52
3:E:58:PHE:HB2	3:E:65:MET:HE3	1.89	0.52
1:C:50:GLU:OE2	1:C:516:ARG:NH2	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:430:ARG:O	1:C:433:PRO:HD2	2.09	0.52
1:C:707:ARG:NH1	1:C:726:ASP:OD2	2.42	0.52
1:C:647:SER:O	1:C:651:ARG:HG3	2.09	0.52
1:A:91:VAL:O	1:A:124:PRO:HB3	2.10	0.52
2:D:26:U:H2'	2:D:27:U:C6	2.44	0.52
1:A:635:PHE:O	1:A:639:VAL:HG13	2.08	0.52
2:D:98:G:H2'	2:D:99:C:H6	1.75	0.52
3:E:104:PHE:HD1	3:E:105:LEU:HD12	1.74	0.52
1:A:559:TYR:HE1	1:A:575:LEU:HD12	1.74	0.52
1:A:114:ARG:HG2	1:A:130:VAL:HG13	1.91	0.52
2:B:27:U:H2'	2:B:28:G:H8	1.74	0.52
3:E:4:ARG:NH1	3:E:91:GLU:OE1	2.36	0.52
2:D:37:U:H2'	2:D:38:U:C6	2.45	0.51
3:E:39:TYR:O	3:E:40:LEU:HG	2.10	0.51
2:B:26:U:H2'	2:B:27:U:C6	2.44	0.51
1:C:712:LEU:HD22	1:C:725:LEU:HD21	1.92	0.51
1:C:867:LEU:HD11	1:C:872:LEU:HB2	1.92	0.51
2:D:48:A:H2'	2:D:49:G:H8	1.75	0.51
1:A:422:GLN:O	1:A:423:ILE:HG13	2.10	0.51
1:A:432:VAL:HG23	1:A:433:PRO:CD	2.40	0.51
1:C:541:PHE:HZ	1:C:578:LEU:HD12	1.75	0.51
1:C:693:LYS:O	1:C:695:VAL:N	2.43	0.51
1:A:289:LEU:HD12	1:A:289:LEU:O	2.11	0.51
2:B:24:C:C2	2:B:25:G:C8	2.99	0.51
1:C:1065:GLU:CD	1:C:1065:GLU:H	2.19	0.51
2:D:60:A:H2'	2:D:61:C:H6	1.76	0.51
2:B:61:C:H2'	2:B:62:A:C8	2.46	0.51
2:B:104:U:H2'	2:B:105:A:H8	1.76	0.51
1:C:414:HIS:O	1:C:415:ILE:HG13	2.10	0.51
1:C:954:GLU:OE2	1:C:998:LYS:NZ	2.37	0.51
1:A:193:ARG:NH1	2:B:24:C:OP2	2.44	0.51
1:A:328:ASP:O	1:A:329:THR:HG22	2.12	0.51
2:B:25:G:H2'	2:B:26:U:H6	1.76	0.51
2:D:76:A:H2'	2:D:77:A:C8	2.46	0.51
3:F:11:PHE:CD1	3:F:50:ASP:HB2	2.46	0.51
1:A:70:ARG:NH2	2:B:19:C:OP1	2.44	0.50
1:C:194:ASN:OD1	1:C:201:HIS:HB2	2.10	0.50
1:C:498:PRO:HG2	1:C:687:LEU:HD23	1.93	0.50
3:E:95:VAL:HA	3:E:102:ARG:HH22	1.75	0.50
1:A:567:LEU:HD11	1:A:639:VAL:HG12	1.93	0.50
1:C:593:SER:OG	1:C:749:GLU:OE2	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:HIS:CD2	2:B:85:G:C8	2.99	0.50
2:B:27:U:H2'	2:B:28:G:C8	2.47	0.50
2:D:51:A:H2'	2:D:52:C:C6	2.46	0.50
1:A:575:LEU:HD13	3:F:2:PHE:CD2	2.47	0.50
1:A:575:LEU:HD23	1:A:576:GLY:H	1.75	0.50
2:D:39:U:H2'	2:D:40:C:H5''	1.94	0.50
2:D:54:G:H2'	2:D:55:U:O4'	2.12	0.50
2:D:117:U:H2'	2:D:118:U:C6	2.47	0.50
3:E:95:VAL:HA	3:E:102:ARG:NH2	2.26	0.50
1:A:35:PRO:HG2	1:A:775:PRO:HB3	1.93	0.50
1:C:435:MET:SD	1:C:441:TYR:HA	2.52	0.50
1:A:317:TYR:HD2	1:A:345:SER:HB2	1.77	0.50
2:D:47:G:O2'	2:D:48:A:P	2.70	0.50
3:E:7:ILE:HD12	3:E:56:PHE:CE2	2.47	0.50
3:E:25:LEU:O	3:E:29:ILE:HG13	2.12	0.50
1:A:252:GLN:O	1:A:255:LEU:HD12	2.11	0.49
1:A:463:TYR:OH	1:A:493:ARG:HG2	2.12	0.49
1:A:693:LYS:C	1:A:695:VAL:H	2.19	0.49
2:D:75:G:H22	2:D:77:A:H3'	1.76	0.49
2:D:128:C:H2'	2:D:129:G:H8	1.77	0.49
1:A:30:ASP:HB3	1:A:32:GLU:H	1.76	0.49
2:D:52:C:H2'	2:D:53:C:O4'	2.12	0.49
2:B:45:A:H2'	2:B:46:A:C8	2.47	0.49
1:C:678:CYS:HA	1:C:695:VAL:HG11	1.94	0.49
1:A:258:CYS:SG	1:A:424:SER:N	2.83	0.49
3:E:0:SER:O	3:E:0:SER:OG	2.27	0.49
1:C:243:ARG:NH1	2:D:85:G:O5'	2.44	0.49
3:F:30:ASN:ND2	3:F:41:ARG:HD3	2.27	0.49
2:B:47:G:O2'	2:B:48:A:P	2.70	0.49
1:C:17:ILE:HB	1:C:505:THR:HG22	1.95	0.49
1:C:386:PHE:CZ	1:C:412:LEU:HD12	2.46	0.49
1:C:961:LEU:O	1:C:979:VAL:HG23	2.12	0.49
3:E:5:ALA:HB2	3:E:65:MET:HE2	1.95	0.49
3:E:29:ILE:HG23	3:E:56:PHE:HZ	1.78	0.49
2:D:95:U:H2'	2:D:96:C:H6	1.78	0.49
1:C:584:VAL:HG23	1:C:605:LEU:HD11	1.95	0.49
1:C:138:ARG:HB2	1:C:205:ARG:HE	1.78	0.49
1:C:324:LEU:HA	3:F:104:PHE:CD2	2.47	0.49
1:A:28:GLU:HB3	1:A:37:ARG:HB3	1.95	0.48
1:A:41:LEU:HD13	1:A:828:LEU:HD12	1.96	0.48
1:A:565:LYS:HG2	1:A:570:GLY:HA2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:24:ARG:O	3:E:28:ILE:HG13	2.14	0.48
2:B:72:U:H2'	2:B:73:C:H6	1.78	0.48
2:D:25:G:H2'	2:D:26:U:H6	1.78	0.48
2:D:75:G:H5'	2:D:76:A:OP2	2.13	0.48
3:F:97:LEU:HD22	3:F:105:LEU:HB3	1.95	0.48
1:A:194:ASN:OD1	1:A:201:HIS:HB2	2.14	0.48
1:A:563:HIS:HB3	3:F:102:ARG:HH11	1.79	0.48
1:A:939:ASN:O	2:B:27:U:O2'	2.31	0.48
1:C:86:LEU:HD21	1:C:131:LEU:HD11	1.94	0.48
2:D:75:G:N2	2:D:78:A:C8	2.81	0.48
1:A:185:PHE:CE2	1:A:201:HIS:HB3	2.49	0.48
1:A:785:ILE:HD11	1:A:791:PRO:HA	1.95	0.48
1:A:1010:ILE:HG12	1:A:1015:ARG:HD2	1.96	0.48
1:C:346:THR:HG22	1:C:346:THR:O	2.14	0.48
1:A:492:VAL:HG22	1:A:498:PRO:HD3	1.94	0.48
1:C:65:ALA:O	1:C:69:ARG:HG3	2.14	0.48
1:C:369:LYS:HD3	3:F:80:THR:HA	1.95	0.48
2:D:102:C:H2'	2:D:103:U:C6	2.48	0.48
1:A:108:ASN:C	1:A:110:PRO:HD3	2.39	0.48
2:B:98:G:H2'	2:B:99:C:H6	1.79	0.48
3:F:24:ARG:O	3:F:28:ILE:HG13	2.13	0.48
1:A:337:TYR:OH	1:A:343:GLU:OE2	2.26	0.47
1:A:567:LEU:HD22	1:A:604:VAL:HG11	1.96	0.47
2:B:26:U:H2'	2:B:27:U:H6	1.79	0.47
1:C:112:GLN:HA	1:C:182:LEU:HD11	1.95	0.47
3:E:7:ILE:HD13	3:E:72:LEU:CD1	2.44	0.47
3:E:99:VAL:HA	3:E:102:ARG:CG	2.34	0.47
2:B:70:C:H2'	2:B:71:G:C8	2.44	0.47
2:B:83:G:O2'	2:B:84:U:H5'	2.14	0.47
1:C:174:PHE:HB3	1:C:180:LEU:HB2	1.95	0.47
1:C:908:TYR:HD1	1:C:917:THR:O	1.97	0.47
3:F:53:TYR:HB3	3:F:103:ASP:HB3	1.96	0.47
1:C:138:ARG:HE	1:C:205:ARG:NH2	2.12	0.47
1:C:718:GLU:CB	1:C:1039:HIS:H	2.27	0.47
1:C:138:ARG:HH22	1:C:242:GLN:HG2	1.78	0.47
1:C:900:ALA:O	1:C:904:ALA:HB2	2.14	0.47
1:A:310:TYR:CZ	1:A:350:MET:HE2	2.49	0.47
2:B:95:U:H2'	2:B:96:C:H6	1.80	0.47
2:D:61:C:H2'	2:D:62:A:C8	2.49	0.47
3:E:11:PHE:CE2	3:E:17:VAL:HG21	2.49	0.47
1:C:29:ILE:HD12	1:C:33:GLU:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1063:ILE:HD11	1:A:1067:GLY:HA2	1.97	0.47
1:C:844:MET:HE3	1:C:945:ALA:HA	1.97	0.47
2:D:83:G:O2'	2:D:84:U:H5'	2.15	0.47
1:A:139:GLY:O	2:B:22:U:H4'	2.15	0.47
1:A:448:ILE:HG22	1:A:449:TYR:CD2	2.50	0.47
1:A:563:HIS:HB3	3:F:102:ARG:NH1	2.30	0.47
1:C:174:PHE:CD2	1:C:180:LEU:HD12	2.50	0.47
1:A:26:MET:HE1	1:A:495:TYR:CG	2.50	0.46
1:A:563:HIS:NE2	1:A:641:THR:O	2.44	0.46
1:A:864:LEU:HA	1:A:867:LEU:HB2	1.97	0.46
1:A:317:TYR:CD1	1:A:347:LEU:HD22	2.49	0.46
1:A:324:LEU:HA	3:E:104:PHE:CD2	2.50	0.46
1:A:709:PHE:HB3	1:A:772:PHE:CD2	2.51	0.46
2:B:56:U:H2'	2:B:57:G:C8	2.50	0.46
1:C:323:LEU:O	3:F:104:PHE:HD2	1.98	0.46
1:A:875:MET:HG2	1:A:877:ASN:H	1.79	0.46
1:C:965:TYR:O	1:C:969:VAL:HG23	2.16	0.46
2:D:91:A:H2'	2:D:92:C:C6	2.50	0.46
1:C:131:LEU:HD12	1:C:132:LEU:N	2.31	0.46
1:C:867:LEU:CD1	1:C:872:LEU:HB2	2.46	0.46
3:E:16:LYS:HD2	3:E:16:LYS:HA	1.68	0.46
1:A:77:HIS:CD2	2:B:85:G:H8	2.33	0.46
1:A:953:PHE:CE2	1:A:997:PHE:HB2	2.49	0.46
1:A:1032:ILE:HD11	1:A:1047:LEU:HD12	1.96	0.46
2:B:56:U:H2'	2:B:57:G:H8	1.81	0.46
1:C:11:TYR:CE2	1:C:26:MET:HG2	2.51	0.46
1:C:953:PHE:HD2	1:C:997:PHE:HA	1.79	0.46
2:D:37:U:H2'	2:D:38:U:H6	1.81	0.46
1:A:772:PHE:HD1	1:A:773:PRO:CD	2.27	0.46
1:A:978:ALA:O	1:A:988:TRP:HB3	2.16	0.46
1:C:212:LEU:HD21	1:C:235:ILE:HG22	1.98	0.46
1:A:100:ASN:O	1:A:102:LEU:N	2.48	0.46
1:A:182:LEU:O	1:A:186:GLU:HB2	2.15	0.46
1:A:317:TYR:CE1	1:A:347:LEU:HD13	2.51	0.46
2:B:61:C:H2'	2:B:62:A:H8	1.80	0.46
1:C:50:GLU:CD	1:C:516:ARG:HH22	2.24	0.46
3:E:98:PRO:O	3:E:102:ARG:HG2	2.16	0.46
1:A:63:ARG:HH21	2:B:17:A:H5'	1.80	0.46
1:A:288:ILE:HG13	1:A:331:PHE:O	2.15	0.46
2:B:39:U:H2'	2:B:40:C:H5''	1.98	0.46
1:C:361:GLU:HA	1:C:366:LYS:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:699:ASN:O	1:C:702:ILE:HG13	2.16	0.46
1:C:181:ALA:O	1:C:185:PHE:HB2	2.16	0.45
1:C:875:MET:HE2	1:C:923:VAL:HG11	1.98	0.45
2:D:45:A:H2'	2:D:46:A:C8	2.50	0.45
1:A:687:LEU:HG	1:A:694:ARG:HE	1.80	0.45
1:C:360:LEU:HB2	1:C:366:LYS:HB2	1.98	0.45
1:C:924:ARG:HG3	2:D:56:U:OP1	2.16	0.45
2:B:25:G:H2'	2:B:26:U:C6	2.51	0.45
1:C:575:LEU:HG	3:E:2:PHE:CE2	2.51	0.45
1:C:745:VAL:O	1:C:749:GLU:HG2	2.16	0.45
2:D:25:G:H2'	2:D:26:U:C6	2.52	0.45
1:A:109:THR:N	1:A:110:PRO:HD3	2.32	0.45
1:A:268:ALA:H	1:A:420:PHE:HD2	1.62	0.45
1:A:432:VAL:HG23	1:A:433:PRO:HD3	1.97	0.45
1:A:442:ASP:OD1	1:A:443:GLU:N	2.50	0.45
1:A:472:ILE:HD11	1:A:676:PHE:CD2	2.50	0.45
1:C:583:TYR:CE1	1:C:607:LEU:HD21	2.52	0.45
1:A:194:ASN:HB3	1:A:198:ASP:O	2.16	0.45
3:F:92:SER:O	3:F:96:ARG:HG3	2.16	0.45
1:C:246:LEU:HG	2:D:19:C:O2'	2.17	0.45
1:A:317:TYR:CD2	1:A:345:SER:HB2	2.52	0.45
1:C:45:VAL:CG2	1:C:832:ARG:HG2	2.47	0.45
1:C:109:THR:O	1:C:113:LEU:HG	2.17	0.45
1:C:865:THR:O	1:C:899:PRO:HB2	2.16	0.45
3:E:2:PHE:HB3	3:E:57:LEU:HD11	1.99	0.45
3:E:40:LEU:HD12	3:E:40:LEU:O	2.17	0.45
1:A:267:ALA:HA	1:A:420:PHE:HD2	1.81	0.45
1:A:377:GLU:N	1:A:377:GLU:OE1	2.50	0.45
2:B:40:C:H5''	2:B:40:C:H6	1.81	0.45
2:B:75:G:N2	2:B:77:A:H3'	2.31	0.45
2:B:102:C:H2'	2:B:103:U:C6	2.52	0.45
1:C:328:ASP:O	1:C:329:THR:HG22	2.17	0.45
1:C:366:LYS:NZ	1:C:372:LEU:HA	2.32	0.45
2:D:17:A:H3'	2:D:18:A:H5''	1.99	0.45
3:F:99:VAL:HA	3:F:102:ARG:CG	2.44	0.45
2:B:30:A:H2'	2:B:31:G:H8	1.82	0.44
1:C:610:GLU:N	1:C:610:GLU:OE1	2.50	0.44
3:E:1:MET:HE3	3:E:1:MET:HB2	1.76	0.44
2:B:36:C:O2'	2:B:37:U:H5'	2.18	0.44
1:C:141:LEU:HB2	1:C:205:ARG:HH12	1.82	0.44
2:D:75:G:H5''	2:D:75:G:C8	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:584:VAL:HB	1:A:605:LEU:HD11	1.99	0.44
1:A:961:LEU:O	1:A:979:VAL:HG22	2.17	0.44
1:C:109:THR:N	1:C:110:PRO:HD3	2.32	0.44
2:D:40:C:H5''	2:D:40:C:H6	1.82	0.44
1:C:305:LEU:HD22	1:C:315:LEU:HD21	1.98	0.44
1:C:306:MET:HE1	1:C:436:GLU:O	2.17	0.44
1:C:882:PRO:O	1:C:886:GLU:HG2	2.17	0.44
1:A:559:TYR:HD2	1:A:573:ILE:HD11	1.82	0.44
1:C:575:LEU:HD21	3:E:57:LEU:HD21	1.99	0.44
1:C:939:ASN:O	2:D:27:U:O2'	2.34	0.44
1:C:320:ALA:O	1:C:324:LEU:HG	2.17	0.44
2:B:31:G:H2'	2:B:32:C:C6	2.53	0.44
2:B:124:G:H2'	2:B:125:C:O4'	2.18	0.44
1:C:564:GLY:O	1:C:573:ILE:HG12	2.17	0.44
1:A:318:ALA:HB2	1:A:343:GLU:HB3	2.00	0.44
1:A:433:PRO:HA	1:A:436:GLU:OE1	2.17	0.44
1:C:563:HIS:O	3:E:102:ARG:NH1	2.48	0.44
1:C:310:TYR:CZ	1:C:350:MET:HE2	2.53	0.43
1:C:435:MET:HE3	1:C:435:MET:HB3	1.72	0.43
1:C:967:TRP:HA	2:D:66:A:N3	2.33	0.43
2:B:75:G:H5''	2:B:75:G:H8	1.83	0.43
1:C:12:ILE:HG13	1:C:29:ILE:HG21	2.00	0.43
1:C:584:VAL:CG2	1:C:605:LEU:HD11	2.48	0.43
3:E:11:PHE:CD1	3:E:50:ASP:HB2	2.53	0.43
3:F:8:PHE:HE2	3:F:10:SER:HB2	1.83	0.43
3:F:30:ASN:O	3:F:33:VAL:HG22	2.18	0.43
1:A:396:THR:O	1:A:400:LYS:CB	2.65	0.43
1:A:504:GLU:HG2	1:A:726:ASP:OD2	2.18	0.43
1:A:741:ILE:HG23	1:A:772:PHE:CE1	2.54	0.43
1:A:880:ARG:C	1:A:882:PRO:HD3	2.43	0.43
1:C:311:ARG:HH22	3:F:85:ASP:CG	2.27	0.43
1:C:559:TYR:OH	3:E:91:GLU:OE2	2.36	0.43
1:A:844:MET:HA	1:A:844:MET:HE2	2.00	0.43
1:C:298:THR:HG22	1:C:301:GLU:OE1	2.18	0.43
1:C:500:ARG:NH1	1:C:694:ARG:HA	2.33	0.43
1:C:878:ARG:HG3	1:C:885:TYR:CB	2.48	0.43
1:A:541:PHE:HZ	1:A:578:LEU:HD12	1.83	0.43
1:C:369:LYS:HD3	3:F:81:GLY:H	1.83	0.43
1:A:57:SER:HB3	2:B:17:A:OP2	2.19	0.43
1:A:71:LEU:HD23	1:A:71:LEU:HA	1.87	0.43
3:F:95:VAL:HA	3:F:102:ARG:NH2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:MET:HE1	1:A:495:TYR:CD1	2.54	0.43
1:A:875:MET:HG2	1:A:876:VAL:N	2.33	0.43
1:C:517:LYS:HD2	1:C:517:LYS:HA	1.93	0.43
1:C:725:LEU:HD11	1:C:784:MET:HE2	2.01	0.43
1:C:979:VAL:HG13	1:C:988:TRP:CZ2	2.53	0.43
1:A:16:ASP:HA	1:A:504:GLU:O	2.19	0.43
1:A:583:TYR:CD2	1:A:584:VAL:HG13	2.54	0.43
1:C:775:PRO:HG2	1:C:779:PHE:CG	2.54	0.43
1:A:449:TYR:O	1:A:449:TYR:CD1	2.72	0.43
1:A:573:ILE:HG22	1:A:607:LEU:HD21	2.01	0.43
1:A:559:TYR:OH	3:F:91:GLU:OE2	2.19	0.42
1:A:125:LEU:HD23	1:A:125:LEU:HA	1.90	0.42
1:A:670:THR:O	1:A:670:THR:OG1	2.33	0.42
1:A:785:ILE:HG21	1:A:810:LEU:HG	2.00	0.42
1:C:69:ARG:HD2	2:D:88:G:C6	2.55	0.42
1:C:267:ALA:N	1:C:423:ILE:O	2.46	0.42
1:C:360:LEU:CB	1:C:366:LYS:HB2	2.49	0.42
1:C:678:CYS:SG	1:C:697:ALA:HB2	2.59	0.42
2:D:30:A:H2'	2:D:31:G:C8	2.54	0.42
1:A:58:LEU:HD12	1:A:58:LEU:H	1.85	0.42
1:A:86:LEU:HD22	1:A:91:VAL:HG11	2.01	0.42
1:A:87:LYS:HB2	1:A:97:PHE:CE2	2.54	0.42
1:A:110:PRO:CG	2:B:60:A:H4'	2.47	0.42
1:A:562:GLN:HE22	1:A:605:LEU:HB3	1.84	0.42
1:C:139:GLY:O	2:D:22:U:H4'	2.19	0.42
1:C:566:CYS:SG	1:C:569:SER:HB3	2.59	0.42
1:A:158:LEU:HD12	1:A:199:TYR:CE1	2.55	0.42
1:A:348:MET:HE2	1:A:350:MET:SD	2.59	0.42
1:C:111:TRP:CE2	1:C:133:HIS:HE1	2.38	0.42
1:C:607:LEU:HD23	1:C:607:LEU:HA	1.78	0.42
2:D:105:A:H2'	2:D:106:A:C8	2.55	0.42
3:E:11:PHE:HB2	3:E:50:ASP:OD1	2.19	0.42
1:A:174:PHE:HE1	1:A:183:ASN:HD22	1.68	0.42
1:A:745:VAL:O	1:A:749:GLU:HG2	2.20	0.42
1:C:462:ILE:HG23	1:C:463:TYR:CD2	2.53	0.42
1:C:618:THR:HG22	1:C:619:PRO:CD	2.46	0.42
2:B:47:G:O2'	2:B:48:A:O5'	2.37	0.42
3:F:18:SER:O	3:F:21:GLU:N	2.52	0.42
1:A:114:ARG:HE	1:A:134:LEU:HD21	1.84	0.42
1:A:154:GLU:O	1:A:156:GLY:C	2.62	0.42
1:A:235:ILE:HD13	1:A:235:ILE:HA	1.94	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:LEU:HD22	1:A:266:LYS:HG3	2.02	0.42
1:A:1043:LYS:HG3	1:A:1043:LYS:O	2.19	0.42
1:C:26:MET:HE1	1:C:495:TYR:CG	2.54	0.42
2:D:124:G:H2'	2:D:125:C:O4'	2.20	0.42
3:E:39:TYR:CE1	3:E:40:LEU:HD23	2.55	0.42
3:F:19:ARG:HB2	3:F:19:ARG:NH1	2.34	0.42
3:F:91:GLU:O	3:F:95:VAL:HG22	2.20	0.42
1:A:287:ARG:HD3	1:A:296:PRO:HB3	2.02	0.42
1:C:135:ILE:HD13	1:C:135:ILE:HA	1.88	0.42
3:F:17:VAL:CG1	3:F:22:LYS:HE2	2.44	0.42
1:C:24:TRP:CZ2	1:C:1066:LEU:HG	2.55	0.41
2:D:116:U:H2'	2:D:117:U:H6	1.84	0.41
1:A:781:GLN:HG2	1:A:814:LEU:HD21	2.02	0.41
1:C:108:ASN:C	1:C:110:PRO:HD3	2.45	0.41
3:F:30:ASN:ND2	3:F:42:ALA:H	2.07	0.41
1:A:11:TYR:CD2	1:A:26:MET:HG2	2.55	0.41
1:C:1041:ILE:HG22	1:C:1047:LEU:HD21	2.02	0.41
2:D:91:A:H2'	2:D:92:C:H6	1.84	0.41
3:E:7:ILE:HA	3:E:85:ASP:O	2.20	0.41
1:C:13:LEU:HD13	1:C:491:VAL:HG11	2.01	0.41
2:D:81:A:H2'	2:D:82:U:O4'	2.21	0.41
3:E:62:SER:HA	3:E:63:PRO:HD3	1.92	0.41
3:F:30:ASN:HA	3:F:33:VAL:HG22	2.03	0.41
1:A:746:ARG:O	1:A:750:MET:HE3	2.20	0.41
1:C:559:TYR:CE1	1:C:564:GLY:HA2	2.55	0.41
2:D:126:A:C6	2:D:128:C:C2	3.08	0.41
1:A:92:LEU:CD1	1:A:103:ILE:HD12	2.48	0.41
1:A:256:GLY:O	1:A:266:LYS:HE3	2.20	0.41
1:A:463:TYR:HE1	1:A:492:VAL:HG12	1.86	0.41
1:A:624:ASN:C	1:A:626:LYS:H	2.28	0.41
2:B:91:A:H2'	2:B:92:C:C6	2.55	0.41
3:F:7:ILE:O	3:F:7:ILE:HG13	2.20	0.41
1:A:624:ASN:ND2	1:A:627:ASP:OD2	2.51	0.41
1:A:959:TYR:HB2	1:A:1052:VAL:O	2.21	0.41
1:A:965:TYR:O	1:A:969:VAL:HG23	2.20	0.41
2:B:81:A:H2'	2:B:82:U:O4'	2.21	0.41
1:C:318:ALA:O	1:C:322:LYS:HG3	2.20	0.41
1:C:574:ASN:O	1:C:575:LEU:HB3	2.21	0.41
1:C:931:THR:HG21	1:C:1025:ARG:NH2	2.35	0.41
2:D:102:C:H2'	2:D:103:U:H6	1.84	0.41
1:A:78:ARG:HG3	1:A:238:LEU:HD22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:575:LEU:CD1	3:F:57:LEU:HD21	2.51	0.41
1:C:842:GLY:HA2	1:C:966:SER:OG	2.20	0.41
1:A:65:ALA:O	1:A:69:ARG:HG3	2.21	0.41
1:A:109:THR:O	1:A:113:LEU:HG	2.20	0.41
1:A:326:LEU:HD12	1:A:326:LEU:HA	1.80	0.41
1:A:1008:GLU:OE1	1:A:1061:TYR:OH	2.33	0.41
1:A:1041:ILE:HG22	1:A:1047:LEU:HD21	2.02	0.41
1:C:127:TRP:O	1:C:131:LEU:HG	2.20	0.41
1:A:24:TRP:CZ2	1:A:1066:LEU:HG	2.55	0.41
1:A:324:LEU:HA	3:E:104:PHE:CE2	2.56	0.41
1:A:588:HIS:HA	1:A:602:ASN:O	2.20	0.41
1:A:844:MET:HE3	1:A:945:ALA:HA	2.03	0.41
2:D:47:G:O2'	2:D:48:A:O5'	2.39	0.41
3:F:25:LEU:O	3:F:29:ILE:HG13	2.22	0.41
1:C:1009:VAL:HG22	1:C:1058:PHE:CD1	2.56	0.40
2:B:128:C:H2'	2:B:129:G:C8	2.55	0.40
1:C:525:GLU:HA	1:C:528:LYS:HD2	2.04	0.40
1:A:396:THR:HA	1:A:408:LEU:HD21	2.03	0.40
1:A:997:PHE:CE2	1:A:1078:ARG:HA	2.56	0.40
2:D:105:A:H2'	2:D:106:A:H8	1.87	0.40
1:A:521:LYS:O	1:A:525:GLU:HG2	2.21	0.40
1:A:620:TYR:CD1	1:A:655:GLN:HG2	2.57	0.40
2:B:105:A:H2'	2:B:106:A:C8	2.56	0.40
1:C:243:ARG:NH1	2:D:84:U:H3'	2.36	0.40
1:C:396:THR:HG23	1:C:412:LEU:HD21	2.04	0.40
1:A:493:ARG:NH2	2:B:123:G:O3'	2.55	0.40
1:C:278:ILE:O	1:C:281:THR:OG1	2.34	0.40
1:C:992:ASP:H	1:C:995:PHE:HD2	1.67	0.40
2:D:92:C:C2	2:D:93:G:C8	3.10	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	1048/1092 (96%)	964 (92%)	73 (7%)	11 (1%)	12	44
1	C	1024/1092 (94%)	949 (93%)	69 (7%)	6 (1%)	21	53
3	E	115/117 (98%)	110 (96%)	5 (4%)	0	100	100
3	F	115/117 (98%)	105 (91%)	10 (9%)	0	100	100
All	All	2302/2418 (95%)	2128 (92%)	157 (7%)	17 (1%)	18	51

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	440	ARG
1	A	625	GLY
1	C	578	LEU
1	A	94	ALA
1	A	101	GLY
1	A	153	LYS
1	A	155	LEU
1	A	328	ASP
1	A	694	ARG
1	C	694	ARG
1	A	95	ALA
1	C	347	LEU
1	A	687	LEU
1	C	197	SER
1	A	197	SER
1	C	328	ASP
1	C	584	VAL

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	774/942 (82%)	773 (100%)	1 (0%)	88	87
1	C	670/942 (71%)	670 (100%)	0	100	100
3	E	84/99 (85%)	84 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	F	79/99 (80%)	79 (100%)	0	100	100
All	All	1607/2082 (77%)	1606 (100%)	1 (0%)	88	87

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	191	HIS
1	A	835	ASN
1	C	34	ASN
1	C	108	ASN
1	C	137	HIS
1	C	270	ASN
1	C	319	GLN
1	C	835	ASN
1	C	895	HIS
1	C	939	ASN
1	C	989	GLN
3	E	30	ASN
3	F	30	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	113/135 (83%)	32 (28%)	3 (2%)
2	D	112/135 (82%)	29 (25%)	4 (3%)
All	All	225/270 (83%)	61 (27%)	7 (3%)

All (61) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	17	A
2	B	18	A
2	B	19	C

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Mol	Chain	Res	Type
2	B	30	A
2	B	36	C
2	B	37	U
2	B	40	C
2	B	42	C
2	B	44	A
2	B	45	A
2	B	46	A
2	B	47	G
2	B	48	A
2	B	50	A
2	B	51	A
2	B	65	A
2	B	74	U
2	B	76	A
2	B	77	A
2	B	79	A
2	B	83	G
2	B	84	U
2	B	85	G
2	B	89	C
2	B	97	U
2	B	98	G
2	B	110	U
2	B	115	C
2	B	125	C
2	B	126	A
2	B	127	U
2	B	133	A
2	D	17	A
2	D	18	A
2	D	19	C
2	D	30	A
2	D	37	U
2	D	40	C
2	D	42	C
2	D	44	A
2	D	45	A
2	D	46	A
2	D	47	G
2	D	48	A
2	D	50	A

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Mol	Chain	Res	Type
2	D	51	A
2	D	65	A
2	D	74	U
2	D	76	A
2	D	79	A
2	D	83	G
2	D	84	U
2	D	85	G
2	D	89	C
2	D	97	U
2	D	98	G
2	D	115	C
2	D	125	C
2	D	126	A
2	D	127	U
2	D	133	A

All (7) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	45	A
2	B	47	G
2	B	83	G
2	D	16	U
2	D	45	A
2	D	47	G
2	D	83	G

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	1056/1092 (96%)	0.37	42 (3%)	42	25	6, 44, 90, 119	0
1	C	1036/1092 (94%)	0.65	83 (8%)	18	13	6, 61, 110, 146	0
2	B	115/135 (85%)	0.47	9 (7%)	19	13	12, 52, 193, 225	0
2	D	113/135 (83%)	0.62	9 (7%)	18	13	15, 71, 153, 195	0
3	E	117/117 (100%)	0.21	1 (0%)	81	58	13, 37, 76, 86	0
3	F	117/117 (100%)	0.29	4 (3%)	48	29	7, 37, 76, 80	0
All	All	2554/2688 (95%)	0.49	148 (5%)	29	18	6, 51, 108, 225	0

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	342	ALA	5.6
1	C	340	ASP	5.6
1	C	918	GLN	5.3
1	C	143	GLN	4.9
1	C	445	CYS	4.5
1	A	451	ASP	4.0
1	C	624	ASN	4.0
1	A	611	ASN	3.8
2	D	44	A	3.6
1	C	101	GLY	3.5
1	C	189	SER	3.4
1	C	389	PHE	3.4
1	C	871	ASP	3.4
2	D	114	G	3.4
1	C	338	GLY	3.4
1	C	337	TYR	3.3
1	C	292	GLY	3.2
1	C	438	GLY	3.2
1	A	257	HIS	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	458	THR	3.1
1	C	343	GLU	3.1
1	C	334	GLY	3.0
1	C	196	ARG	3.0
1	A	769	LYS	3.0
1	A	186	GLU	3.0
1	A	1083	SER	3.0
1	C	286	LEU	2.9
1	C	770	THR	2.9
1	C	459	GLU	2.9
1	A	765	VAL	2.9
1	C	130	VAL	2.9
1	C	1083	SER	2.9
2	D	46	A	2.9
1	A	256	GLY	2.9
1	A	449	TYR	2.9
1	C	399	LEU	2.9
3	F	60	ASP	2.8
1	C	149	GLU	2.8
1	A	888	LEU	2.8
1	C	277	PHE	2.8
3	E	60	ASP	2.8
1	C	248	GLY	2.8
1	C	294	GLU	2.8
1	C	392	ASP	2.7
1	A	887	ALA	2.7
2	B	114	G	2.7
1	C	575	LEU	2.7
2	B	41	U	2.7
1	C	716	ARG	2.7
2	B	40	C	2.7
1	A	752	ALA	2.7
1	C	391	THR	2.7
1	C	418	ASP	2.7
1	C	919	GLN	2.6
1	A	247	SER	2.6
1	C	897	ASP	2.6
1	C	154	GLU	2.6
1	C	214	LEU	2.6
1	C	260	PHE	2.6
1	C	47	GLU	2.6
1	A	877	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	926	GLU	2.6
1	C	678	CYS	2.5
1	C	142	SER	2.5
1	A	267	ALA	2.5
1	A	259	THR	2.5
1	A	1053	LYS	2.5
1	A	996	ASN	2.5
1	C	186	GLU	2.5
1	C	279	TRP	2.5
1	C	333	LYS	2.5
2	B	45	A	2.5
1	A	143	GLN	2.5
1	C	922	ALA	2.5
1	C	865	THR	2.5
1	A	609	SER	2.4
3	F	0	SER	2.4
1	C	174	PHE	2.4
3	F	12	ASN	2.4
1	C	82	THR	2.4
2	B	47	G	2.4
1	A	342	ALA	2.4
2	D	43	G	2.4
1	C	913	ALA	2.4
1	A	156	GLY	2.4
1	A	94	ALA	2.4
1	A	181	ALA	2.4
1	C	621	GLU	2.4
1	C	906	PRO	2.3
1	C	234	GLY	2.3
1	C	94	ALA	2.3
1	C	446	ALA	2.3
1	C	894	ALA	2.3
1	C	441	TYR	2.3
1	A	884	LEU	2.3
1	C	598	ASP	2.3
1	A	610	GLU	2.3
2	D	115	C	2.3
1	C	328	ASP	2.3
1	A	441	TYR	2.2
1	A	782	GLU	2.2
1	A	709	PHE	2.2
1	C	339	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	176	THR	2.2
1	C	247	SER	2.2
2	D	117	U	2.2
1	C	210	ALA	2.2
1	C	940	GLY	2.2
1	C	205	ARG	2.2
1	C	171	THR	2.2
1	C	204	SER	2.2
2	D	37	U	2.2
2	D	116	U	2.2
2	B	36	C	2.1
1	A	21	SER	2.1
1	A	148	GLY	2.1
1	C	655	GLN	2.1
1	C	236	GLU	2.1
1	C	634	GLU	2.1
1	A	1073	CYS	2.1
1	C	162	VAL	2.1
1	C	885	TYR	2.1
1	A	1051	GLY	2.1
2	B	46	A	2.1
2	D	32	C	2.1
1	A	198	ASP	2.1
1	C	667	LEU	2.1
1	C	336	ARG	2.1
1	C	458	THR	2.1
1	C	223	GLY	2.1
1	A	886	GLU	2.1
1	C	341	ASN	2.1
1	C	439	LYS	2.1
1	C	872	LEU	2.1
3	F	1	MET	2.1
1	A	768	GLN	2.1
1	A	262	PRO	2.1
1	C	395	ILE	2.0
1	C	168	ALA	2.0
2	B	133	A	2.0
1	A	190	GLY	2.0
1	A	438	GLY	2.0
1	C	325	GLY	2.0
1	C	211	GLU	2.0
2	B	42	C	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	392	ASP	2.0
1	C	335	LEU	2.0
1	C	408	LEU	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.