Full wwPDB X-ray Structure Validation Report (i)

Nov 15, 2023 – 02:57 PM JST

PDB ID	:	6JE3
Title	:	Crystal structure of Nme2Cas9 in complex with sgRNA and target DNA (AG-GCCC PAM) with 5 nt overhang
Authors	:	Sun, W.; Yang, J.; Cheng, Z.; Liu, C.; Wang, K.; Huang, X.; Wang, Y.
Deposited on	:	2019-02-03
Resolution	:	2.93 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 (i) 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 (i)) were used in the production of this report:

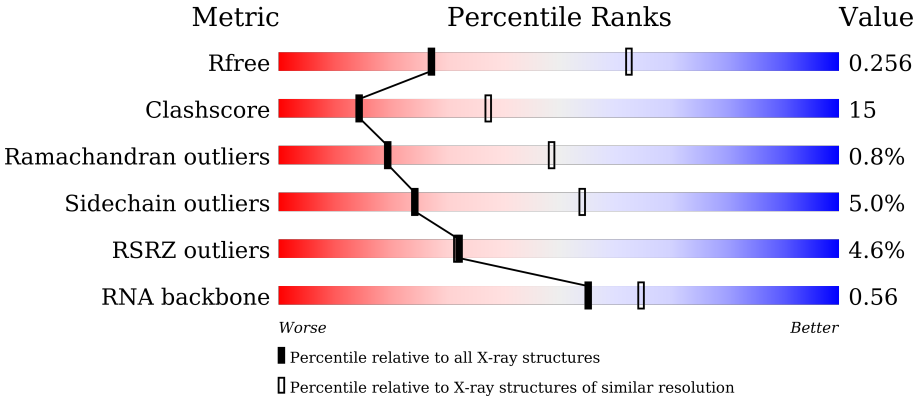
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.93 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}	130704	2969 (2.98-2.90)
Clashscore	141614	3218 (2.98-2.90)
Ramachandran outliers	138981	3122 (2.98-2.90)
Sidechain outliers	138945	3124 (2.98-2.90)
RSRZ outliers	127900	2902 (2.98-2.90)
RNA backbone	3102	1060 (3.20-2.68)

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 <=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	1083	4%, 53%, 24%, •, 21%
2	B	135	4%, 32%, 43%, 11%, •, 13%
3	C	35	40%, 60%
4	D	16	31%, 50%, 19%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 9806 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
			Total	C	N	O	S			
1	A	858	6327	3994	1154	1159	20	0	0	0

- Molecule 2 is a RNA chain called sgRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	117	2472	1106	424	825	117	0	0	0

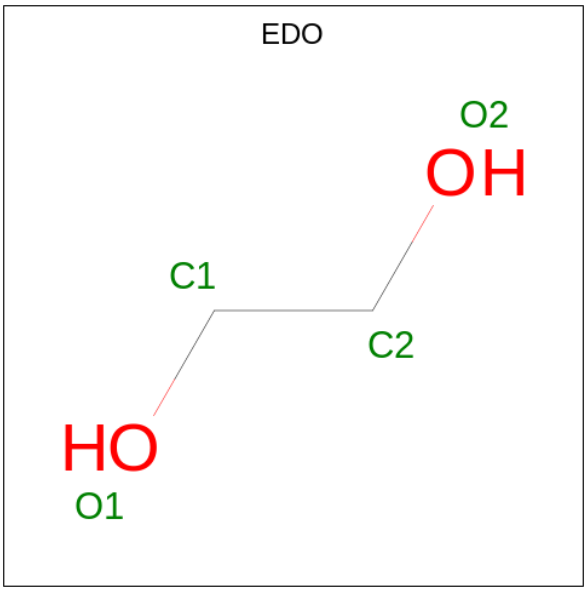
- Molecule 3 is a DNA chain called target DNA strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	C	35	720	344	139	203	34	0	0	0

- Molecule 4 is a DNA chain called non-target DNA strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
4	D	13	268	127	50	78	13	0	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	4	2	2	0	0

- Molecule 6 is water.

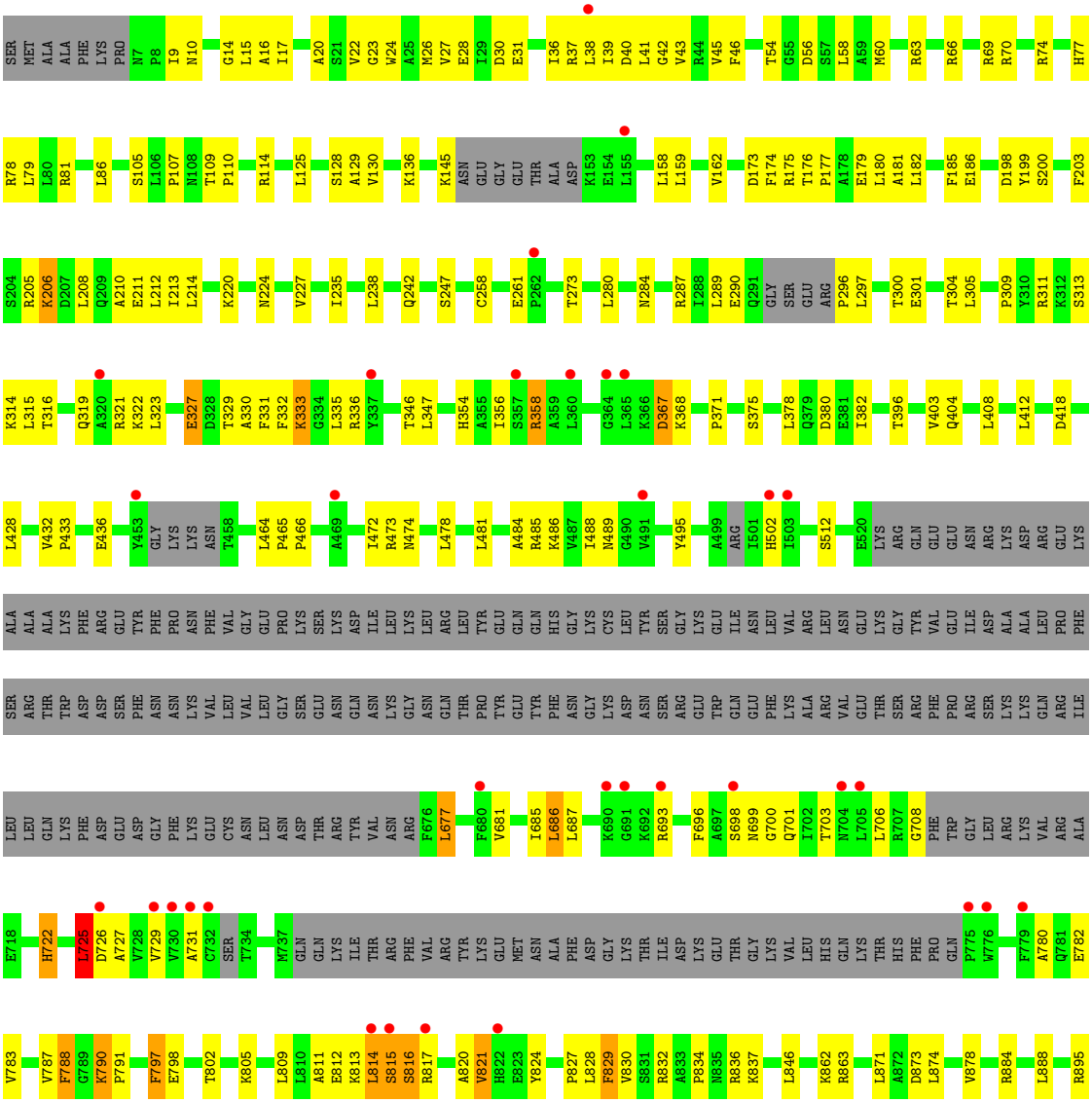
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	7	7	7	0	0
			Total	O		
6	B	5	5	5	0	0
			Total	O		
6	C	3	3	3	0	0

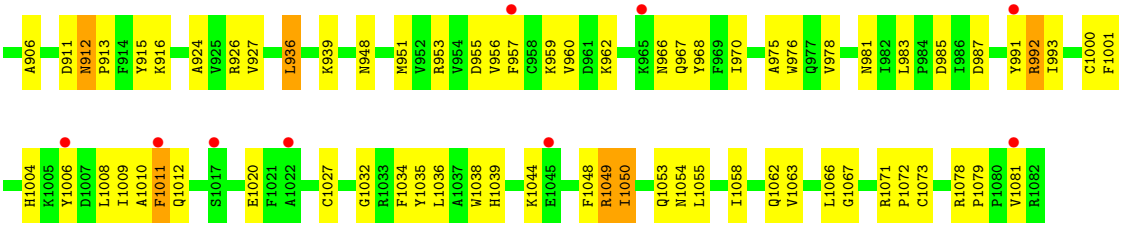
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: CRISPR-associated endonuclease Cas9

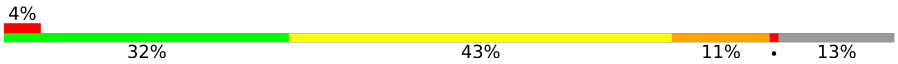
Chain A:

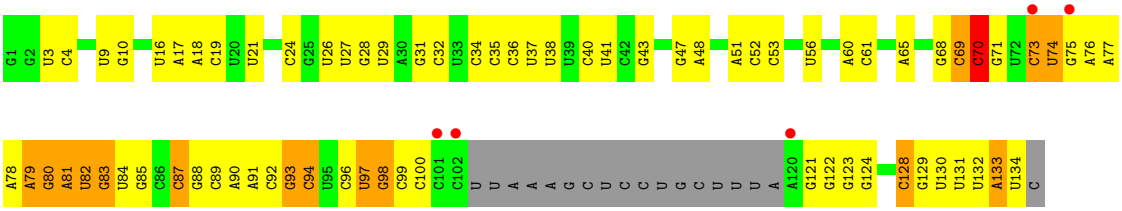




- Molecule 2: sgRNA

Chain B:



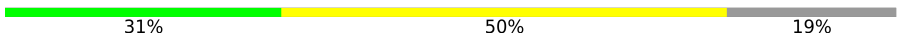


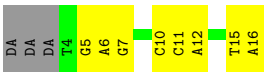
- Molecule 3: target DNA strand

Chain C:

- Molecule 4: non-target DNA strand

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	166.39Å 165.97Å 178.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.60 – 2.93 41.60 – 2.93	Depositor EDS
% Data completeness (in resolution range)	75.3 (41.60-2.93) 75.3 (41.60-2.93)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.95Å)	Xtriage
Refinement program	PHENIX (1.14_3247: ???)	Depositor
R, R_{free}	0.227 , 0.256 0.227 , 0.256	Depositor DCC
R_{free} test set	1906 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å^2)	49.1	Xtriage
Anisotropy	0.548	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.25 , 54.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.038 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	9806	wwPDB-VP
Average B, all atoms (Å^2)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.30	0/6437	0.56	2/8721 (0.0%)
2	B	0.32	0/2757	0.91	4/4287 (0.1%)
3	C	0.61	0/810	0.93	1/1250 (0.1%)
4	D	0.66	0/300	0.97	0/461
All	All	0.35	0/10304	0.73	7/14719 (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	A	0	5

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	725	LEU	CA-CB-CG	6.63	130.54	115.30
2	B	70	C	C5-C6-N1	6.12	124.06	121.00
1	A	791	PRO	N-CA-CB	5.78	110.23	103.30
2	B	87	C	C2-N1-C1'	5.33	124.66	118.80
3	C	17	DG	O4'-C4'-C3'	-5.30	102.38	104.50
2	B	70	C	C6-N1-C2	-5.18	118.23	120.30
2	B	87	C	C5-C6-N1	5.09	123.55	121.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1010	ALA	Peptide
1	A	512	SER	Peptide
1	A	814	LEU	Peptide
1	A	820	ALA	Peptide
1	A	821	VAL	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6327	0	5883	213	0
2	B	2472	0	1256	48	0
3	C	720	0	392	17	0
4	D	268	0	147	5	0
5	A	4	0	6	0	0
6	A	7	0	0	0	0
6	B	5	0	0	0	0
6	C	3	0	0	0	0
All	All	9806	0	7684	269	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 15.

All (269) 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:687:LEU:HD22	1:A:693:ARG:HA	1.35	1.08
1:A:305:LEU:HD23	1:A:323:LEU:HD13	1.37	1.02
1:A:358:ARG:NH1	1:A:368:LYS:HE3	1.77	0.99
1:A:289:LEU:O	1:A:331:PHE:N	2.01	0.94
1:A:687:LEU:HD12	1:A:687:LEU:O	1.67	0.93
1:A:42:GLY:HA2	1:A:828:LEU:HD13	1.50	0.93
1:A:305:LEU:CD2	1:A:323:LEU:HD13	1.99	0.92
1:A:358:ARG:HH12	1:A:368:LYS:HE3	1.33	0.89
1:A:289:LEU:N	1:A:331:PHE:O	2.07	0.87
1:A:333:LYS:HD3	3:C:28:DG:H5"	1.57	0.87
1:A:290:GLU:CB	1:A:330:ALA:HB2	2.17	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:LEU:HD23	1:A:323:LEU:O	1.89	0.72
1:A:321:ARG:NH2	1:A:330:ALA:O	2.24	0.70
1:A:834:PRO:HA	1:A:1006:TYR:HB2	1.74	0.70
2:B:69:C:H2'	2:B:70:C:H5"	1.72	0.70
1:A:63:ARG:NH2	2:B:18:A:OP2	2.24	0.70
1:A:481:LEU:HD13	1:A:677:LEU:HB2	1.73	0.70
1:A:725:LEU:HD23	1:A:726:ASP:H	1.57	0.70
1:A:722:HIS:O	1:A:725:LEU:HD22	1.92	0.70
1:A:284:ASN:O	1:A:287:ARG:NH2	2.25	0.69
3:C:34:DC:H2"	3:C:35:DC:H5'	1.74	0.69
1:A:955:ASP:HB3	1:A:1078:ARG:HH12	1.58	0.69
1:A:15:LEU:HA	1:A:23:GLY:O	1.93	0.68
1:A:212:LEU:HD11	1:A:235:ILE:HG22	1.76	0.67
1:A:978:VAL:HG23	1:A:1081:VAL:HB	1.77	0.67
1:A:948:ASN:HD22	1:A:976:TRP:HE1	1.43	0.66
1:A:884:ARG:NH2	2:B:53:C:O2	2.29	0.66
1:A:884:ARG:NH1	2:B:36:C:O2	2.29	0.66
1:A:54:THR:HG23	1:A:56:ASP:H	1.59	0.66
1:A:301:GLU:O	1:A:304:THR:OG1	2.12	0.66
1:A:60:MET:HG3	1:A:63:ARG:HH11	1.61	0.65
1:A:9:ILE:HG22	1:A:10:ASN:N	2.12	0.65
1:A:287:ARG:HD2	1:A:296:PRO:HA	1.78	0.65
1:A:28:GLU:H	1:A:36:ILE:HG22	1.60	0.65
1:A:797:PHE:HD2	1:A:798:GLU:H	1.45	0.64
1:A:39:ILE:HG23	1:A:824:TYR:HB3	1.80	0.64
1:A:685:ILE:HG12	1:A:686:LEU:HD23	1.80	0.64
1:A:327:GLU:CB	1:A:329:THR:HG22	2.27	0.64
1:A:20:ALA:O	1:A:45:VAL:HG23	1.97	0.64
1:A:177:PRO:HG2	1:A:208:LEU:HD23	1.80	0.64
1:A:280:LEU:HD11	1:A:436:GLU:HG3	1.80	0.63
2:B:82:U:H2'	2:B:83:G:C8	2.33	0.63
1:A:300:THR:O	1:A:304:THR:HG23	1.99	0.63
1:A:953:ARG:NH2	1:A:1079:PRO:O	2.29	0.63
1:A:290:GLU:CB	1:A:330:ALA:CB	2.76	0.62
1:A:396:THR:HG22	1:A:408:LEU:HD21	1.79	0.62
1:A:313:SER:HA	1:A:354:HIS:CE1	2.35	0.62
2:B:90:A:H2'	2:B:91:A:C8	2.35	0.62
1:A:829:PHE:HB2	1:A:1039:HIS:HB3	1.82	0.61
1:A:181:ALA:O	1:A:185:PHE:HB2	2.00	0.61
1:A:862:LYS:NZ	1:A:863:ARG:O	2.29	0.61
1:A:478:LEU:HD13	1:A:478:LEU:C	2.21	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ALA:HB2	1:A:727:ALA:N	2.15	0.61
1:A:464:LEU:HD12	1:A:465:PRO:HD2	1.83	0.61
1:A:174:PHE:CE2	1:A:180:LEU:HD12	2.35	0.61
1:A:486:LYS:HB3	1:A:1066:LEU:HD21	1.83	0.61
1:A:802:THR:HG23	1:A:805:LYS:H	1.66	0.61
1:A:17:ILE:HD11	1:A:484:ALA:HB2	1.82	0.60
1:A:830:VAL:HG22	1:A:1039:HIS:HA	1.83	0.60
1:A:74:ARG:NH1	2:B:21:U:OP2	2.34	0.60
3:C:26:DC:H2'	3:C:27:DA:H8	1.67	0.60
1:A:814:LEU:O	1:A:816:SER:N	2.35	0.59
1:A:780:ALA:O	1:A:783:VAL:HG12	2.02	0.59
1:A:464:LEU:HB2	1:A:685:ILE:HD12	1.85	0.58
1:A:403:VAL:HG22	1:A:404:GLN:H	1.67	0.58
1:A:706:LEU:O	1:A:708:GLY:N	2.37	0.58
1:A:79:LEU:HD11	1:A:136:LYS:HG3	1.86	0.58
1:A:210:ALA:O	1:A:213:ILE:HG13	2.04	0.57
1:A:159:LEU:HG	3:C:16:DA:H5"	1.86	0.57
1:A:1034:PHE:O	1:A:1050:ILE:HA	2.05	0.57
2:B:51:A:H2'	2:B:52:C:C6	2.38	0.56
1:A:836:ARG:HG2	1:A:1004:HIS:HB2	1.88	0.56
3:C:5:DT:H2"	3:C:6:DG:C8	2.40	0.56
2:B:60:A:H2'	2:B:61:C:C6	2.40	0.56
2:B:79:A:O2'	2:B:80:G:OP1	2.23	0.56
1:A:174:PHE:CD2	1:A:180:LEU:HD13	2.41	0.55
1:A:830:VAL:H	1:A:1039:HIS:HB2	1.72	0.55
1:A:953:ARG:NE	1:A:955:ASP:OD1	2.34	0.55
1:A:975:ALA:O	1:A:978:VAL:HG12	2.07	0.54
1:A:24:TRP:CZ2	1:A:1067:GLY:HA2	2.42	0.54
1:A:687:LEU:HD12	1:A:687:LEU:C	2.28	0.54
2:B:80:G:H2'	2:B:81:A:H8	1.72	0.54
1:A:309:PRO:HG3	1:A:315:LEU:HG	1.89	0.54
1:A:464:LEU:HB3	1:A:489:ASN:HD21	1.73	0.53
1:A:888:LEU:HD22	1:A:916:LYS:HE2	1.89	0.53
2:B:97:U:H2'	2:B:98:G:C8	2.43	0.53
1:A:14:GLY:O	1:A:24:TRP:HA	2.09	0.53
1:A:686:LEU:HD23	1:A:686:LEU:H	1.73	0.53
1:A:797:PHE:HD2	1:A:798:GLU:N	2.07	0.53
1:A:9:ILE:CG2	1:A:10:ASN:N	2.71	0.53
1:A:472:ILE:HD12	1:A:473:ARG:H	1.73	0.53
2:B:97:U:H2'	2:B:98:G:H8	1.74	0.53
1:A:332:PHE:O	1:A:335:LEU:HD23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:983:LEU:HD21	1:A:1078:ARG:NH2	2.24	0.53
1:A:783:VAL:O	1:A:787:VAL:HG23	2.09	0.52
1:A:802:THR:CG2	1:A:805:LYS:H	2.22	0.52
1:A:956:VAL:HG22	1:A:1001:PHE:H	1.73	0.52
1:A:811:ALA:O	1:A:815:SER:HA	2.09	0.52
1:A:199:TYR:OH	3:C:14:DA:N3	2.38	0.52
1:A:43:VAL:HG13	1:A:830:VAL:HG12	1.91	0.52
1:A:220:LYS:HG3	1:A:227:VAL:HG21	1.91	0.52
1:A:174:PHE:CD2	1:A:180:LEU:CD1	2.93	0.52
1:A:14:GLY:HA3	1:A:731:ALA:HB2	1.93	0.51
1:A:24:TRP:CH2	1:A:42:GLY:HA3	2.45	0.51
1:A:829:PHE:HB2	1:A:1039:HIS:CB	2.40	0.51
1:A:956:VAL:HG22	1:A:1001:PHE:O	2.11	0.51
1:A:983:LEU:HD21	1:A:1078:ARG:HH21	1.74	0.51
3:C:32:DG:H2'	3:C:33:DA:C8	2.46	0.51
2:B:130:U:H2'	2:B:131:U:C6	2.46	0.51
3:C:23:DT:H2'	3:C:24:DA:H8	1.74	0.51
2:B:99:C:H2'	2:B:100:C:C6	2.46	0.51
1:A:311:ARG:HG2	1:A:371:PRO:HB3	1.92	0.51
1:A:992:ARG:HB3	1:A:992:ARG:CZ	2.40	0.51
1:A:959:LYS:O	1:A:966:ASN:ND2	2.44	0.50
1:A:957:PHE:CD2	1:A:993:ILE:HG21	2.47	0.50
4:D:15:DT:H2"	4:D:16:DA:C8	2.46	0.50
1:A:16:ALA:O	1:A:22:VAL:HA	2.11	0.50
1:A:273:THR:OG1	1:A:380:ASP:OD2	2.27	0.50
3:C:22:DA:H2'	3:C:23:DT:C6	2.46	0.50
1:A:911:ASP:O	1:A:912:ASN:ND2	2.45	0.50
1:A:955:ASP:CB	1:A:1078:ARG:HH12	2.23	0.50
1:A:956:VAL:CG2	1:A:1001:PHE:H	2.24	0.50
3:C:26:DC:H2'	3:C:27:DA:C8	2.47	0.50
2:B:70:C:H2'	2:B:71:G:H8	1.77	0.49
1:A:40:ASP:OD2	1:A:495:TYR:OH	2.20	0.49
1:A:158:LEU:O	1:A:162:VAL:HG12	2.11	0.49
1:A:815:SER:O	1:A:816:SER:O	2.31	0.49
2:B:98:G:H2'	2:B:99:C:C6	2.47	0.49
1:A:790:LYS:HA	1:A:797:PHE:HB3	1.94	0.48
1:A:956:VAL:HG21	1:A:1000:CYS:HB2	1.96	0.48
1:A:780:ALA:O	1:A:782:GLU:N	2.46	0.48
1:A:77:HIS:O	1:A:81:ARG:HG3	2.13	0.48
2:B:82:U:H4'	2:B:83:G:OP1	2.12	0.48
1:A:677:LEU:O	1:A:681:VAL:HG12	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:985:ASP:OD1	1:A:985:ASP:N	2.45	0.48
2:B:9:U:H2'	2:B:10:G:H8	1.79	0.48
2:B:31:G:H2'	2:B:32:C:C6	2.48	0.48
1:A:224:ASN:HB3	1:A:227:VAL:HG13	1.95	0.47
2:B:93:G:H2'	2:B:94:C:C6	2.50	0.47
3:C:19:DT:H2'	3:C:20:DA:C8	2.50	0.47
1:A:356:ILE:HD13	1:A:382:ILE:HD11	1.96	0.47
1:A:481:LEU:HD22	1:A:677:LEU:HD13	1.96	0.47
1:A:314:LYS:HG3	1:A:346:THR:HG23	1.96	0.47
1:A:970:ILE:N	1:A:987:ASP:OD1	2.42	0.47
2:B:132:U:H4'	2:B:133:A:O5'	2.14	0.47
1:A:315:LEU:HD12	1:A:347:LEU:HD23	1.97	0.47
1:A:375:SER:O	1:A:378:LEU:N	2.48	0.47
2:B:16:U:H2'	2:B:17:A:C8	2.49	0.47
2:B:73:C:H2'	2:B:74:U:C6	2.50	0.47
2:B:79:A:H2'	2:B:80:G:H8	1.80	0.47
4:D:10:DC:H2"	4:D:11:DC:H5'	1.97	0.47
1:A:66:ARG:NH1	2:B:19:C:OP2	2.48	0.47
1:A:1011:PHE:HB3	1:A:1020:GLU:O	2.15	0.47
2:B:28:G:H2'	2:B:29:U:C6	2.49	0.47
1:A:699:ASN:O	1:A:701:GLN:N	2.47	0.47
1:A:175:ARG:NH2	1:A:179:GLU:OE2	2.45	0.46
4:D:11:DC:H2"	4:D:12:DA:C8	2.50	0.46
1:A:466:PRO:HA	1:A:485:ARG:CZ	2.45	0.46
1:A:205:ARG:NH1	1:A:242:GLN:OE1	2.47	0.46
2:B:79:A:HO2'	2:B:80:G:P	2.37	0.46
1:A:198:ASP:OD2	1:A:200:SER:OG	2.29	0.46
1:A:114:ARG:HG3	1:A:130:VAL:HG22	1.97	0.46
1:A:177:PRO:HD2	1:A:211:GLU:HB2	1.97	0.46
1:A:1044:LYS:HA	1:A:1048:PHE:CD1	2.50	0.46
1:A:895:ARG:HD3	1:A:906:ALA:O	2.15	0.46
2:B:79:A:H2'	2:B:80:G:C8	2.51	0.46
1:A:41:LEU:HD11	1:A:787:VAL:HA	1.97	0.46
1:A:465:PRO:O	1:A:485:ARG:HD2	2.16	0.46
1:A:798:GLU:H	1:A:798:GLU:CD	2.19	0.46
1:A:109:THR:N	1:A:110:PRO:HD3	2.30	0.45
1:A:176:THR:HG23	1:A:214:LEU:HD23	1.98	0.45
1:A:913:PRO:HB2	1:A:915:TYR:CE1	2.51	0.45
2:B:96:C:N3	2:B:128:C:O2'	2.47	0.45
1:A:26:MET:HE2	1:A:26:MET:HB2	1.83	0.45
1:A:378:LEU:O	1:A:382:ILE:HG23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:LEU:O	1:A:432:VAL:HG23	2.17	0.45
1:A:412:LEU:HD12	1:A:412:LEU:HA	1.82	0.45
1:A:805:LYS:O	1:A:809:LEU:HB2	2.17	0.45
1:A:871:LEU:HD23	1:A:871:LEU:O	2.16	0.45
1:A:206:LYS:H	1:A:206:LYS:HD2	1.82	0.45
1:A:182:LEU:O	1:A:186:GLU:HB2	2.17	0.45
1:A:936:LEU:O	1:A:939:LYS:NZ	2.45	0.45
1:A:1035:TYR:O	1:A:1036:LEU:HD23	2.16	0.45
1:A:145:LYS:HG2	3:C:18:DT:OP1	2.17	0.45
1:A:836:ARG:HG2	1:A:1004:HIS:CB	2.46	0.45
3:C:25:DG:H2'	3:C:26:DC:C6	2.51	0.45
1:A:107:PRO:HD2	1:A:129:ALA:HB2	1.99	0.44
2:B:99:C:H2'	2:B:100:C:H6	1.81	0.44
1:A:60:MET:HG3	1:A:63:ARG:NH1	2.31	0.44
2:B:92:C:H2'	2:B:93:G:C8	2.51	0.44
1:A:26:MET:HE1	1:A:495:TYR:CD1	2.53	0.44
1:A:813:LYS:HE3	1:A:813:LYS:HB3	1.80	0.44
1:A:258:CYS:HB3	1:A:261:GLU:O	2.18	0.44
1:A:926:ARG:HG3	2:B:56:U:OP1	2.18	0.44
2:B:37:U:H2'	2:B:38:U:O4'	2.18	0.44
1:A:464:LEU:CD2	1:A:488:ILE:HD11	2.47	0.44
1:A:787:VAL:HG12	1:A:788:PHE:CE1	2.53	0.44
1:A:951:MET:HE1	1:A:1027:CYS:HB3	1.99	0.44
1:A:703:THR:CB	1:A:729:VAL:HG11	2.48	0.43
1:A:966:ASN:ND2	1:A:967:GLN:N	2.66	0.43
1:A:1062:GLN:HA	1:A:1072:PRO:HA	1.99	0.43
2:B:92:C:H2'	2:B:93:G:H8	1.83	0.43
1:A:296:PRO:O	1:A:297:LEU:HD12	2.17	0.43
1:A:726:ASP:HA	1:A:729:VAL:HG12	2.00	0.43
1:A:846:LEU:O	2:B:27:U:H5'	2.18	0.43
1:A:968:TYR:HB2	1:A:1053:GLN:O	2.18	0.43
1:A:316:THR:OG1	1:A:319:GLN:HG3	2.19	0.43
1:A:816:SER:O	1:A:817:ARG:C	2.55	0.43
1:A:828:LEU:HD12	1:A:829:PHE:N	2.34	0.43
1:A:433:PRO:HA	1:A:436:GLU:OE2	2.19	0.43
1:A:913:PRO:HB2	1:A:915:TYR:HE1	1.81	0.43
1:A:9:ILE:CG2	1:A:10:ASN:H	2.30	0.43
1:A:24:TRP:CZ2	1:A:42:GLY:HA3	2.54	0.43
1:A:174:PHE:CE2	1:A:180:LEU:CD1	3.02	0.43
1:A:180:LEU:HD23	1:A:203:PHE:HE1	1.84	0.43
1:A:43:VAL:HG12	1:A:828:LEU:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:GLU:HA	1:A:330:ALA:HA	2.01	0.43
1:A:69:ARG:NH1	2:B:88:G:N7	2.67	0.43
1:A:485:ARG:HA	1:A:488:ILE:HG12	2.01	0.42
1:A:502:HIS:HA	1:A:696:PHE:O	2.19	0.42
1:A:846:LEU:O	2:B:26:U:O2'	2.32	0.42
1:A:863:ARG:HE	1:A:924:ALA:HB3	1.84	0.42
2:B:34:C:H2'	2:B:35:C:O4'	2.19	0.42
1:A:78:ARG:HG3	1:A:238:LEU:HD22	2.01	0.42
1:A:316:THR:HA	1:A:347:LEU:H	1.82	0.42
1:A:981:ASN:ND2	1:A:981:ASN:O	2.52	0.42
1:A:418:ASP:OD1	1:A:418:ASP:N	2.53	0.42
1:A:1008:LEU:HD12	1:A:1009:ILE:H	1.84	0.42
2:B:90:A:H2'	2:B:91:A:H8	1.82	0.42
1:A:464:LEU:HD23	1:A:488:ILE:HD11	2.01	0.42
1:A:1032:GLY:O	1:A:1053:GLN:NE2	2.50	0.42
1:A:36:ILE:HG12	1:A:37:ARG:H	1.85	0.42
1:A:367:ASP:OD2	1:A:367:ASP:N	2.53	0.42
1:A:878:VAL:HG23	1:A:926:ARG:O	2.18	0.42
2:B:128:C:H2'	2:B:129:G:H8	1.83	0.42
1:A:832:ARG:HB3	1:A:832:ARG:NH1	2.35	0.42
2:B:40:C:H2'	2:B:41:U:C6	2.55	0.42
1:A:30:ASP:OD1	1:A:31:GLU:N	2.53	0.41
2:B:97:U:O2'	2:B:98:G:OP1	2.29	0.41
2:B:98:G:H2'	2:B:99:C:H6	1.85	0.41
3:C:1:DT:H2"	3:C:2:DA:C8	2.55	0.41
1:A:9:ILE:HG22	1:A:10:ASN:H	1.85	0.41
4:D:5:DG:H2"	4:D:6:DA:C8	2.55	0.41
1:A:125:LEU:HD23	1:A:125:LEU:HA	1.95	0.41
1:A:220:LYS:HE3	1:A:220:LYS:HB3	1.92	0.41
1:A:836:ARG:HG2	1:A:1004:HIS:CG	2.54	0.41
1:A:960:VAL:O	1:A:962:LYS:N	2.44	0.41
3:C:23:DT:H2'	3:C:24:DA:C8	2.53	0.41
3:C:32:DG:H2'	3:C:33:DA:H8	1.85	0.41
1:A:86:LEU:HD13	1:A:128:SER:HB3	2.01	0.41
1:A:472:ILE:HD13	3:C:23:DT:H5'	2.03	0.41
1:A:874:LEU:HD12	1:A:874:LEU:HA	1.85	0.41
1:A:1063:VAL:HG22	1:A:1071:ARG:O	2.21	0.41
2:B:70:C:H2'	2:B:71:G:C8	2.53	0.41
1:A:27:VAL:HG12	1:A:38:LEU:HA	2.01	0.41
1:A:966:ASN:HD22	1:A:967:GLN:N	2.18	0.41
1:A:20:ALA:HA	1:A:46:PHE:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:LEU:HD23	1:A:323:LEU:C	2.41	0.41
4:D:6:DA:H2"	4:D:7:DG:C8	2.56	0.41
1:A:22:VAL:O	1:A:43:VAL:HG23	2.20	0.41
1:A:58:LEU:H	1:A:58:LEU:HD12	1.86	0.41
2:B:3:U:H2'	2:B:4:C:C6	2.56	0.41
1:A:70:ARG:O	1:A:74:ARG:HG3	2.20	0.40
1:A:1054:ASN:O	1:A:1055:LEU:HD13	2.21	0.40
1:A:1049:ARG:HE	1:A:1049:ARG:HB2	1.76	0.40
2:B:80:G:H2'	2:B:81:A:C8	2.54	0.40
1:A:1012:GLN:CB	1:A:1058:ILE:HG22	2.51	0.40
1:A:873:ASP:HB3	1:A:927:VAL:HG21	2.04	0.40
2:B:123:G:H2'	2:B:124:G:C8	2.56	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	840/1083 (78%)	758 (90%)	75 (9%)	7 (1%)	19 49

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	816	SER
1	A	327	GLU
1	A	700	GLY
1	A	815	SER
1	A	821	VAL
1	A	790	LYS
1	A	1050	ILE

5.3.2 Protein sidechains (i)

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	576/935 (62%)	547 (95%)	29 (5%)	24 54

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	SER
1	A	173	ASP
1	A	206	LYS
1	A	247	SER
1	A	322	LYS
1	A	333	LYS
1	A	336	ARG
1	A	358	ARG
1	A	367	ASP
1	A	474	ASN
1	A	677	LEU
1	A	686	LEU
1	A	698	SER
1	A	722	HIS
1	A	725	LEU
1	A	788	PHE
1	A	797	PHE
1	A	812	GLU
1	A	827	PRO
1	A	829	PHE
1	A	837	LYS
1	A	912	ASN
1	A	936	LEU
1	A	991	TYR
1	A	992	ARG
1	A	1011	PHE
1	A	1038	TRP
1	A	1049	ARG
1	A	1073	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	722	HIS
1	A	966	ASN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	115/135 (85%)	29 (25%)	6 (5%)

All (29) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	24	C
2	B	43	G
2	B	47	G
2	B	48	A
2	B	65	A
2	B	69	C
2	B	70	C
2	B	73	C
2	B	74	U
2	B	75	G
2	B	76	A
2	B	77	A
2	B	78	A
2	B	79	A
2	B	80	G
2	B	81	A
2	B	83	G
2	B	84	U
2	B	85	G
2	B	87	C
2	B	89	C
2	B	94	C
2	B	97	U
2	B	98	G
2	B	121	G
2	B	122	G
2	B	128	C
2	B	133	A
2	B	134	U

All (6) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	68	G
2	B	73	C
2	B	79	A
2	B	82	U
2	B	93	G
2	B	97	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	EDO	A	1101	-	3,3,3	0.47	0	2,2,2	0.35	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	1101	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (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 (i)

6.1 Protein, DNA and RNA chains (i)

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	858/1083 (79%)	0.11	42 (4%) 29 29	12, 70, 136, 166	0
2	B	117/135 (86%)	0.03	5 (4%) 35 34	12, 53, 161, 207	0
3	C	35/35 (100%)	-0.55	0 100 100	17, 39, 75, 79	0
4	D	13/16 (81%)	-0.36	0 100 100	32, 43, 74, 110	0
All	All	1023/1269 (80%)	0.08	47 (4%) 32 32	12, 65, 141, 207	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	730	VAL	6.0
1	A	731	ALA	5.8
1	A	704	ASN	5.6
1	A	38	LEU	4.5
1	A	691	GLY	4.4
1	A	957	PHE	4.1
2	B	102	C	4.0
1	A	365	LEU	3.8
1	A	320	ALA	3.6
1	A	732	CYS	3.4
1	A	698	SER	3.4
1	A	776	TRP	3.3
1	A	693	ARG	3.1
2	B	120	A	3.0
1	A	502	HIS	3.0
1	A	155	LEU	3.0
1	A	705	LEU	2.9
1	A	822	HIS	2.9
1	A	965	LYS	2.8
1	A	337	TYR	2.8
1	A	729	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	469	ALA	2.6
1	A	817	ARG	2.5
1	A	357	SER	2.5
1	A	815	SER	2.5
1	A	364	GLY	2.5
1	A	503	ILE	2.4
1	A	991	TYR	2.4
1	A	690	LYS	2.4
1	A	779	PHE	2.4
1	A	1017	SER	2.4
1	A	360	LEU	2.4
1	A	680	PHE	2.3
1	A	775	PRO	2.3
1	A	814	LEU	2.2
1	A	1022	ALA	2.2
1	A	726	ASP	2.2
1	A	1081	VAL	2.2
1	A	262	PRO	2.1
1	A	453	TYR	2.1
2	B	73	C	2.1
2	B	101	C	2.0
1	A	1006	TYR	2.0
1	A	1011	PHE	2.0
2	B	75	G	2.0
1	A	1045	GLU	2.0
1	A	491	VAL	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 monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	EDO	A	1101	4/4	0.95	0.12	33,40,54,56	0

6.5 Other polymers (i)

There are no such residues in this entry.