



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:01 PM UTC

PDB ID : 4OO8 / pdb_00004oo8
Title : Crystal structure of Streptococcus pyogenes Cas9 in complex with guide RNA and target DNA
Authors : Nishimasu, H.; Ishitani, R.; Nureki, O.
Deposited on : 2014-01-31
Resolution : 2.50 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

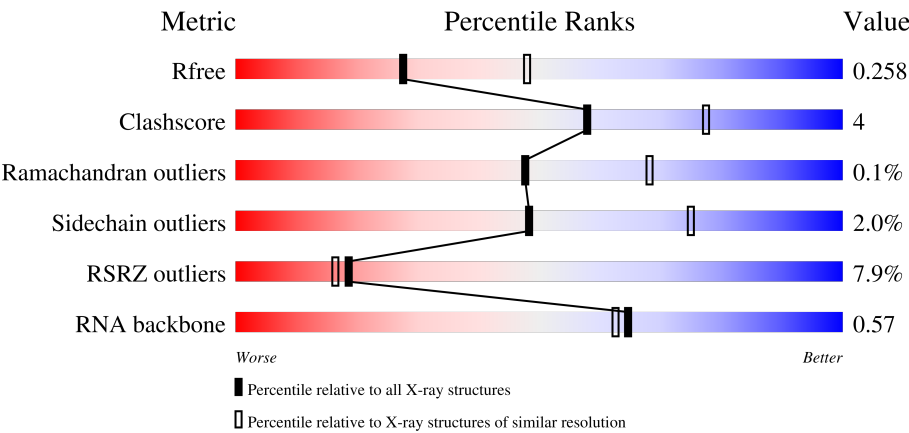
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)
RNA backbone	3983	1003 (2.78-2.22)

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	1372	
1	D	1372	
2	B	98	
2	E	98	

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Mol	Chain	Length	Quality of chain
3	C	23	4% 57% 35% 9%
3	F	23	65% 35%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24153 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/Csn1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1301	Total	C	N	O	S	0	0	0
			9999	6388	1708	1883	20			
1	D	1163	Total	C	N	O	S	0	0	0
			8998	5754	1529	1698	17			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q99ZW2
A	-2	SER	-	expression tag	UNP Q99ZW2
A	-1	GLY	-	expression tag	UNP Q99ZW2
A	0	HIS	-	expression tag	UNP Q99ZW2
A	10	ALA	ASP	engineered mutation	UNP Q99ZW2
A	80	LEU	CYS	engineered mutation	UNP Q99ZW2
A	574	GLU	CYS	engineered mutation	UNP Q99ZW2
A	840	ALA	HIS	engineered mutation	UNP Q99ZW2
D	-3	GLY	-	expression tag	UNP Q99ZW2
D	-2	SER	-	expression tag	UNP Q99ZW2
D	-1	GLY	-	expression tag	UNP Q99ZW2
D	0	HIS	-	expression tag	UNP Q99ZW2
D	10	ALA	ASP	engineered mutation	UNP Q99ZW2
D	80	LEU	CYS	engineered mutation	UNP Q99ZW2
D	574	GLU	CYS	engineered mutation	UNP Q99ZW2
D	840	ALA	HIS	engineered mutation	UNP Q99ZW2

- Molecule 2 is a RNA chain called RNA (97-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	97	Total	C	N	O	P	0	0	0
			2082	930	380	675	97			
2	E	97	Total	C	N	O	P	0	0	0
			2082	930	380	675	97			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	21	Total 404	C 192	N 72	O 120	P 20	0	0	1
3	F	23	Total 444	C 211	N 80	O 131	P 22	0	0	1

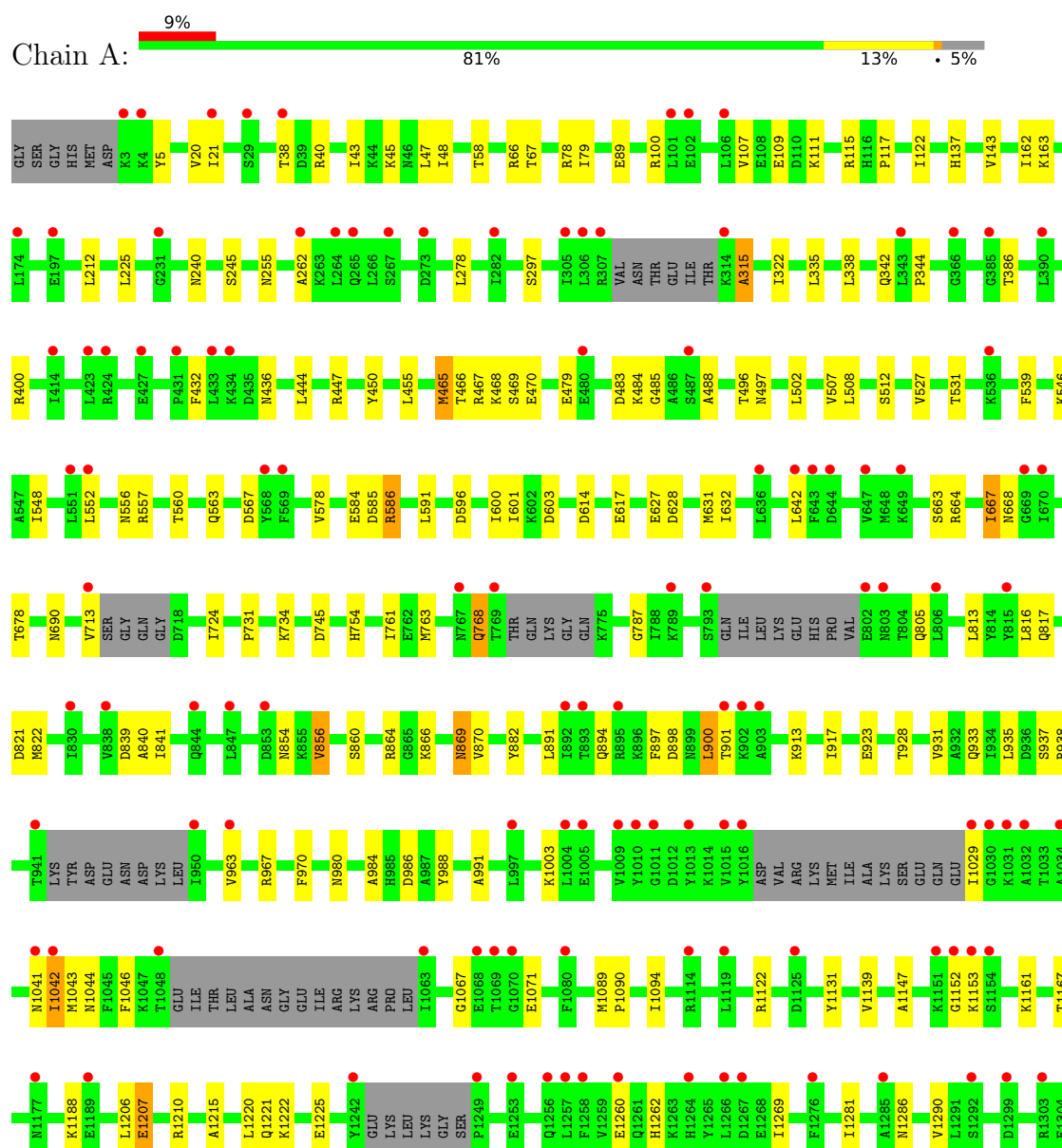
- Molecule 4 is water.

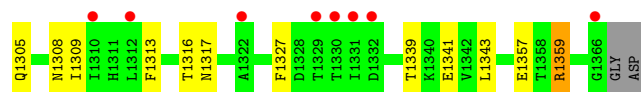
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total 32	O 32	0	0
4	B	28	Total 28	O 28	0	0
4	C	2	Total 2	O 2	0	0
4	D	49	Total 49	O 49	0	0
4	E	32	Total 32	O 32	0	0
4	F	1	Total 1	O 1	0	0

3 Residue-property plots

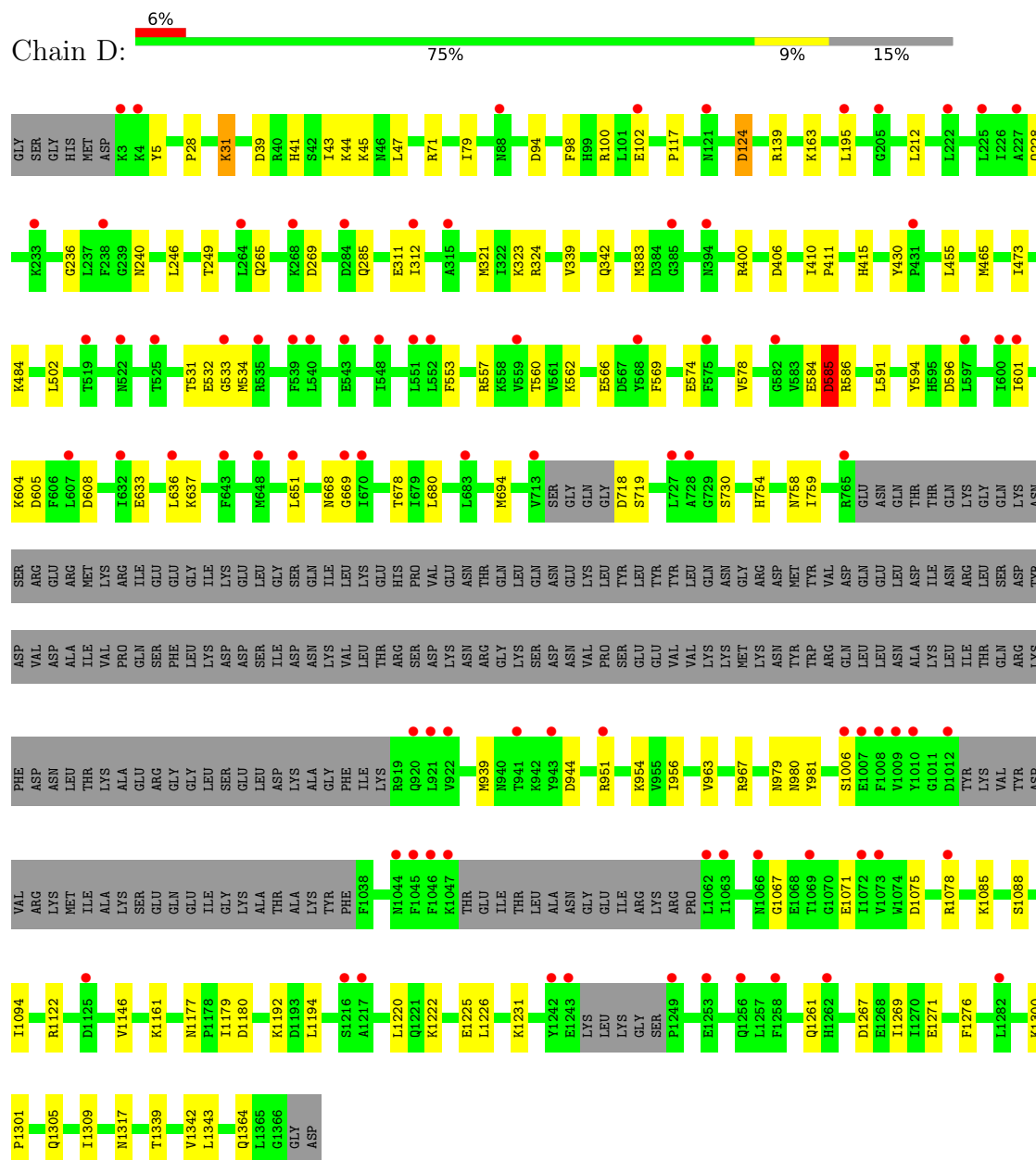
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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/Csn1

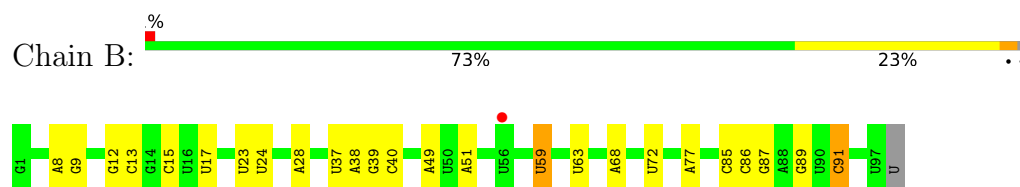




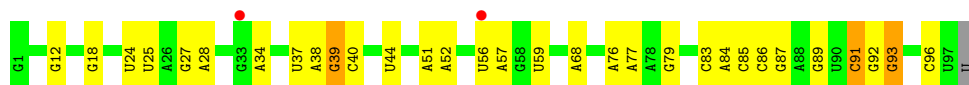
• Molecule 1: CRISPR-associated endonuclease Cas9/Csn1



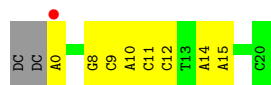
• Molecule 2: RNA (97-MER)



• Molecule 2: RNA (97-MER)



● Molecule 3: DNA (5'-D(*CP*CP*AP*GP*CP*CP*AP*AP*GP*CP*GP*CP*AP*CP*CP*TP*AP*AP*TP*TP*TP*CP*C)-3')



● Molecule 3: DNA (5'-D(*CP*CP*AP*GP*CP*CP*AP*AP*GP*CP*GP*CP*AP*CP*CP*TP*AP*AP*TP*TP*TP*CP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	76.71Å 105.69Å 126.82Å 97.68° 98.43° 100.31°	Depositor
Resolution (Å)	46.70 – 2.50 46.70 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.5 (46.70-2.50) 98.6 (46.70-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.8.3_1479	Depositor
R, R_{free}	0.222 , 0.254 0.230 , 0.258	Depositor DCC
R_{free} test set	6523 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	59.8	Xtriage
Anisotropy	0.201	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	24153	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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	0.27	0/10177	0.67	4/13792 (0.0%)
1	D	0.27	0/9164	0.67	4/12424 (0.0%)
2	B	0.11	0/2332	0.25	0/3633
2	E	0.11	0/2332	0.25	0/3633
3	C	0.43	1/451 (0.2%)	0.36	0/693
3	F	0.42	1/496 (0.2%)	0.38	0/762
All	All	0.26	2/24952 (0.0%)	0.60	8/34937 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	-2	DC	O3'-P	-8.45	1.48	1.61
3	C	0	DA	O3'-P	-8.42	1.48	1.61

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	430	TYR	CA-C-N	6.11	125.59	119.24
1	D	430	TYR	C-N-CA	6.11	125.59	119.24
1	A	502	LEU	CA-C-N	5.27	126.43	119.84
1	A	502	LEU	C-N-CA	5.27	126.43	119.84
1	D	502	LEU	CA-C-N	5.17	124.96	119.28
1	D	502	LEU	C-N-CA	5.17	124.96	119.28
1	A	315	ALA	CA-C-N	5.10	124.28	118.97
1	A	315	ALA	C-N-CA	5.10	124.28	118.97

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	9999	0	9573	102	1
1	D	8998	0	8657	77	1
2	B	2082	0	1043	8	0
2	E	2082	0	1043	14	0
3	C	404	0	225	5	0
3	F	444	0	247	5	0
4	A	32	0	0	2	0
4	B	28	0	0	0	0
4	C	2	0	0	0	0
4	D	49	0	0	0	0
4	E	32	0	0	0	0
4	F	1	0	0	0	0
All	All	24153	0	20788	198	1

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 4.

All (198) 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:D:139:ARG:HH22	1:D:415:HIS:HD2	1.29	0.78
1:D:28:PRO:HG2	1:D:47:LEU:HD12	1.73	0.69
1:D:100:ARG:NH1	1:D:117:PRO:O	2.26	0.69
1:A:967:ARG:NH1	1:A:986:ASP:OD1	2.26	0.69
1:A:21:ILE:HD12	1:A:991:ALA:HB3	1.76	0.68
1:D:980:ASN:HB2	1:D:1094:ILE:HD13	1.76	0.67
1:D:1222:LYS:NZ	1:D:1317:ASN:O	2.26	0.66
1:D:31:LYS:HE2	2:E:84:A:H62	1.61	0.66
1:A:1042:ILE:HD12	1:A:1043:MET:HG2	1.79	0.65
1:A:1222:LYS:HD2	1:A:1281:ILE:HD12	1.79	0.63
1:D:1179:ILE:HD11	1:D:1192:LYS:HD2	1.80	0.63
1:D:311:GLU:HG2	1:D:312:ILE:HG23	1.80	0.63
1:A:400:ARG:NH1	4:A:1417:HOH:O	2.32	0.62
1:D:94:ASP:OD2	1:D:100:ARG:NH2	2.33	0.62
1:D:1067:GLY:N	1:D:1071:GLU:O	2.29	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:668:ASN:ND2	1:A:678:THR:OG1	2.33	0.61
1:A:1207:GLU:OE1	1:A:1210:ARG:NH2	2.31	0.60
1:A:745:ASP:OD2	1:A:938:ARG:NH2	2.33	0.60
1:D:553:PHE:HB3	1:D:591:LEU:HD13	1.85	0.59
2:E:38:A:H3'	2:E:39:G:H5''	1.83	0.59
1:A:560:THR:HG23	1:A:563:GLN:H	1.67	0.59
1:A:980:ASN:HB2	1:A:1094:ILE:HD13	1.84	0.57
1:A:66:ARG:NH1	4:A:1405:HOH:O	2.36	0.57
1:A:432:PHE:O	1:A:436:ASN:ND2	2.36	0.57
1:D:236:GLY:O	1:D:240:ASN:ND2	2.35	0.57
1:D:730:SER:HB3	2:E:12:G:H4'	1.86	0.57
1:D:339:VAL:HG22	1:D:383:MET:HE1	1.87	0.57
1:D:956:ILE:HG23	1:D:1006:SER:HB3	1.87	0.56
2:E:92:G:HO2'	2:E:93:G:H8	1.53	0.56
1:A:467:ARG:NH1	2:B:59:U:OP1	2.38	0.56
1:D:758:ASN:HD22	1:D:954:LYS:HB2	1.71	0.56
1:D:531:THR:HG22	1:D:534:MET:HG3	1.88	0.55
1:A:584:GLU:O	1:A:586:ARG:N	2.40	0.54
1:D:718:ASP:OD1	1:D:719:SER:N	2.40	0.54
1:A:78:ARG:NH1	1:A:162:ILE:O	2.41	0.54
1:D:1300:LYS:HG3	1:D:1301:PRO:HD2	1.91	0.53
1:A:479:GLU:HG2	1:A:484:LYS:HE3	1.90	0.53
1:A:1122:ARG:NH2	2:B:49:A:N3	2.57	0.53
1:A:894:GLN:NE2	1:A:898:ASP:OD2	2.42	0.52
1:D:1300:LYS:O	1:D:1305:GLN:NE2	2.42	0.52
1:A:614:ASP:OD1	1:A:664:ARG:NH2	2.43	0.52
1:D:43:ILE:HD13	1:D:45:LYS:HE3	1.91	0.52
1:D:531:THR:OG1	1:D:532:GLU:OE1	2.27	0.52
1:D:124:ASP:OD2	1:D:124:ASP:N	2.42	0.51
1:D:71:ARG:NH1	2:E:18:G:N7	2.49	0.51
1:A:869:ASN:HD22	1:A:870:VAL:HG12	1.74	0.51
1:D:604:LYS:NZ	1:D:608:ASP:OD2	2.40	0.51
1:A:1222:LYS:HE2	1:A:1317:ASN:O	2.11	0.51
1:D:633:GLU:O	1:D:637:LYS:NZ	2.44	0.51
1:A:1067:GLY:N	1:A:1071:GLU:O	2.40	0.50
1:A:1041:ASN:HA	1:A:1044:ASN:HB2	1.94	0.50
1:D:249:THR:HG23	1:D:265:GLN:HB2	1.93	0.50
1:A:79:ILE:HD11	1:A:163:LYS:HB2	1.94	0.49
1:D:531:THR:HG23	1:D:533:GLY:H	1.76	0.49
1:D:569:PHE:O	1:D:574:GLU:N	2.44	0.49
3:C:8:DG:H2'	3:C:9:DC:C6	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:981:TYR:HD1	1:D:1094:ILE:HD11	1.77	0.49
1:A:89:GLU:HG3	1:A:432:PHE:CD1	2.48	0.49
1:D:585:ASP:OD2	1:D:585:ASP:N	2.45	0.49
3:C:9:DC:H2'	3:C:10:DA:C8	2.48	0.49
1:D:5:TYR:OH	1:D:754:HIS:O	2.30	0.49
1:A:923:GLU:HG2	1:A:928:THR:HG21	1.94	0.48
1:A:488:ALA:HB3	1:A:631:MET:HE1	1.95	0.48
1:A:1220:LEU:HG	1:A:1339:THR:HG22	1.96	0.48
1:A:531:THR:HG22	1:A:578:VAL:HG12	1.95	0.48
1:A:1207:GLU:O	1:A:1210:ARG:HG2	2.14	0.48
1:A:447:ARG:HG3	2:B:17:U:H5'	1.96	0.48
1:D:1146:VAL:HG21	1:D:1194:LEU:HD23	1.96	0.48
1:A:338:LEU:HD13	1:A:386:THR:HG22	1.96	0.47
1:A:821:ASP:OD1	1:A:822:MET:N	2.47	0.47
1:D:212:LEU:HD22	1:D:246:LEU:HD21	1.96	0.47
1:A:5:TYR:OH	1:A:754:HIS:O	2.32	0.47
1:A:763:MET:HE1	1:A:931:VAL:HG21	1.96	0.47
1:A:240:ASN:ND2	1:A:255:ASN:OD1	2.46	0.47
1:D:400:ARG:NH2	1:D:406:ASP:OD2	2.47	0.47
2:E:24:U:H2'	2:E:25:U:C6	2.50	0.47
2:E:86:C:H2'	2:E:87:G:O4'	2.15	0.47
1:A:21:ILE:HD11	1:A:988:TYR:CD1	2.50	0.47
2:E:27:G:N2	2:E:44:U:OP2	2.48	0.47
1:D:44:LYS:HD2	2:E:92:G:N7	2.30	0.46
1:A:497:ASN:HD21	3:C:11:DC:P	2.37	0.46
1:A:1308:ASN:ND2	1:A:1327:PHE:H	2.13	0.46
1:A:560:THR:HA	1:A:586:ARG:HA	1.96	0.46
1:A:897:PHE:O	1:A:901:THR:HG22	2.15	0.46
1:D:694:MET:HA	1:D:694:MET:HE2	1.98	0.46
1:A:839:ASP:OD2	1:A:840:ALA:N	2.48	0.46
1:D:39:ASP:O	1:D:41:HIS:ND1	2.36	0.46
1:A:864:ARG:NH2	1:A:866:LYS:O	2.49	0.46
1:D:1085:LYS:O	1:D:1088:SER:OG	2.28	0.46
2:E:92:G:O2'	2:E:93:G:H8	1.98	0.46
1:A:933:GLN:NE2	1:A:937:SER:OG	2.48	0.46
2:E:85:C:H2'	2:E:86:C:C6	2.51	0.46
1:A:100:ARG:HB3	1:A:117:PRO:HA	1.97	0.45
1:A:450:TYR:OH	1:A:627:GLU:OE2	2.26	0.45
1:A:468:LYS:HG3	1:A:483:ASP:HB2	1.98	0.45
1:D:560:THR:HG22	1:D:562:LYS:H	1.80	0.45
1:A:245:SER:HA	1:A:297:SER:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:854:ASN:O	1:A:854:ASN:ND2	2.48	0.45
1:D:269:ASP:OD1	1:D:269:ASP:N	2.49	0.45
1:A:768:GLN:H	1:A:768:GLN:HG3	1.56	0.45
1:D:1146:VAL:HG22	1:D:1161:LYS:HG3	1.99	0.45
1:D:1161:LYS:HE3	1:D:1343:LEU:HB3	1.99	0.45
3:C:11:DC:H2'	3:C:12:DC:C6	2.52	0.45
1:D:668:ASN:ND2	1:D:678:THR:OG1	2.50	0.45
1:A:817:GLN:O	1:A:882:TYR:OH	2.34	0.45
1:A:45:LYS:NZ	1:A:1357:GLU:OE2	2.48	0.44
1:A:663:SER:O	1:A:667:ILE:HG22	2.17	0.44
1:A:1139:VAL:HA	1:A:1167:THR:HA	1.99	0.44
1:D:1220:LEU:HG	1:D:1339:THR:HG22	1.99	0.44
1:D:1305:GLN:O	1:D:1309:ILE:HG12	2.17	0.44
1:A:841:ILE:HD11	1:A:856:VAL:HG23	1.99	0.44
3:F:12:DC:H2'	3:F:13:DT:C6	2.53	0.44
1:D:139:ARG:HH22	1:D:415:HIS:CD2	2.20	0.44
1:D:321:MET:HE1	1:D:324:ARG:NH2	2.33	0.44
1:D:951:ARG:HD3	1:D:951:ARG:HA	1.84	0.44
1:D:1177:ASN:HB3	1:D:1180:ASP:HB3	2.00	0.44
1:A:628:ASP:O	1:A:632:ILE:HG12	2.17	0.44
1:D:605:ASP:N	1:D:605:ASP:OD1	2.50	0.44
1:A:38:THR:HG23	1:A:40:ARG:H	1.83	0.44
1:A:107:VAL:HG23	1:A:1131:TYR:CZ	2.52	0.44
1:A:512:SER:OG	1:A:617:GLU:OE2	2.28	0.44
1:A:1161:LYS:HD2	1:A:1343:LEU:HB3	1.99	0.44
1:A:465:MET:HG2	1:A:466:THR:N	2.33	0.44
3:C:14:DA:H2''	3:C:15:DA:C8	2.53	0.44
1:D:1269:ILE:HD12	1:D:1309:ILE:HG21	2.00	0.43
1:A:485:GLY:HA2	1:A:631:MET:HE3	2.01	0.43
1:D:584:GLU:O	1:D:586:ARG:N	2.51	0.43
1:D:1075:ASP:OD1	1:D:1078:ARG:N	2.52	0.43
1:A:970:PHE:HZ	1:A:1046:PHE:HB3	1.83	0.43
1:D:591:LEU:HB3	1:D:594:TYR:HB3	2.00	0.43
1:D:979:ASN:HB2	1:D:1225:GLU:HG3	2.00	0.43
1:A:48:ILE:HG12	1:A:984:ALA:HB1	2.01	0.43
1:A:508:LEU:HD21	1:A:664:ARG:HB2	2.00	0.43
1:A:539:PHE:HB3	1:A:690:ASN:ND2	2.33	0.43
1:A:557:ARG:NH2	1:A:596:ASP:OD1	2.28	0.43
1:A:601:ILE:HG13	1:A:603:ASP:H	1.84	0.43
1:D:557:ARG:HH22	1:D:596:ASP:CG	2.23	0.43
1:A:43:ILE:HD11	2:B:91:C:H5'	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:813:LEU:HD23	1:A:816:LEU:HD12	2.00	0.43
1:D:981:TYR:CD1	1:D:1094:ILE:HD11	2.54	0.43
1:A:724:ILE:O	1:A:734:LYS:HE3	2.19	0.43
1:A:841:ILE:HG12	1:A:900:LEU:HD22	2.01	0.43
1:A:913:LYS:O	1:A:917:ILE:HG12	2.19	0.43
1:D:759:ILE:HD11	1:D:939:MET:HE3	2.01	0.43
1:A:787:GLY:HA3	1:A:891:LEU:HD21	2.00	0.42
1:D:963:VAL:O	1:D:967:ARG:HG3	2.19	0.42
1:A:761:ILE:HD11	1:A:935:LEU:HD12	2.00	0.42
1:D:43:ILE:HD11	2:E:91:C:H5'	2.00	0.42
3:F:2:DC:H2'	3:F:3:DC:C6	2.54	0.42
1:A:66:ARG:NH2	2:B:15:C:OP2	2.52	0.42
1:A:465:MET:HE2	1:A:465:MET:HB3	1.81	0.42
1:D:98:PHE:O	1:D:102:GLU:HG2	2.20	0.42
1:D:465:MET:HE3	1:D:465:MET:HB3	1.97	0.42
1:A:342:GLN:C	1:A:344:PRO:HD3	2.45	0.42
1:A:1357:GLU:OE1	1:A:1359:ARG:NH1	2.46	0.42
1:D:31:LYS:HD3	2:E:83:C:N4	2.34	0.42
1:A:137:HIS:HA	1:A:322:ILE:HD11	2.01	0.42
1:A:143:VAL:HG11	1:A:315:ALA:HB2	2.01	0.42
1:D:455:LEU:HB3	1:D:473:ILE:HD12	2.01	0.42
1:A:58:THR:HA	1:A:731:PRO:HG2	2.01	0.42
1:A:212:LEU:HD21	1:A:225:LEU:HD13	2.01	0.42
1:A:596:ASP:O	1:A:600:ILE:HG12	2.19	0.42
1:D:79:ILE:HD11	1:D:163:LYS:HB2	2.01	0.42
1:D:1226:LEU:HD13	1:D:1276:PHE:CG	2.55	0.42
1:A:548:ILE:HG23	1:A:552:LEU:HD12	2.02	0.42
3:F:2:DC:H2'	3:F:3:DC:H6	1.84	0.42
1:A:1089:MET:HA	1:A:1090:PRO:HD3	1.94	0.42
1:A:1152:GLY:HA2	1:A:1153:LYS:HA	1.81	0.41
1:A:262:ALA:HB1	1:A:278:LEU:HD13	2.01	0.41
2:B:12:G:H2'	2:B:13:C:C6	2.55	0.41
1:D:668:ASN:HA	1:D:669:GLY:HA3	1.87	0.41
1:D:1122:ARG:O	2:E:52:A:H5''	2.20	0.41
1:A:162:ILE:HD13	1:A:444:LEU:HD12	2.02	0.41
1:A:1147:ALA:HB1	1:A:1188:LYS:O	2.21	0.41
1:A:1206:LEU:HD21	1:A:1341:GLU:HB3	2.01	0.41
1:D:342:GLN:HE22	1:D:383:MET:HA	1.85	0.41
3:F:15:DA:H2'	3:F:16:DT:C6	2.54	0.41
1:A:1215:ALA:HB2	1:A:1221:GLN:HG3	2.03	0.41
1:D:410:ILE:HA	1:D:411:PRO:HD3	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:LYS:HD2	1:A:115:ARG:HA	2.03	0.41
1:D:195:LEU:HD21	1:D:285:GLN:O	2.20	0.41
3:F:11:DC:H2'	3:F:12:DC:H6	1.85	0.41
1:A:469:SER:OG	1:A:470:GLU:N	2.53	0.41
1:A:963:VAL:O	1:A:967:ARG:HG3	2.21	0.41
1:A:1286:ASN:O	1:A:1290:VAL:HG23	2.21	0.41
1:D:636:LEU:HD21	1:D:651:LEU:HD23	2.03	0.41
1:D:1231:LYS:H	1:D:1231:LYS:HG3	1.66	0.41
1:A:20:VAL:HB	1:A:47:LEU:HB3	2.03	0.41
1:A:1094:ILE:HD12	1:A:1225:GLU:HG3	2.03	0.41
1:A:1305:GLN:O	1:A:1309:ILE:HG12	2.21	0.41
1:D:668:ASN:HD22	1:D:680:LEU:HB3	1.86	0.41
1:A:805:GLN:HE21	1:A:805:GLN:HB2	1.71	0.40
1:D:1267:ASP:O	1:D:1271:GLU:HG2	2.21	0.40
1:A:109:GLU:OE2	1:A:109:GLU:N	2.52	0.40
1:A:1313:PHE:HA	1:A:1316:THR:HG22	2.03	0.40
2:B:23:U:H2'	2:B:24:U:C6	2.56	0.40
2:B:85:C:H2'	2:B:86:C:C6	2.57	0.40
1:D:1364:GLN:N	1:D:1364:GLN:OE1	2.54	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:860:SER:OG	1:D:944:ASP:O[1_454]	2.16	0.04

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	1283/1372 (94%)	1252 (98%)	30 (2%)	1 (0%)	48 68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	D	1151/1372 (84%)	1108 (96%)	42 (4%)	1 (0%)	48 68
All	All	2434/2744 (89%)	2360 (97%)	72 (3%)	2 (0%)	48 68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	585	ASP
1	A	585	ASP

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	1013/1227 (83%)	985 (97%)	28 (3%)	38 66
1	D	921/1227 (75%)	910 (99%)	11 (1%)	63 83
All	All	1934/2454 (79%)	1895 (98%)	39 (2%)	48 75

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	THR
1	A	122	ILE
1	A	335	LEU
1	A	455	LEU
1	A	465	MET
1	A	496	THR
1	A	507	VAL
1	A	527	VAL
1	A	546	LYS
1	A	556	ASN
1	A	567	ASP
1	A	586	ARG
1	A	591	LEU
1	A	642	LEU

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Mol	Chain	Res	Type
1	A	667	ILE
1	A	713	VAL
1	A	768	GLN
1	A	856	VAL
1	A	869	ASN
1	A	900	LEU
1	A	1003	LYS
1	A	1029	ILE
1	A	1042	ILE
1	A	1207	GLU
1	A	1260	GLU
1	A	1262	HIS
1	A	1269	ILE
1	A	1359	ARG
1	D	31	LYS
1	D	124	ASP
1	D	228	GLN
1	D	323	LYS
1	D	484	LYS
1	D	566	GLU
1	D	578	VAL
1	D	585	ASP
1	D	601	ILE
1	D	1261	GLN
1	D	1342	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	GLN
1	A	281	GLN
1	A	285	GLN
1	A	407	ASN
1	A	497	ASN
1	A	668	ASN
1	A	690	ASN
1	A	726	ASN
1	A	758	ASN
1	A	805	GLN
1	A	869	ASN
1	A	899	ASN
1	A	933	GLN

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Mol	Chain	Res	Type
1	A	982	HIS
1	A	1308	ASN
1	D	342	GLN
1	D	354	GLN
1	D	412	HIS
1	D	415	HIS
1	D	426	GLN
1	D	501	ASN
1	D	668	ASN
1	D	723	HIS
1	D	758	ASN
1	D	985	HIS
1	D	1221	GLN
1	D	1254	GLN
1	D	1311	HIS
1	D	1350	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	96/98 (97%)	15 (15%)	1 (1%)
2	E	96/98 (97%)	17 (17%)	0
All	All	192/196 (97%)	32 (16%)	1 (0%)

All (32) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	8	A
2	B	9	G
2	B	28	A
2	B	37	U
2	B	39	G
2	B	40	C
2	B	51	A
2	B	59	U
2	B	63	U
2	B	68	A
2	B	72	U
2	B	77	A
2	B	87	G
2	B	89	G

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Mol	Chain	Res	Type
2	B	91	C
2	E	28	A
2	E	34	A
2	E	37	U
2	E	39	G
2	E	40	C
2	E	51	A
2	E	56	U
2	E	57	A
2	E	59	U
2	E	68	A
2	E	76	A
2	E	77	A
2	E	79	G
2	E	89	G
2	E	91	C
2	E	93	G
2	E	96	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	38	A

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 ⓘ

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	1301/1372 (94%)	0.84	124 (9%) 14 12	36, 82, 120, 160	0
1	D	1163/1372 (84%)	0.70	85 (7%) 21 19	40, 75, 117, 166	0
2	B	97/98 (98%)	0.14	1 (1%) 79 76	40, 75, 146, 168	0
2	E	97/98 (98%)	0.14	2 (2%) 63 59	44, 77, 172, 189	0
3	C	21/23 (91%)	0.13	1 (4%) 35 31	49, 65, 102, 108	0
3	F	23/23 (100%)	0.08	0 100 100	53, 64, 131, 152	0
All	All	2702/2986 (90%)	0.72	213 (7%) 18 16	36, 78, 121, 189	0

All (213) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1046	PHE	6.0
1	A	1242	TYR	5.5
1	D	1045	PHE	5.4
1	A	950	ILE	5.3
1	A	431	PRO	4.8
1	D	1047	LYS	4.8
1	A	1029	ILE	4.7
1	A	643	PHE	4.1
1	A	1009	VAL	4.0
1	D	1007	GLU	4.0
1	D	3	LYS	4.0
1	A	1042	ILE	3.8
1	A	1004	LEU	3.8
1	A	769	THR	3.7
1	A	1011	GLY	3.7
1	D	552	LEU	3.6
1	D	1062	LEU	3.6
1	D	597	LEU	3.5
1	A	3	LYS	3.4

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Mol	Chain	Res	Type	RSRZ
1	D	713	VAL	3.4
1	D	921	LEU	3.3
1	D	600	ILE	3.3
1	D	1256	GLN	3.3
1	A	815	TYR	3.2
1	D	1010	TYR	3.2
1	A	669	GLY	3.2
1	D	4	LYS	3.2
1	A	265	GLN	3.2
1	D	533	GLY	3.2
1	D	1063	ILE	3.2
1	D	284	ASP	3.1
1	D	1009	VAL	3.1
1	D	582	GLY	3.1
1	A	803	ASN	3.0
1	A	1152	GLY	3.0
1	A	1015	VAL	3.0
1	A	1005	GLU	3.0
1	A	427	GLU	3.0
1	D	551	LEU	3.0
1	A	4	LYS	3.0
1	A	1257	LEU	2.9
1	A	1312	LEU	2.9
1	D	651	LEU	2.9
1	A	1253	GLU	2.9
1	D	648	MET	2.9
1	A	366	GLY	2.9
1	A	307	ARG	2.9
1	A	433	LEU	2.9
1	A	1030	GLY	2.9
1	D	568	TYR	2.9
1	A	262	ALA	2.8
1	A	1016	TYR	2.8
1	D	225	LEU	2.8
1	A	853	ASP	2.8
1	A	649	LYS	2.8
1	D	643	PHE	2.8
1	D	385	GLY	2.8
1	A	1010	TYR	2.8
1	A	1249	PRO	2.8
1	D	264	LEU	2.8
1	A	1031	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	1243	GLU	2.8
1	A	231	GLY	2.7
1	D	195	LEU	2.7
1	D	233	LYS	2.7
1	D	205	GLY	2.7
1	A	1063	ILE	2.7
1	A	901	THR	2.7
1	D	1008	PHE	2.7
1	D	431	PRO	2.7
1	A	1153	LYS	2.7
1	A	174	LEU	2.7
1	D	268	LYS	2.6
1	D	670	ILE	2.6
1	D	1044	ASN	2.6
1	D	1217	ALA	2.6
1	D	522	ASN	2.6
1	A	1032	ALA	2.6
1	A	1330	THR	2.6
1	A	1267	ASP	2.6
1	D	1125	ASP	2.6
1	A	264	LEU	2.6
1	D	88	ASN	2.6
1	A	793	SER	2.6
1	D	1006	SER	2.5
1	A	1125	ASP	2.5
1	A	789	LYS	2.5
1	D	1012	ASP	2.5
1	A	552	LEU	2.5
1	A	487	SER	2.5
1	D	941	THR	2.5
1	D	669	GLY	2.5
1	A	390	LEU	2.5
1	A	197	GLU	2.4
1	A	1299	ASP	2.4
1	A	844	GLN	2.4
1	D	238	PHE	2.4
1	A	1303	ARG	2.4
1	D	765	ARG	2.4
1	D	1066	ASN	2.4
2	E	56	U	2.4
1	A	267	SER	2.4
1	A	802	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	644	ASP	2.4
1	A	830	ILE	2.4
1	D	539	PHE	2.4
1	A	963	VAL	2.4
1	D	922	VAL	2.4
1	A	997	LEU	2.4
1	D	1262	HIS	2.4
1	A	1154	SER	2.4
1	A	21	ILE	2.4
3	C	0	DA	2.4
1	A	414	ILE	2.3
1	A	1256	GLN	2.3
1	A	1264	HIS	2.3
1	D	540	LEU	2.3
1	A	1151	LYS	2.3
1	A	941	THR	2.3
1	D	920	GLN	2.3
1	A	551	LEU	2.3
1	D	636	LEU	2.3
1	A	536	LYS	2.3
1	D	1242	TYR	2.3
1	A	767	ASN	2.3
1	D	315	ALA	2.3
1	D	1282	LEU	2.3
1	A	273	ASP	2.3
1	A	314	LYS	2.3
1	A	1080	PHE	2.3
1	A	1189	GLU	2.3
1	D	1073	VAL	2.3
1	A	1322	ALA	2.3
1	D	951	ARG	2.3
1	A	343	LEU	2.3
1	A	670	ILE	2.3
1	D	1072	ILE	2.3
1	A	569	PHE	2.3
1	A	1114	ARG	2.3
1	A	1177	ASN	2.2
1	D	1078	ARG	2.3
1	A	1069	THR	2.2
1	A	306	LEU	2.2
1	A	423	LEU	2.2
1	A	38	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	1069	THR	2.2
1	A	1070	GLY	2.2
1	A	305	ILE	2.2
1	A	1331	ILE	2.2
1	A	1068	GLU	2.2
1	A	895	ARG	2.2
1	A	106	LEU	2.2
1	A	642	LEU	2.2
1	D	607	LEU	2.2
1	A	1285	ALA	2.2
1	D	727	LEU	2.2
1	D	632	ILE	2.2
1	A	1260	GLU	2.2
1	A	424	ARG	2.2
1	D	1249	PRO	2.2
1	A	902	LYS	2.2
1	D	121	ASN	2.1
1	A	636	LEU	2.1
1	D	102	GLU	2.1
1	A	29	SER	2.1
1	A	1258	PHE	2.1
1	A	1276	PHE	2.1
1	D	575	PHE	2.1
1	D	227	ALA	2.1
1	A	101	LEU	2.1
1	A	1266	LEU	2.1
1	A	1310	ILE	2.1
1	D	543	GLU	2.1
1	A	1292	SER	2.1
1	A	1332	ASP	2.1
1	A	903	ALA	2.1
1	D	394	ASN	2.1
1	D	222	LEU	2.1
1	A	480	GLU	2.1
1	A	892	ILE	2.1
1	A	1013	TYR	2.1
1	A	434	LYS	2.1
1	A	838	VAL	2.1
1	D	559	VAL	2.1
1	D	728	ALA	2.1
2	B	56	U	2.1
1	A	893	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	1329	THR	2.1
1	D	519	THR	2.1
1	D	525	THR	2.1
1	D	1253	GLU	2.1
1	D	548	ILE	2.1
1	D	601	ILE	2.1
2	E	33	G	2.0
1	A	647	VAL	2.0
1	D	1258	PHE	2.0
1	A	1034	ALA	2.0
1	D	535	ARG	2.0
1	A	806	LEU	2.0
1	A	847	LEU	2.0
1	A	1041	ASN	2.0
1	A	102	GLU	2.0
1	A	1366	GLY	2.0
1	D	312	ILE	2.0
1	A	568	TYR	2.0
1	A	713	VAL	2.0
1	A	1119	LEU	2.0
1	D	683	LEU	2.0
1	A	385	GLY	2.0
1	A	282	ILE	2.0
1	A	1048	THR	2.0
1	D	943	TYR	2.0
1	D	1216	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.