



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 06:40 pm BST

PDB ID : 5U31
Title : Crystal structure of AacC2c1-sgRNA-8mer substrate DNA ternary complex
Authors : Yang, H.; Gao, P.; Rajashankar, K.R.; Patel, D.J.
Deposited on : 2016-12-01
Resolution : 2.89 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

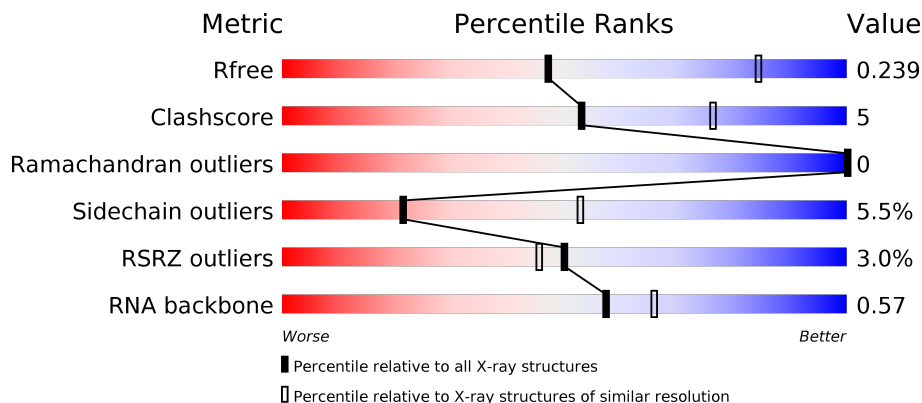
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)
RNA backbone	3102	1007 (3.16-2.64)

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 on 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	1130	 3% 80% 15% . .
2	B	112	 2% 49% 27% 13% 12%
3	C	28	 75% 25%
4	D	8	 50% 50%

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Mol	Chain	Length	Quality of chain
4	E	8	 50% 38% 13%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11871 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 C2c1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	S				Se
1	A	1085	8864	5569	1653	1611	10	21	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP T0D7A2
A	570	ALA	ASP	engineered mutation	UNP T0D7A2
A	848	ALA	GLU	engineered mutation	UNP T0D7A2
A	977	ALA	ASP	engineered mutation	UNP T0D7A2

- Molecule 2 is a RNA chain called sgRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	99	2127	949	396	683	99	0	0	0

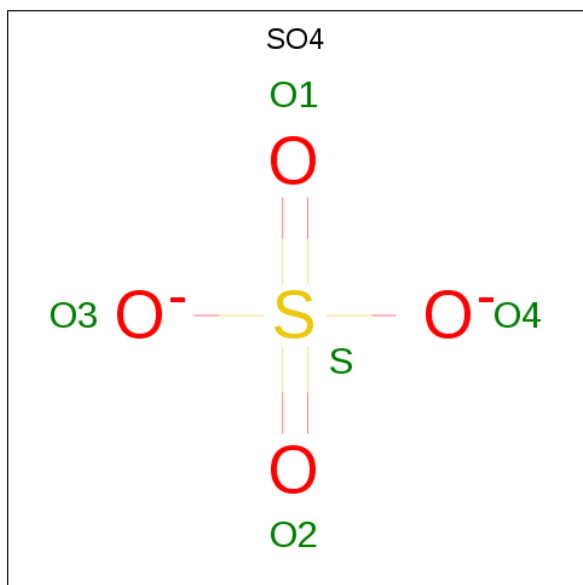
- Molecule 3 is a DNA chain called Target DNA strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	C	28	565	271	98	169	27	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	8	162	79	26	50	7	0	0	0
4	E	7	143	70	23	44	6	0	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3: Target DNA strand

Chain C:  75% 25%



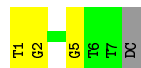
- Molecule 4: Non-target DNA strand

Chain D:  50% 50%



- Molecule 4: Non-target DNA strand

Chain E:  50% 38% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	117.25Å 182.15Å 216.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.30 – 2.89 139.46 – 2.89	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.30-2.89) 100.0 (139.46-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.198 , 0.239 0.199 , 0.239	Depositor DCC
R_{free} test set	2556 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtrriage
Anisotropy	0.338	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11871	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.29	1/9036 (0.0%)	0.43	0/12139
2	B	0.28	0/2380	0.82	1/3706 (0.0%)
3	C	0.59	0/631	0.94	0/971
4	D	0.62	0/180	1.02	0/277
4	E	0.63	0/159	1.05	0/245
All	All	0.32	1/12386 (0.0%)	0.60	1/17338 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	58	GLU	CD-OE2	6.98	1.33	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	20	U	P-O3'-C3'	-5.05	113.64	119.70

There are no chirality outliers.

There are no planarity outliers.

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	8864	0	8763	102	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2127	0	1075	29	0
3	C	565	0	318	5	0
4	D	162	0	94	3	0
4	E	143	0	83	2	0
5	A	10	0	0	1	0
All	All	11871	0	10333	117	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 5.

All (117) 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:754:GLY:HA3	1:A:760:GLN:HG3	1.61	0.82
1:A:742:ARG:NH2	2:B:35:A:OP1	2.16	0.79
1:A:130:ASP:HB3	1:A:133:ALA:HB2	1.63	0.78
1:A:1085:ILE:HD13	1:A:1100:VAL:HG22	1.64	0.77
1:A:991:ASP:O	1:A:1020:ARG:NH1	2.21	0.73
1:A:397:LEU:HD12	1:A:458:PRO:HG2	1.72	0.72
1:A:15:MSE:HE1	1:A:440:PRO:HD2	1.71	0.72
1:A:332:ARG:NH2	4:E:5:DG:O6	2.26	0.69
1:A:678:VAL:HG21	1:A:693:GLU:HG2	1.75	0.69
1:A:905:THR:HG22	1:A:907:ALA:H	1.61	0.66
1:A:745:ILE:HD11	2:B:34:A:OP1	1.97	0.65
1:A:397:LEU:HG	1:A:405:THR:HG22	1.80	0.63
1:A:218:ARG:NH1	4:D:5:DG:OP2	2.26	0.63
1:A:614:HIS:ND1	2:B:6:G:OP1	2.32	0.63
3:C:26:DA:H2''	3:C:27:DC:H5''	1.81	0.63
1:A:15:MSE:HE1	1:A:439:VAL:HA	1.81	0.62
1:A:569:VAL:HB	1:A:847:LEU:HD23	1.81	0.61
1:A:260:GLN:HB3	1:A:263:LEU:HD13	1.84	0.60
1:A:824:VAL:HB	1:A:838:LYS:HD3	1.82	0.59
1:A:766:ARG:NH2	5:A:1202:SO4:O4	2.34	0.59
1:A:893:MSE:HE3	1:A:988:LEU:HA	1.85	0.57
1:A:507:ARG:NH2	3:C:21:DG:OP1	2.37	0.57
1:A:734:ARG:HD2	1:A:771:LEU:HD21	1.86	0.57
1:A:405:THR:HG23	1:A:421:HIS:NE2	2.20	0.57
1:A:50:ARG:NH2	1:A:60:GLU:OE1	2.37	0.57
1:A:994:ILE:HD12	1:A:998:ARG:HB3	1.86	0.57
2:B:85:G:N1	2:B:86:U:O4	2.38	0.57
1:A:616:ARG:NH1	1:A:819:GLU:OE1	2.37	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:ARG:NH1	1:A:177:ASP:OD1	2.37	0.56
1:A:607:ASN:N	1:A:607:ASN:OD1	2.39	0.56
1:A:245:LEU:HB3	1:A:371:LEU:HD13	1.88	0.55
1:A:918:ARG:HG3	2:B:13:G:OP1	2.07	0.55
1:A:54:ASN:HA	1:A:789:LYS:HE2	1.89	0.54
1:A:238:VAL:HG22	2:B:98:C:H4'	1.89	0.54
1:A:588:ASP:OD1	1:A:588:ASP:N	2.41	0.54
1:A:765:GLU:OE1	1:A:798:ARG:NH2	2.39	0.54
1:A:15:MSE:CE	1:A:440:PRO:HD2	2.39	0.53
2:B:69:U:O2'	2:B:70:U:O5'	2.24	0.53
1:A:746:ARG:NH2	2:B:60:C:OP1	2.41	0.53
1:A:1000:ARG:HG2	1:A:1076:VAL:HG22	1.91	0.52
1:A:33:VAL:HG12	1:A:383:PRO:HG3	1.91	0.52
1:A:914:ARG:HE	1:A:959:ILE:HD13	1.75	0.52
1:A:400:ASN:ND2	3:C:23:DA:H62	2.07	0.52
1:A:745:ILE:HD12	2:B:35:A:H62	1.75	0.51
1:A:108:PRO:HD2	1:A:229:MSE:HE2	1.91	0.51
1:A:745:ILE:HD12	1:A:745:ILE:H	1.75	0.51
4:E:1:DT:H2'	4:E:2:DG:N7	2.24	0.51
1:A:464:ARG:HG2	1:A:471:HIS:CE1	2.46	0.51
1:A:453:ARG:NH2	1:A:474:GLY:O	2.44	0.51
1:A:188:LYS:HB3	1:A:189:PRO:HD3	1.93	0.51
1:A:738:ARG:HB3	2:B:37:G:H5''	1.94	0.50
1:A:305:LYS:HA	1:A:308:LYS:HE2	1.93	0.50
1:A:847:LEU:HD11	1:A:879:LEU:HD12	1.94	0.50
1:A:83:ASN:ND2	1:A:189:PRO:HD2	2.27	0.50
1:A:444:SER:HA	2:B:29:G:C8	2.47	0.49
1:A:22:LEU:HD22	1:A:481:ILE:HD13	1.94	0.49
1:A:270:LEU:HD21	1:A:341:LYS:HG2	1.93	0.49
1:A:109:GLN:OE1	1:A:109:GLN:N	2.45	0.49
1:A:123:LYS:HE2	1:A:175:THR:HG21	1.95	0.49
1:A:4:LYS:HB3	2:B:92:C:H5'	1.95	0.48
1:A:821:LEU:HG	1:A:888:LEU:HD11	1.95	0.48
2:B:68:C:H42	2:B:85:G:H1	1.61	0.48
1:A:865:ASN:HA	1:A:868:MSE:HE3	1.94	0.48
1:A:868:MSE:HE1	2:B:102:G:C4'	2.44	0.48
2:B:21:C:C2'	2:B:22:A:H5'	2.44	0.47
1:A:305:LYS:HB3	1:A:321:GLU:CG	2.43	0.47
1:A:69:LYS:HD3	1:A:96:LEU:HD13	1.95	0.47
1:A:1030:PHE:CE1	1:A:1040:GLU:HG2	2.49	0.47
1:A:868:MSE:HE1	2:B:102:G:H4'	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:TRP:HZ2	1:A:322:ILE:HD11	1.80	0.47
1:A:629:LYS:NZ	2:B:15:A:H62	2.13	0.47
1:A:418:ILE:HD12	1:A:441:ILE:HD11	1.97	0.47
4:D:8:DC:H6	4:D:8:DC:H5'	1.80	0.46
1:A:1009:GLU:HG3	1:A:1011:VAL:HG13	1.98	0.46
2:B:21:C:H2'	2:B:22:A:H5'	1.96	0.46
2:B:106:A:H2'	2:B:107:A:C8	2.51	0.46
1:A:252:PHE:CE1	1:A:364:TYR:HB2	2.51	0.46
1:A:635:ARG:HA	1:A:638:ARG:HD2	1.97	0.46
2:B:68:C:N4	2:B:69:U:O4	2.49	0.46
1:A:768:TYR:CE1	1:A:798:ARG:HG2	2.51	0.45
1:A:418:ILE:HD13	1:A:502:LEU:HD23	1.98	0.45
1:A:344:GLU:O	1:A:348:GLN:HG3	2.17	0.45
1:A:828:ASP:C	1:A:830:ARG:H	2.20	0.45
1:A:226:GLU:HG2	3:C:19:DT:H5'	1.99	0.45
1:A:12:LEU:HD23	1:A:15:MSE:HE3	1.98	0.45
1:A:415:ARG:NH2	2:B:27:G:OP1	2.50	0.45
1:A:904:ARG:NH2	2:B:5:A:OP2	2.50	0.44
1:A:1028:LYS:NZ	1:A:1040:GLU:OE1	2.44	0.44
1:A:8:VAL:HG13	1:A:441:ILE:HG23	1.99	0.44
1:A:29:VAL:HA	1:A:220:MSE:HE2	1.98	0.44
1:A:165:LYS:HA	1:A:168:THR:HG22	1.99	0.43
1:A:22:LEU:HD13	1:A:500:VAL:HG11	2.00	0.43
1:A:800:HIS:ND1	2:B:32:C:OP1	2.44	0.43
2:B:56:A:H2'	2:B:57:A:O4'	2.19	0.43
1:A:832:LYS:NZ	2:B:85:G:H5''	2.34	0.42
1:A:746:ARG:HH22	2:B:60:C:P	2.42	0.42
1:A:397:LEU:HD23	1:A:397:LEU:HA	1.84	0.42
1:A:679:ASP:HB3	1:A:689:ARG:HH11	1.85	0.42
1:A:82:GLU:OE1	1:A:192:ARG:NH2	2.53	0.42
2:B:94:G:H2'	2:B:95:A:C8	2.55	0.42
1:A:167:GLU:H	1:A:167:GLU:HG3	1.74	0.41
1:A:624:GLY:HA3	1:A:754:GLY:O	2.19	0.41
1:A:1029:VAL:HG12	1:A:1041:ARG:HB2	2.01	0.41
1:A:459:ILE:HA	1:A:459:ILE:HD13	1.74	0.41
1:A:490:HIS:HB3	1:A:492:HIS:CD2	2.55	0.41
1:A:638:ARG:HH22	1:A:760:GLN:HG2	1.86	0.41
1:A:819:GLU:OE2	2:B:8:G:N1	2.43	0.41
2:B:87:G:O2'	2:B:88:G:H5''	2.20	0.41
1:A:780:VAL:O	1:A:783:GLN:HG2	2.20	0.41
3:C:24:DC:H2''	3:C:25:DC:H5'	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:3:DT:H2"	4:D:4:DG:C8	2.56	0.41
1:A:34:ARG:NH1	1:A:383:PRO:O	2.50	0.41
1:A:879:LEU:HD23	1:A:879:LEU:HA	1.89	0.41
1:A:319:ASP:HA	1:A:322:ILE:HD12	2.04	0.40
1:A:487:GLN:O	1:A:498:ARG:NH1	2.54	0.40
1:A:568:SER:O	1:A:579:ILE:HA	2.22	0.40
1:A:522:TYR:HB3	1:A:545:TYR:OH	2.21	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	1077/1130 (95%)	1025 (95%)	52 (5%)	0	100 100

There are no Ramachandran outliers to report.

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	930/945 (98%)	879 (94%)	51 (6%)	21 53

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	72	LEU
1	A	98	LEU
1	A	134	VAL
1	A	190	LEU
1	A	251	ARG
1	A	253	GLU
1	A	257	PHE
1	A	270	LEU
1	A	309	LEU
1	A	315	PHE
1	A	338	LEU
1	A	392	THR
1	A	397	LEU
1	A	436	ASP
1	A	459	ILE
1	A	480	LYS
1	A	490	HIS
1	A	491	MSE
1	A	517	GLU
1	A	588	ASP
1	A	590	LEU
1	A	595	LYS
1	A	597	ARG
1	A	598	VAL
1	A	607	ASN
1	A	629	LYS
1	A	649	LEU
1	A	652	LEU
1	A	678	VAL
1	A	711	TRP
1	A	722	VAL
1	A	738	ARG
1	A	786	ARG
1	A	798	ARG
1	A	812	LEU
1	A	855	PHE
1	A	890	VAL
1	A	893	MSE
1	A	932	LEU
1	A	950	ASP
1	A	956	GLU
1	A	988	LEU

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Mol	Chain	Res	Type
1	A	994	ILE
1	A	997	ILE
1	A	1000	ARG
1	A	1007	ASP
1	A	1029	VAL
1	A	1032	THR
1	A	1079	ARG
1	A	1092	ARG

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

Mol	Chain	Res	Type
1	A	400	ASN
1	A	490	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	97/112 (86%)	25 (25%)	4 (4%)

All (25) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	5	A
2	B	6	G
2	B	12	A
2	B	13	G
2	B	16	U
2	B	21	C
2	B	22	A
2	B	29	G
2	B	33	C
2	B	34	A
2	B	36	U
2	B	46	C
2	B	58	G
2	B	60	C
2	B	66	A
2	B	67	G
2	B	68	C

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Mol	Chain	Res	Type
2	B	69	U
2	B	70	U
2	B	83	A
2	B	86	U
2	B	87	G
2	B	93	A
2	B	101	A
2	B	110	U

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	5	A
2	B	32	C
2	B	82	G
2	B	86	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are 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	SO4	A	1201	-	4,4,4	0.14	0	6,6,6	0.15	0
5	SO4	A	1202	-	4,4,4	0.13	0	6,6,6	0.15	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1202	SO4	1	0

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	1064/1130 (94%)	0.41	34 (3%) 47 43	28, 51, 108, 153	0
2	B	99/112 (88%)	0.29	2 (2%) 65 63	30, 58, 156, 174	0
3	C	28/28 (100%)	-0.05	0 100 100	32, 40, 57, 69	0
4	D	8/8 (100%)	-0.01	0 100 100	32, 34, 39, 55	0
4	E	7/8 (87%)	0.67	0 100 100	44, 56, 97, 99	0
All	All	1206/1286 (93%)	0.39	36 (2%) 50 45	28, 51, 115, 174	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	495	ARG	7.3
1	A	1071	ARG	6.6
1	A	831	GLY	5.7
1	A	54	ASN	5.6
1	A	160	GLU	5.5
1	A	1072	GLU	4.7
1	A	830	ARG	4.2
1	A	161	GLU	4.2
1	A	608	ASP	3.8
1	A	494	ARG	3.8
1	A	164	GLU	3.5
1	A	492	HIS	3.1
1	A	332	ARG	3.1
1	A	490	HIS	3.0
1	A	0	SER	3.0
1	A	493	ARG	2.9
1	A	750	LYS	2.8
1	A	832	LYS	2.8
1	A	1044	GLY	2.8
2	B	67	G	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	168	THR	2.5
1	A	923	HIS	2.5
1	A	829	GLU	2.5
1	A	58	GLU	2.5
1	A	162	GLU	2.4
1	A	827	LEU	2.3
1	A	166	ALA	2.2
1	A	845	ILE	2.1
1	A	739	SER	2.1
1	A	291	VAL	2.1
1	A	463	PHE	2.1
1	A	317	LEU	2.0
1	A	328	ARG	2.0
1	A	908	PRO	2.0
2	B	66	A	2.0
1	A	789	LYS	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 carbohydrates 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(\AA^2)	Q<0.9
5	SO4	A	1202	5/5	0.78	0.30	95,111,117,121	0
5	SO4	A	1201	5/5	0.98	0.17	42,51,55,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.