



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 28, 2016 – 11:08 AM EDT

PDB ID : 5B2R
Title : Crystal structure of the Streptococcus pyogenes Cas9 VQR variant in complex with sgRNA and target DNA (TGA PAM)
Authors : Hirano, S.; Nishimasu, H.; Ishitani, R.; Nureki, O.
Deposited on : 2016-02-02
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

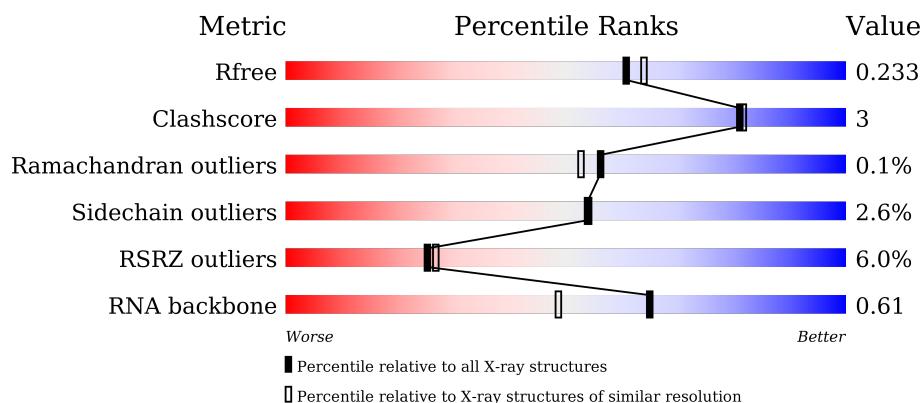
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)
RNA backbone	2183	1002 (2.72-1.30)

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. 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	81	<div> <div></div> <div> <div></div> <div>80%</div> <div>15%</div> <div>5%</div> </div> </div>
2	B	1372	<div> <div>6%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>...</div> </div> </div>
3	C	28	<div> <div></div> <div> <div></div> <div>71%</div> <div>29%</div> </div> </div>
4	D	8	<div> <div>13%</div> <div> <div></div> <div>75%</div> <div>25%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	K	A	101	-	-	-	X
5	K	A	108	-	-	-	X
5	K	B	1407	-	-	-	X
8	ACT	A	107	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 13799 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 RNA chain called Guide RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	81	Total	C	N	O	P	0	0	0
			1739	778	319	561	81			

- Molecule 2 is a protein called CRISPR-associated endonuclease Cas9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1319	Total	C	N	O	S	0	1	0
			10547	6734	1826	1967	20			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP Q99ZW2
B	-2	SER	-	expression tag	UNP Q99ZW2
B	-1	GLY	-	expression tag	UNP Q99ZW2
B	0	HIS	-	expression tag	UNP Q99ZW2
B	10	ALA	ASP	engineered mutation	UNP Q99ZW2
B	80	LEU	CYS	engineered mutation	UNP Q99ZW2
B	574	GLU	CYS	engineered mutation	UNP Q99ZW2
B	840	ALA	HIS	engineered mutation	UNP Q99ZW2
B	1135	VAL	ASP	engineered mutation	UNP Q99ZW2
B	1335	GLN	ARG	engineered mutation	UNP Q99ZW2
B	1337	ARG	THR	engineered mutation	UNP Q99ZW2

- Molecule 3 is a DNA chain called Target DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	28	Total	C	N	O	P	0	0	0
			560	269	100	164	27			

- Molecule 4 is a DNA chain called Non-target DNA, DNA (5'-D(*TP*GP*AP*GP*AP*TP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	8	Total	C	N	O	P	0	0	0
			165	80	31	47	7			

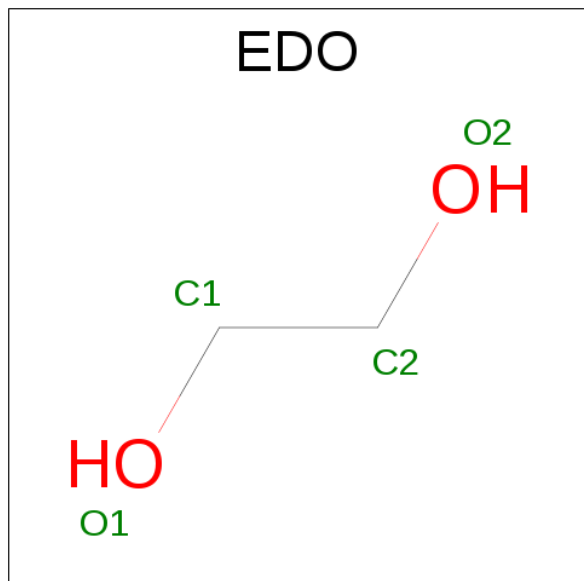
- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	7	Total	K	0	0
			7	7		
5	A	4	Total	K	0	0
			4	4		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Mg	0	0
			1	1		
6	A	2	Total	Mg	0	0
			2	2		

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



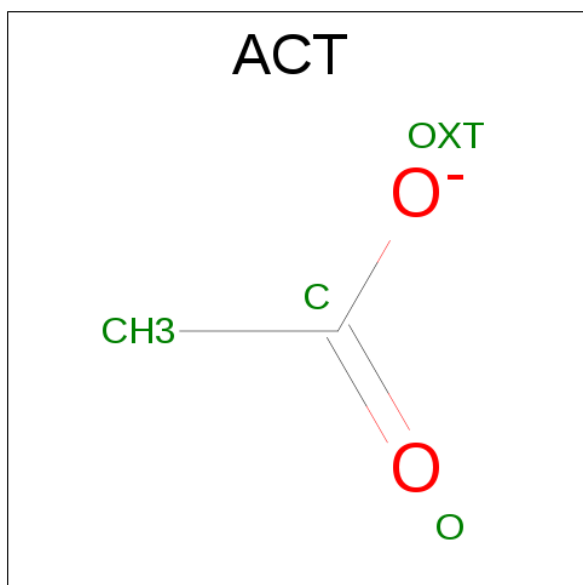
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



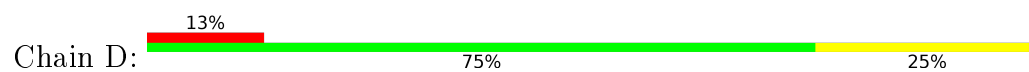
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	199	Total	O	0	0
			199	199		
9	B	492	Total	O	0	0
			492	492		
9	C	50	Total	O	0	0
			50	50		
9	D	13	Total	O	0	0
			13	13		



- Molecule 4: Non-target DNA, DNA (5'-D(*TP*GP*AP*GP*AP*TP*TP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.76Å 67.74Å 187.59Å 90.00° 111.23° 90.00°	Depositor
Resolution (Å)	47.94 – 2.00 47.94 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.2 (47.94-2.00) 97.3 (47.94-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.00Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.201 , 0.230 0.205 , 0.233	Depositor DCC
R_{free} test set	12952 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.586	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 137046 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13799	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, K, EDO, ACT

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	0.43	0/1949	0.90	0/3037
2	B	0.31	0/10738	0.47	0/14474
3	C	0.75	0/626	0.99	0/961
4	D	0.66	0/185	1.03	1/285 (0.4%)
All	All	0.37	0/13498	0.61	1/18757 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	7	DA	O4'-C1'-N9	5.54	111.88	108.00

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	1739	0	870	5	0
2	B	10547	0	10509	69	0
3	C	560	0	316	7	0
4	D	165	0	93	1	0
5	A	4	0	0	0	0
5	B	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	2	0	0	0	0
6	B	1	0	0	0	0
7	A	4	0	6	0	0
7	B	12	0	18	0	0
8	A	4	0	3	0	0
9	A	199	0	0	0	0
9	B	492	0	0	6	0
9	C	50	0	0	2	0
9	D	13	0	0	0	0
All	All	13799	0	11815	79	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 3.

All (79) 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:C:2:DA:H2'	3:C:3:DA:C8	2.19	0.78
3:C:19:DC:OP2	9:C:101:HOH:O	2.07	0.72
2:B:586:ARG:NH1	9:B:1501:HOH:O	2.26	0.67
2:B:558:LYS:HD2	2:B:586:ARG:HH21	1.62	0.64
2:B:241:LEU:O	2:B:244:LEU:N	2.24	0.64
2:B:222:LEU:HD23	2:B:234:LYS:HE3	1.83	0.61
2:B:253:LYS:NZ	2:B:261:ASP:OD1	2.30	0.59
1:A:71:U:H2'	1:A:72:U:C6	2.37	0.58
2:B:790:GLU:HG2	2:B:889:ALA:HA	1.85	0.57
2:B:427:GLU:OE1	2:B:437[B]:ARG:NH1	2.37	0.57
2:B:842:VAL:HB	2:B:847:LEU:HD22	1.87	0.56
2:B:784:ILE:HD12	2:B:806:LEU:HD13	1.88	0.56
2:B:302:LEU:HD12	2:B:414:ILE:HD11	1.88	0.56
2:B:468:LYS:NZ	9:B:1504:HOH:O	2.32	0.56
3:C:19:DC:H5''	3:C:19:DC:H6	1.71	0.55
2:B:787:GLY:HA3	2:B:891:LEU:HD21	1.88	0.55
2:B:1267:ASP:OD1	2:B:1294:TYR:OH	2.22	0.54
2:B:1179:ILE:HD11	2:B:1192:LYS:HG3	1.90	0.53
2:B:1256:GLN:O	2:B:1260:GLU:HG2	2.08	0.53
2:B:1303:ARG:NH2	2:B:1307:GLU:OE2	2.42	0.53
2:B:999:LYS:HB3	2:B:1073:VAL:HG22	1.91	0.52
2:B:243:ALA:HB1	2:B:248:LEU:HB3	1.92	0.51
2:B:338:LEU:HB3	2:B:383:MET:HE1	1.92	0.51
2:B:1114:ARG:NH1	4:D:9:DA:OP1	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:302:LEU:H	2:B:302:LEU:HD22	1.76	0.51
2:B:826:GLN:NE2	2:B:859:ARG:HD2	2.25	0.51
2:B:776:ASN:N	2:B:776:ASN:OD1	2.43	0.50
2:B:809:GLU:OE2	2:B:855:LYS:NZ	2.32	0.50
2:B:411:PRO:HD2	2:B:414:ILE:HG13	1.94	0.50
2:B:627:GLU:HG3	9:C:102:HOH:O	2.11	0.49
2:B:524:LEU:HD13	2:B:545:LYS:HG2	1.93	0.49
2:B:822:MET:HG3	2:B:856:VAL:HB	1.93	0.49
2:B:249:THR:HG22	2:B:263:LYS:HB2	1.93	0.49
2:B:1339:THR:O	2:B:1342:VAL:HG22	2.14	0.48
2:B:846:PHE:O	2:B:920:GLN:NE2	2.44	0.48
2:B:195:LEU:HD21	2:B:286:TYR:HA	1.94	0.48
3:C:1:DC:H2'	3:C:2:DA:C8	2.48	0.48
2:B:1177:ASN:ND2	2:B:1180:ASP:OD2	2.48	0.47
2:B:901:THR:O	2:B:904:GLU:HG2	2.16	0.46
2:B:336:LYS:NZ	9:B:1511:HOH:O	2.41	0.46
2:B:1215:ALA:HB2	2:B:1221:GLN:HG3	1.98	0.46
2:B:185:PHE:HD2	2:B:296:LEU:HD11	1.80	0.46
2:B:253:LYS:HG3	2:B:261:ASP:HA	1.97	0.46
2:B:302:LEU:HG	2:B:306:LEU:HD12	1.97	0.46
2:B:817:GLN:CD	2:B:822:MET:HG2	2.36	0.45
2:B:972:PHE:HE1	2:B:1084:ARG:HG2	1.81	0.45
2:B:1085:LYS:O	2:B:1089:MET:HG3	2.17	0.45
1:A:14:G:OP2	2:B:63:ARG:NH1	2.39	0.45
2:B:756:PRO:O	2:B:953:VAL:HG22	2.17	0.45
2:B:1004:LEU:HD11	2:B:1042:ILE:HD11	1.98	0.45
2:B:387:GLU:O	2:B:391:VAL:HG23	2.16	0.45
1:A:45:U:H5'	2:B:402:GLN:HE21	1.82	0.44
2:B:788:ILE:HG23	2:B:793:SER:HB3	2.00	0.44
2:B:511:HIS:HD2	9:B:1610:HOH:O	2.00	0.44
2:B:1194:LEU:HD13	2:B:1365:LEU:HD22	2.00	0.44
2:B:1340:LYS:NZ	9:B:1532:HOH:O	2.51	0.44
2:B:248:LEU:HD12	2:B:249:THR:H	1.83	0.43
2:B:288:ASP:HA	2:B:291:LEU:HB3	2.00	0.43
2:B:870:VAL:HG11	2:B:902:LYS:HB3	2.01	0.43
1:A:46:A:H2'	1:A:47:A:C8	2.54	0.43
2:B:1041:ASN:OD1	2:B:1044:ASN:ND2	2.49	0.43
2:B:187:GLN:O	2:B:191:THR:N	2.50	0.43
2:B:390:LEU:HD23	2:B:390:LEU:HA	1.87	0.43
2:B:1231:LYS:HE3	2:B:1232:TYR:CZ	2.54	0.43
2:B:841:ILE:HD13	2:B:900:LEU:HG	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:A:H3'	1:A:75:A:H8	1.84	0.42
2:B:195:LEU:HD23	2:B:289:LEU:HB2	2.01	0.42
2:B:302:LEU:H	2:B:302:LEU:CD2	2.31	0.42
2:B:1063:ILE:HG23	2:B:1072:ILE:HG13	2.02	0.42
2:B:1119:LEU:HB3	2:B:1128:PRO:HB3	2.02	0.42
3:C:20:DC:H2''	3:C:21:DT:H5'	2.02	0.42
3:C:16:DG:H2'	3:C:17:DC:C6	2.54	0.42
2:B:381:GLU:HG2	2:B:390:LEU:HD11	2.02	0.41
2:B:15:SER:HA	2:B:51:LEU:O	2.21	0.41
2:B:962:LEU:HD11	2:B:1043:MET:CE	2.49	0.41
2:B:302:LEU:HG	2:B:306:LEU:CD1	2.50	0.41
3:C:2:DA:H2''	3:C:3:DA:O5'	2.20	0.41
2:B:812:TYR:CZ	2:B:816:LEU:HD11	2.55	0.41
2:B:104:SER:HA	9:B:1774:HOH:O	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
2	B	1308/1372 (95%)	1269 (97%)	38 (3%)	1 (0%)	56 53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	1216	SER

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
2	B	1121/1227 (91%)	1092 (97%)	29 (3%)	54 54

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

Mol	Chain	Res	Type
2	B	101	LEU
2	B	123	VAL
2	B	219	SER
2	B	253	LYS
2	B	263	LYS
2	B	264	LEU
2	B	301	LEU
2	B	302	LEU
2	B	313	THR
2	B	314	LYS
2	B	419	LEU
2	B	476	TRP
2	B	490	SER
2	B	512	SER
2	B	524	LEU
2	B	571	LYS
2	B	638	THR
2	B	642	LEU
2	B	718	ASP
2	B	776	ASN
2	B	801	VAL
2	B	833	LEU
2	B	847	LEU
2	B	853	ASP
2	B	870	VAL
2	B	947	ASP
2	B	1119	LEU
2	B	1154	SER
2	B	1250	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	402	GLN
2	B	412	HIS
2	B	511	HIS
2	B	726	ASN
2	B	926	GLN
2	B	1041	ASN
2	B	1044	ASN
2	B	1261	GLN
2	B	1295	ASN
2	B	1297	HIS
2	B	1317	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	80/81 (98%)	12 (15%)	1 (1%)

All (12) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	17	U
1	A	28	A
1	A	40	C
1	A	51	A
1	A	56	U
1	A	58	G
1	A	59	U
1	A	68	A
1	A	71	U
1	A	72	U
1	A	74	A
1	A	75	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	71	U

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 19 ligands modelled in this entry, 14 are monoatomic - leaving 5 for Mogul analysis.

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	$\# Z > 2$
7	EDO	A	106	-	3,3,3	0.46	0	2,2,2	0.30	0
8	ACT	A	107	-	0,3,3	0.00	-	0,3,3	0.00	-
7	EDO	B	1409	-	3,3,3	0.55	0	2,2,2	0.21	0
7	EDO	B	1410	-	3,3,3	0.52	0	2,2,2	0.22	0
7	EDO	B	1411	-	3,3,3	0.44	0	2,2,2	0.40	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
7	EDO	A	106	-	-	0/1/1/1	0/0/0/0
8	ACT	A	107	-	-	0/0/0/0	0/0/0/0
7	EDO	B	1409	-	-	0/1/1/1	0/0/0/0
7	EDO	B	1410	-	-	0/1/1/1	0/0/0/0
7	EDO	B	1411	-	-	0/1/1/1	0/0/0/0

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.

No monomer is involved in short contacts.

5.7 Other polymers

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	81/81 (100%)	-0.59	1 (1%) 81 81	26, 41, 122, 188	0
2	B	1319/1372 (96%)	0.20	84 (6%) 23 24	24, 52, 95, 123	0
3	C	28/28 (100%)	-0.29	0 100 100	34, 48, 74, 84	0
4	D	8/8 (100%)	0.11	1 (12%) 5 5	37, 52, 98, 119	0
All	All	1436/1489 (96%)	0.15	86 (5%) 25 27	24, 51, 96, 188	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	186	ILE	8.0
2	B	264	LEU	6.4
2	B	247	GLY	6.0
2	B	292	ALA	5.4
2	B	244	LEU	5.1
2	B	190	GLN	5.0
2	B	262	ALA	4.8
2	B	296	LEU	4.8
2	B	245	SER	4.7
2	B	1033	THR	4.6
2	B	194	GLN	4.4
2	B	275	LEU	4.4
2	B	911	LEU	4.4
2	B	286	TYR	4.4
2	B	1257	LEU	4.3
2	B	1039	TYR	4.1
2	B	189	VAL	4.1
2	B	193	ASN	4.0
2	B	1036	TYR	3.8
2	B	287	ALA	3.8
2	B	804	THR	3.7

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Mol	Chain	Res	Type	RSRZ
2	B	200	PRO	3.7
2	B	266	LEU	3.6
2	B	1072	ILE	3.6
2	B	1052	LEU	3.6
2	B	216	LEU	3.6
2	B	178	ASN	3.5
2	B	1054	ASN	3.5
2	B	847	LEU	3.4
2	B	295	ASN	3.3
2	B	271	TYR	3.3
2	B	299	ALA	3.2
1	A	74	A	3.1
2	B	184	LEU	3.1
2	B	192	TYR	3.1
2	B	276	ASP	3.0
2	B	267	SER	3.0
2	B	917	ILE	3.0
2	B	1062	LEU	3.0
2	B	1034	ALA	3.0
2	B	269	ASP	2.9
2	B	201	ILE	2.9
2	B	214	ALA	2.9
2	B	795	ILE	2.9
2	B	195	LEU	2.9
2	B	1032	ALA	2.8
2	B	270	THR	2.8
2	B	279	LEU	2.7
2	B	301	LEU	2.7
2	B	191	THR	2.7
2	B	310	THR	2.7
2	B	187	GLN	2.7
2	B	249	THR	2.6
2	B	222	LEU	2.6
2	B	308	VAL	2.6
2	B	280	ALA	2.5
2	B	211	ILE	2.5
2	B	1251	ASP	2.5
2	B	1261	GLN	2.5
2	B	220	ARG	2.4
2	B	339	VAL	2.4
4	D	12	DG	2.4
2	B	800	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	796	LEU	2.3
2	B	199	ASN	2.3
2	B	343	LEU	2.3
2	B	777	SER	2.3
2	B	185	PHE	2.3
2	B	914	ALA	2.3
2	B	909	SER	2.3
2	B	1051	THR	2.3
2	B	205	GLY	2.2
2	B	830	ILE	2.2
2	B	1264	HIS	2.2
2	B	921	LEU	2.2
2	B	283	GLY	2.2
2	B	1037	PHE	2.2
2	B	852	ILE	2.1
2	B	246	LEU	2.1
2	B	811	LEU	2.1
2	B	1242	TYR	2.1
2	B	215	ARG	2.1
2	B	176	PRO	2.1
2	B	890	LYS	2.1
2	B	285	GLN	2.0
2	B	278	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
5	K	B	1407	1/1	-0.48	0.65	20.21	210,210,210,210	0
5	K	A	108	1/1	0.77	0.31	17.04	114,114,114,114	0
8	ACT	A	107	4/4	0.85	0.19	2.95	49,52,53,55	0
5	K	A	101	1/1	0.99	0.16	2.44	30,30,30,30	0
7	EDO	B	1410	4/4	0.88	0.14	1.02	45,47,47,48	0
7	EDO	B	1411	4/4	0.96	0.16	0.81	34,41,45,48	0
5	K	B	1404	1/1	0.96	0.14	0.52	94,94,94,94	0
7	EDO	A	106	4/4	0.95	0.13	-0.31	33,34,34,37	0
7	EDO	B	1409	4/4	0.96	0.10	-0.50	33,34,35,35	0
5	K	A	102	1/1	0.99	0.09	-2.67	34,34,34,34	0
5	K	B	1401	1/1	0.99	0.05	-2.97	39,39,39,39	0
5	K	B	1403	1/1	0.96	0.08	-3.12	58,58,58,58	0
6	MG	A	104	1/1	0.95	0.07	-3.24	42,42,42,42	0
6	MG	B	1408	1/1	0.96	0.10	-	46,46,46,46	0
5	K	B	1406	1/1	0.82	0.61	-	118,118,118,118	0
5	K	A	103	1/1	0.99	0.08	-	34,34,34,34	0
5	K	B	1402	1/1	0.98	0.12	-	48,48,48,48	0
5	K	B	1405	1/1	0.83	0.07	-	83,83,83,83	0
6	MG	A	105	1/1	0.97	0.11	-	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.