



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 30, 2023 – 02:02 PM JST

PDB ID : 5B2S
Title : Crystal structure of the Streptococcus pyogenes Cas9 EQR variant in complex with sgRNA and target DNA (TGAG PAM)
Authors : Hirano, S.; Nishimasu, H.; Ishitani, R.; Nureki, O.
Deposited on : 2016-02-02
Resolution : 2.20 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 : 4.02b-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

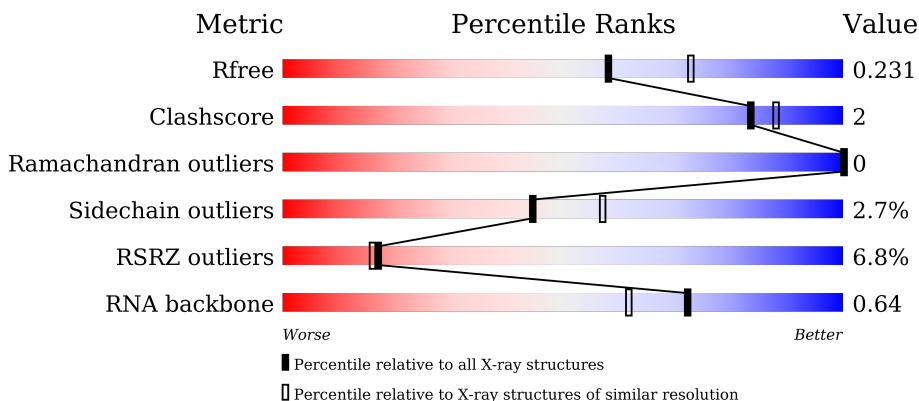
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)
RNA backbone	3102	1032 (2.60-1.80)

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	81	 2% 79% 20%
2	B	1372	 7% 88% 8%
3	C	28	 79% 21%
4	D	8	 100%

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	B	1408	-	-	-	X

2 Entry composition i

There are 9 unique types of molecules in this entry. The entry contains 13535 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
			Total	C	N	O	P			
1	A	81	1739	778	319	561	81	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	1319	10494	6695	1820	1959	20	0	1	0

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	GLU	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
			Total	C	N	O	P			
3	C	28	560	269	100	164	27	0	0	0

- 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			

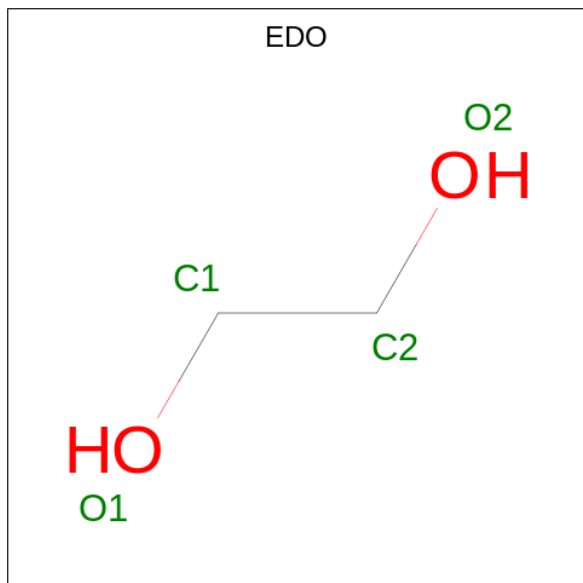
- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	K	0	0
			3	3		
5	B	8	Total	K	0	0
			8	8		

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

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

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



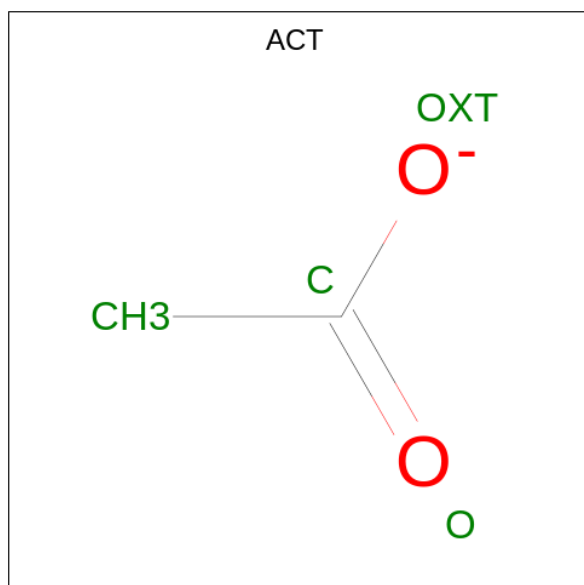
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		

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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: C₂H₃O₂).



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

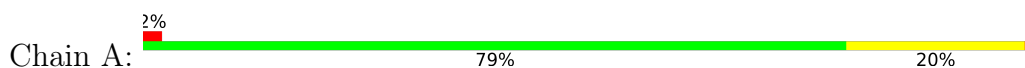
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	161	Total	O	0	0
			161	161		
9	B	343	Total	O	0	0
			343	343		
9	C	30	Total	O	0	0
			30	30		
9	D	9	Total	O	0	0
			9	9		

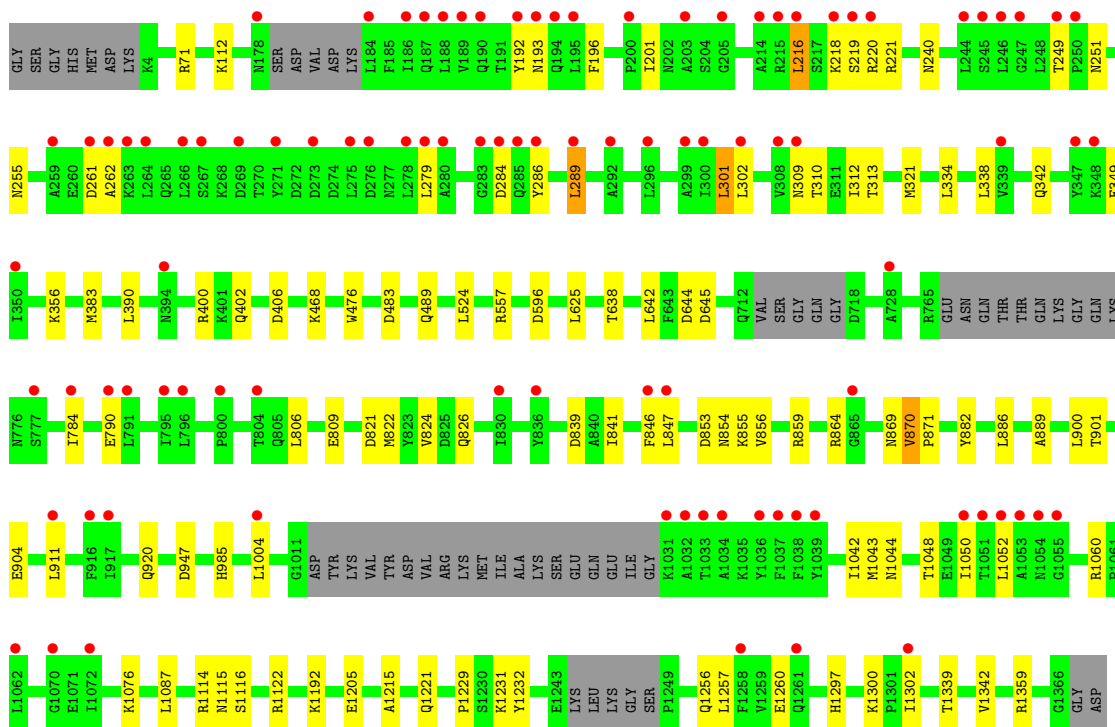
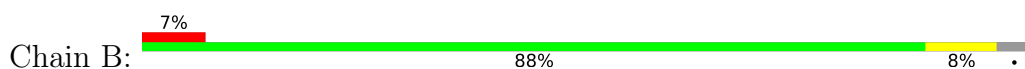
3 Residue-property plots [i](#)

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.

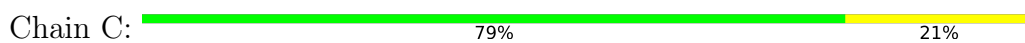
- Molecule 1: Guide RNA



- Molecule 2: CRISPR-associated endonuclease Cas9



- Molecule 3: Target DNA



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

Chain D:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.02Å 67.82Å 187.57Å 90.00° 111.12° 90.00°	Depositor
Resolution (Å)	48.01 – 2.20 48.01 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.01-2.20) 97.5 (48.01-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.10_2155: ???	Depositor
R, R_{free}	0.207 , 0.231 0.207 , 0.231	Depositor DCC
R_{free} test set	5172 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtrriage
Anisotropy	0.540	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13535	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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: EDO, MG, ACT, K

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	0.18	0/1949	0.71	0/3037
2	B	0.24	0/10682	0.40	0/14409
3	C	0.54	0/626	0.96	0/961
4	D	0.50	0/185	0.95	0/285
All	All	0.26	0/13442	0.52	0/18692

There are no bond length outliers.

There are no bond angle outliers.

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	6	0
2	B	10494	0	10384	49	0
3	C	560	0	316	5	0
4	D	165	0	93	0	0
5	A	3	0	0	0	0
5	B	8	0	0	0	0
6	A	2	0	0	0	0
6	B	1	0	0	0	0
7	A	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	12	0	18	0	0
8	A	4	0	3	0	0
9	A	161	0	0	0	0
9	B	343	0	0	0	0
9	C	30	0	0	0	0
9	D	9	0	0	0	0
All	All	13535	0	11690	58	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 2.

All (58) 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:18:G:N7	2:B:71:ARG:NH2	2.29	0.81
2:B:240:ASN:ND2	2:B:255:ASN:OD1	2.31	0.62
2:B:251:ASN:HD21	2:B:262:ALA:H	1.48	0.62
2:B:1004:LEU:HD11	2:B:1042:ILE:HD11	1.83	0.60
3:C:2:DA:H2'	3:C:3:DA:C8	2.36	0.60
1:A:44:U:O2'	2:B:402:GLN:NE2	2.36	0.58
2:B:1205:GLU:OE1	2:B:1359:ARG:NH2	2.36	0.56
2:B:251:ASN:ND2	2:B:262:ALA:H	2.04	0.55
2:B:1256:GLN:NE2	2:B:1260:GLU:OE2	2.37	0.55
2:B:1215:ALA:HB2	2:B:1221:GLN:HG3	1.89	0.54
2:B:279:LEU:HG	2:B:284:ASP:HA	1.90	0.53
2:B:790:GLU:HG2	2:B:889:ALA:HA	1.90	0.53
2:B:193:ASN:ND2	2:B:201:ILE:O	2.34	0.51
2:B:870:VAL:HG22	2:B:871:PRO:HD2	1.93	0.51
2:B:784:ILE:HD12	2:B:806:LEU:HD13	1.92	0.51
2:B:349:GLU:HG3	2:B:356:LYS:HG3	1.93	0.51
2:B:841:ILE:HD12	2:B:854:ASN:HA	1.95	0.49
2:B:809:GLU:OE1	2:B:855:LYS:NZ	2.31	0.48
2:B:1114:ARG:HG3	2:B:1116:SER:H	1.76	0.48
2:B:1297:HIS:HD2	2:B:1300:LYS:HE3	1.77	0.48
1:A:73:G:H3'	1:A:74:A:C5'	2.43	0.48
2:B:826:GLN:NE2	2:B:859:ARG:HD2	2.28	0.48
2:B:869:ASN:OD1	2:B:870:VAL:N	2.34	0.47
2:B:839:ASP:OD2	2:B:864:ARG:NE	2.37	0.47
2:B:1048:THR:HG22	2:B:1076:LYS:HD3	1.95	0.47
2:B:1044:ASN:HB3	2:B:1050:ILE:HD13	1.97	0.47
2:B:557:ARG:NH2	2:B:596:ASP:OD1	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:17:DC:H2'	3:C:18:DA:C8	2.50	0.46
2:B:1256:GLN:O	2:B:1260:GLU:HG2	2.16	0.46
2:B:312:ILE:HG23	2:B:313:THR:HG22	1.97	0.45
2:B:901:THR:O	2:B:904:GLU:HG2	2.16	0.45
2:B:338:LEU:HB3	2:B:383:MET:HE2	1.99	0.45
2:B:489:GLN:HG3	2:B:625:LEU:HD21	1.97	0.44
2:B:1042:ILE:HG23	2:B:1043:MET:HG2	1.99	0.44
2:B:1231:LYS:HE3	2:B:1232:TYR:CZ	2.52	0.44
3:C:1:DC:H2'	3:C:2:DA:C8	2.51	0.44
2:B:216:LEU:HD22	2:B:220:ARG:HG2	2.00	0.44
2:B:846:PHE:O	2:B:920:GLN:NE2	2.46	0.44
2:B:822:MET:HG3	2:B:856:VAL:HB	1.99	0.44
2:B:301:LEU:HG	2:B:302:LEU:HD13	2.00	0.44
2:B:342:GLN:HB2	2:B:383:MET:HE3	1.99	0.43
2:B:1229:PRO:HD2	2:B:1232:TYR:HD2	1.83	0.43
2:B:985:HIS:CG	2:B:1087:LEU:HD22	2.53	0.43
2:B:1060:ARG:HB2	2:B:1076:LYS:NZ	2.33	0.43
2:B:1339:THR:O	2:B:1342:VAL:HG22	2.19	0.43
1:A:71:U:H2'	1:A:72:U:C6	2.54	0.43
2:B:196:PHE:CZ	2:B:286:TYR:HE1	2.36	0.43
3:C:16:DG:H2'	3:C:17:DC:C6	2.54	0.42
2:B:882:TYR:CZ	2:B:886:LEU:HD11	2.54	0.42
2:B:400:ARG:NH2	2:B:406:ASP:OD2	2.51	0.42
2:B:192:TYR:HB2	2:B:289:LEU:HD11	2.01	0.42
1:A:3:A:H2'	1:A:4:A:C8	2.55	0.42
2:B:468:LYS:HE3	2:B:483:ASP:HA	2.02	0.42
2:B:218:LYS:HG3	2:B:221:ARG:NH1	2.35	0.41
2:B:289:LEU:HD12	2:B:289:LEU:HA	1.82	0.41
1:A:38:A:H2'	1:A:39:G:C8	2.55	0.41
3:C:17:DC:H2'	3:C:18:DA:H8	1.86	0.41
2:B:821:ASP:HB3	2:B:824:VAL:O	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	1308/1372 (95%)	1264 (97%)	44 (3%)	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	
2	B	1106/1227 (90%)	1076 (97%)	30 (3%)	44	57

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	112	LYS
2	B	216	LEU
2	B	219	SER
2	B	249	THR
2	B	261	ASP
2	B	289	LEU
2	B	301	LEU
2	B	309	ASN
2	B	310	THR
2	B	321	MET
2	B	334	LEU
2	B	390	LEU
2	B	476	TRP
2	B	524	LEU
2	B	638	THR
2	B	642	LEU
2	B	644	ASP
2	B	645	ASP
2	B	847	LEU
2	B	853	ASP
2	B	870	VAL

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Mol	Chain	Res	Type
2	B	900	LEU
2	B	911	LEU
2	B	947	ASP
2	B	1052	LEU
2	B	1115	ASN
2	B	1122	ARG
2	B	1192	LYS
2	B	1257	LEU
2	B	1302	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	251	ASN
2	B	309	ASN
2	B	402	GLN
2	B	511	HIS
2	B	726	ASN
2	B	926	GLN
2	B	1115	ASN
2	B	1261	GLN
2	B	1297	HIS
2	B	1317	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	80/81 (98%)	8 (10%)	0

All (8) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	28	A
1	A	40	C
1	A	51	A
1	A	56	U
1	A	59	U
1	A	68	A
1	A	74	A
1	A	75	A

There are no RNA pucker outliers to report.

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](#)

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	B	1411	-	3,3,3	0.47	0	2,2,2	0.32	0
7	EDO	A	106	-	3,3,3	0.46	0	2,2,2	0.38	0
8	ACT	A	107	-	3,3,3	0.75	0	3,3,3	1.31	0
7	EDO	B	1412	-	3,3,3	0.44	0	2,2,2	0.37	0
7	EDO	B	1410	-	3,3,3	0.48	0	2,2,2	0.30	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	B	1412	-	-	0/1/1/1	-
7	EDO	B	1411	-	-	0/1/1/1	-
7	EDO	A	106	-	-	0/1/1/1	-
7	EDO	B	1410	-	-	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

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%)	0.14	2 (2%) 57 55	28, 42, 127, 182	0
2	B	1319/1372 (96%)	0.37	96 (7%) 15 14	26, 53, 106, 140	0
3	C	28/28 (100%)	-0.21	0 100 100	35, 47, 72, 86	0
4	D	8/8 (100%)	0.10	0 100 100	37, 51, 97, 114	0
All	All	1436/1489 (96%)	0.34	98 (6%) 17 16	26, 53, 107, 182	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	244	LEU	8.0
2	B	216	LEU	8.0
2	B	247	GLY	7.3
2	B	189	VAL	6.3
2	B	286	TYR	6.2
2	B	1033	THR	5.8
2	B	275	LEU	5.5
2	B	847	LEU	5.5
1	A	74	A	5.5
2	B	1039	TYR	5.5
2	B	266	LEU	5.4
2	B	215	ARG	5.3
2	B	264	LEU	5.0
2	B	262	ALA	4.7
2	B	1053	ALA	4.4
2	B	299	ALA	4.3
2	B	271	TYR	4.3
2	B	283	GLY	4.2
2	B	178	ASN	4.1
2	B	1036	TYR	4.1
2	B	309	ASN	4.0

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Mol	Chain	Res	Type	RSRZ
2	B	276	ASP	3.9
2	B	280	ALA	3.9
2	B	1070	GLY	3.8
2	B	1051	THR	3.7
2	B	279	LEU	3.7
2	B	200	PRO	3.7
2	B	292	ALA	3.6
2	B	194	GLN	3.6
2	B	296	LEU	3.6
2	B	259	ALA	3.5
2	B	1052	LEU	3.4
2	B	1302	ILE	3.4
2	B	804	THR	3.4
2	B	289	LEU	3.4
2	B	192	TYR	3.4
2	B	193	ASN	3.3
2	B	219	SER	3.2
2	B	308	VAL	3.2
2	B	1258	PHE	3.1
2	B	1054	ASN	3.1
2	B	278	LEU	3.0
2	B	917	ILE	3.0
2	B	339	VAL	3.0
2	B	245	SER	3.0
2	B	1261	GLN	3.0
2	B	269	ASP	2.9
2	B	267	SER	2.9
2	B	777	SER	2.9
2	B	1050	ILE	2.9
2	B	195	LEU	2.9
2	B	300	ILE	2.8
2	B	347	TYR	2.8
2	B	1037	PHE	2.8
2	B	1034	ALA	2.8
2	B	214	ALA	2.8
2	B	800	PRO	2.8
2	B	1072	ILE	2.7
2	B	1038	PHE	2.7
2	B	285	GLN	2.7
2	B	784	ILE	2.7
2	B	728	ALA	2.6
2	B	273	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	218	LYS	2.6
2	B	246	LEU	2.6
2	B	348	LYS	2.6
1	A	75	A	2.5
2	B	186	ILE	2.5
2	B	205	GLY	2.5
2	B	302	LEU	2.5
2	B	911	LEU	2.5
2	B	1031	LYS	2.5
2	B	190	GLN	2.4
2	B	916	PHE	2.4
2	B	796	LEU	2.4
2	B	394	ASN	2.4
2	B	261	ASP	2.3
2	B	284	ASP	2.3
2	B	836	TYR	2.3
2	B	188	LEU	2.3
2	B	1032	ALA	2.3
2	B	1004	LEU	2.3
2	B	220	ARG	2.3
2	B	184	LEU	2.3
2	B	791	LEU	2.2
2	B	250	PRO	2.2
2	B	846	PHE	2.2
2	B	249	THR	2.2
2	B	795	ILE	2.1
2	B	203	ALA	2.1
2	B	263	LYS	2.1
2	B	187	GLN	2.1
2	B	350	ILE	2.1
2	B	1062	LEU	2.1
2	B	830	ILE	2.0
2	B	865	GLY	2.0
2	B	790	GLU	2.0
2	B	1055	GLY	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	K	B	1408	1/1	-0.10	0.44	160,160,160,160	0
5	K	B	1406	1/1	0.70	0.13	93,93,93,93	0
6	MG	A	104	1/1	0.76	0.06	43,43,43,43	0
6	MG	A	105	1/1	0.84	0.26	49,49,49,49	0
7	EDO	B	1411	4/4	0.84	0.24	50,50,51,51	0
5	K	B	1403	1/1	0.88	0.09	74,74,74,74	0
5	K	B	1405	1/1	0.91	0.11	81,81,81,81	0
6	MG	B	1409	1/1	0.93	0.14	41,41,41,41	0
8	ACT	A	107	4/4	0.94	0.24	50,50,50,51	0
5	K	B	1407	1/1	0.95	0.15	60,60,60,60	0
7	EDO	B	1412	4/4	0.95	0.18	39,42,45,47	0
7	EDO	A	106	4/4	0.95	0.15	30,32,34,35	0
5	K	B	1404	1/1	0.96	0.08	52,52,52,52	0
7	EDO	B	1410	4/4	0.98	0.16	31,33,34,36	0
5	K	A	103	1/1	0.98	0.10	38,38,38,38	0
5	K	B	1401	1/1	0.99	0.06	43,43,43,43	0
5	K	B	1402	1/1	0.99	0.10	47,47,47,47	0
5	K	A	102	1/1	0.99	0.10	38,38,38,38	0
5	K	A	101	1/1	0.99	0.13	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.