



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:45 AM UTC

PDB ID : 5FW2 / pdb_00005fw2
Title : Crystal structure of SpCas9 variant EQR bound to sgRNA and TGAG PAM target DNA
Authors : Anders, C.; Bargsten, K.; Jinek, M.
Deposited on : 2016-02-11
Resolution : 2.68 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

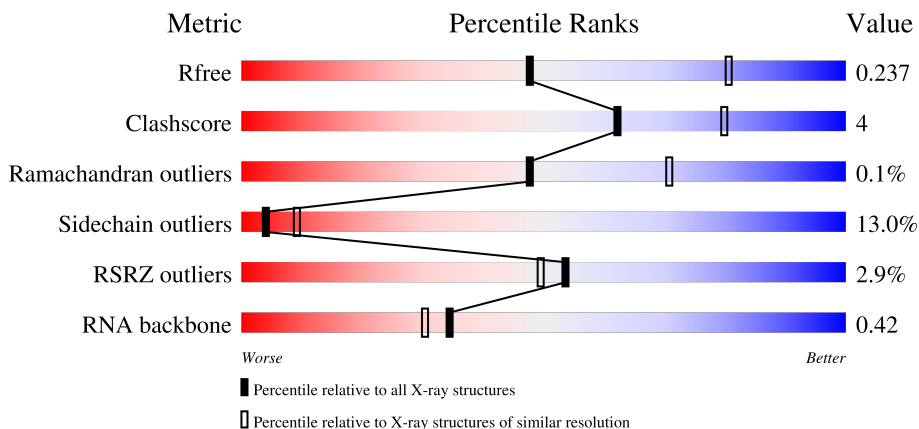
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)
RNA backbone	3983	1075 (2.90-2.46)

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	83	
2	B	1372	
3	C	28	
4	D	12	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 25767 atoms, of which 12261 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 SGRNA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	81	Total	C	H	N	O	P	0	0	0
			2601	778	869	318	555	81			

- Molecule 2 is a protein called CRISPR-ASSOCIATED ENDONUCLEASE CAS9/CSN1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	1316	Total	C	H	N	O	S	0	0	0
			21707	6862	10943	1869	2011	22			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP Q99ZW2
B	-2	ALA	-	expression tag	UNP Q99ZW2
B	-1	ALA	-	expression tag	UNP Q99ZW2
B	0	SER	-	expression tag	UNP Q99ZW2
B	10	ALA	ASP	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 STRAND.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	28	Total	C	H	N	O	P	0	0	0
			889	276	323	93	170	27			

- Molecule 4 is a DNA chain called NON-TARGET DNA STRAND.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	D	11	Total	C	H	N	O	P	0	0	0
			354	110	126	46	62	10			

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		

- Molecule 6 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	K	0	0
			2	2		
6	B	8	Total	K	0	0
			8	8		

- Molecule 7 is water.

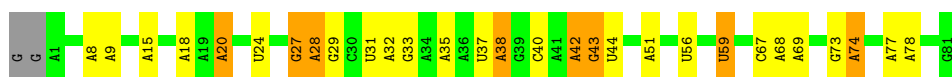
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	87	Total	O	0	0
			87	87		
7	B	107	Total	O	0	0
			107	107		
7	C	10	Total	O	0	0
			10	10		

3 Residue-property plots

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

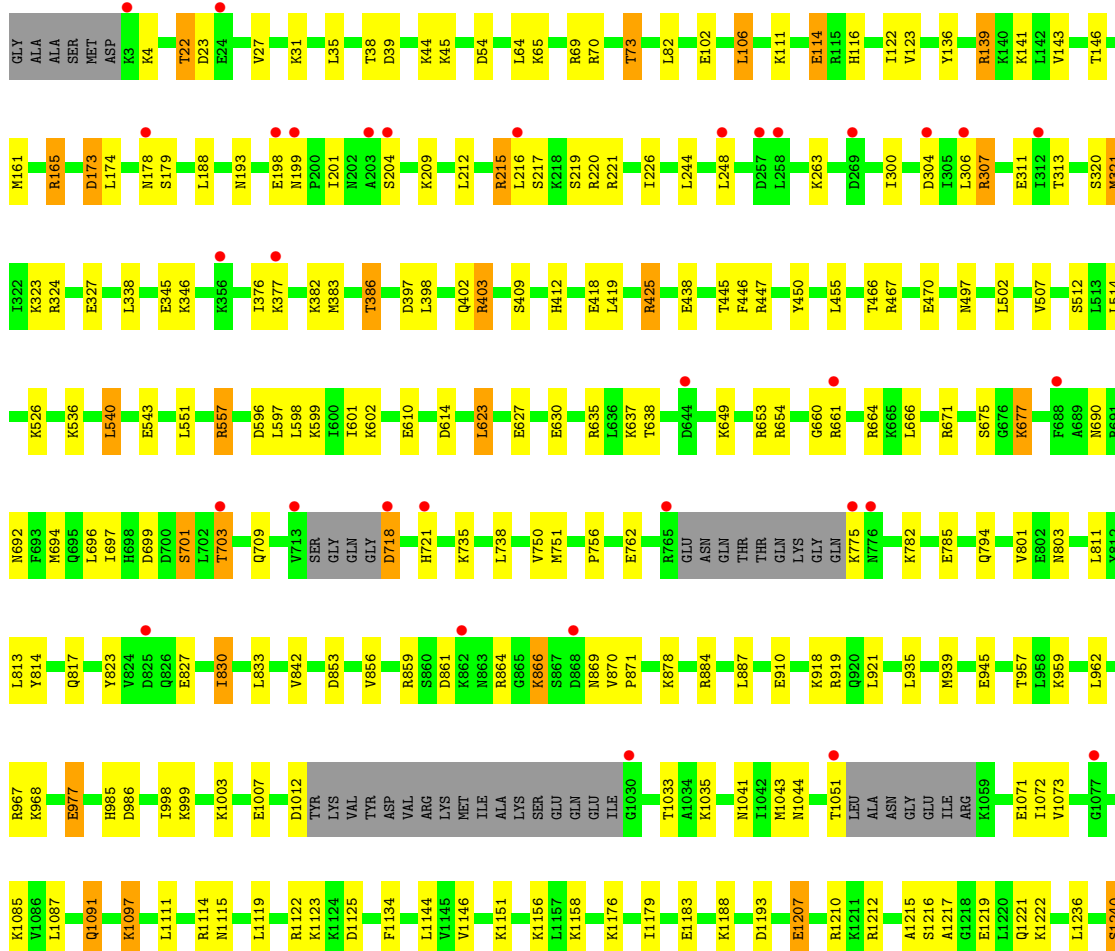
• Molecule 1: SGRNA

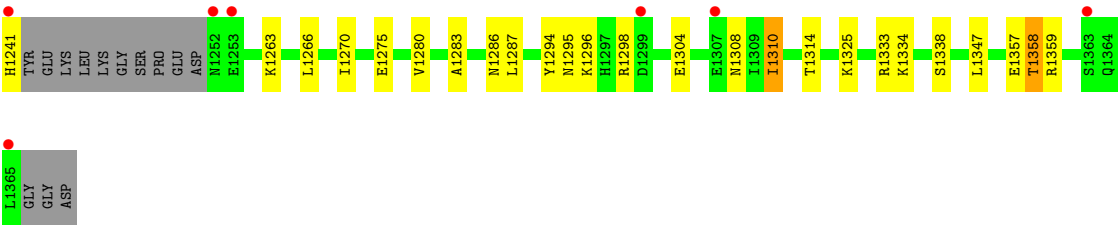
Chain A: 



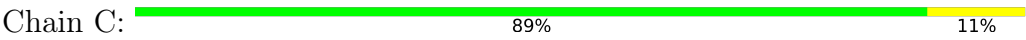
• Molecule 2: CRISPR-ASSOCIATED ENDONUCLEASE CAS9/CSN1

Chain B: 

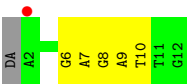




● Molecule 3: TARGET DNA STRAND



● Molecule 4: NON-TARGET DNA STRAND



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.74Å 68.24Å 188.22Å 90.00° 110.95° 90.00°	Depositor
Resolution (Å)	48.25 – 2.68 48.25 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.25-2.68) 99.5 (48.25-2.68)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.205 , 0.231 0.211 , 0.237	Depositor DCC
R_{free} test set	2983 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	44.7	Xtriage
Anisotropy	0.615	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	25767	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 3.19% 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: MG, 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	0/1942	0.40	0/3023
2	B	0.15	0/10952	0.34	0/14713
3	C	0.24	0/632	0.54	0/973
4	D	0.33	0/257	0.55	0/396
All	All	0.17	0/13783	0.37	0/19105

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	1732	869	869	16	0
2	B	10764	10943	10940	87	0
3	C	566	323	323	5	0
4	D	228	126	126	6	0
5	A	2	0	0	0	0
6	A	2	0	0	0	0
6	B	8	0	0	0	0
7	A	87	0	0	3	0
7	B	107	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	10	0	0	2	0
All	All	13506	12261	12258	102	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 4.

All (102) 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)
2:B:557:ARG:NH2	2:B:596:ASP:OD1	2.05	0.90
2:B:1286:ASN:N	7:B:2097:HOH:O	2.05	0.90
2:B:1334:LYS:NZ	3:C:3:DA:OP2	2.04	0.89
2:B:106:LEU:O	2:B:111:LYS:NZ	2.08	0.86
2:B:1283:ALA:O	7:B:2097:HOH:O	1.93	0.86
2:B:1236:LEU:O	2:B:1240:SER:OG	1.97	0.82
1:A:77:A:OP1	2:B:721:HIS:NE2	2.13	0.81
2:B:967:ARG:NH1	2:B:986:ASP:OD1	2.15	0.80
2:B:827:GLU:O	2:B:859:ARG:NH2	2.18	0.76
1:A:20:A:OP2	2:B:403:ARG:NH1	2.19	0.75
2:B:307:ARG:NH2	2:B:397:ASP:OD2	2.19	0.75
2:B:1212:ARG:NH2	2:B:1280:VAL:O	2.19	0.75
7:C:2002:HOH:O	4:D:10:DT:O2	2.08	0.69
2:B:216:LEU:O	2:B:221:ARG:NH1	2.27	0.68
2:B:1125:ASP:OD1	7:B:2079:HOH:O	2.12	0.67
7:A:2071:HOH:O	2:B:69:ARG:NH2	2.24	0.67
1:A:78:A:N3	7:A:2085:HOH:O	2.27	0.66
2:B:699:ASP:OD1	2:B:701:SER:OG	2.14	0.66
2:B:985:HIS:ND1	7:B:2069:HOH:O	2.28	0.66
2:B:709:GLN:NE2	7:B:2061:HOH:O	2.28	0.66
2:B:1304:GLU:O	2:B:1308:ASN:ND2	2.28	0.65
1:A:67:C:OP2	2:B:1097:LYS:NZ	2.31	0.64
2:B:526:LYS:NZ	2:B:690:ASN:O	2.30	0.61
2:B:614:ASP:OD1	2:B:664:ARG:NH2	2.32	0.60
2:B:718:ASP:N	2:B:718:ASP:OD1	2.34	0.60
2:B:143:VAL:O	2:B:425:ARG:NE	2.35	0.58
2:B:1357:GLU:OE1	2:B:1359:ARG:NH1	2.34	0.58
1:A:59:U:OP1	2:B:467:ARG:NH2	2.37	0.57
1:A:27:G:H4'	1:A:28:A:OP2	2.06	0.55
1:A:15:A:OP1	2:B:70:ARG:NH2	2.36	0.55
1:A:18:A:OP1	2:B:165:ARG:HD3	2.07	0.53
2:B:1287:LEU:N	7:B:2097:HOH:O	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1114:ARG:NH1	4:D:9:DA:OP1	2.42	0.52
2:B:813:LEU:O	2:B:817:GLN:HG3	2.10	0.51
2:B:977:GLU:HG3	2:B:1310:ILE:HG23	1.92	0.50
1:A:27:G:N2	1:A:44:U:OP2	2.45	0.50
1:A:73:G:C3'	1:A:74:A:H5''	2.42	0.49
2:B:321:MET:HE1	2:B:324:ARG:CZ	2.42	0.49
1:A:42:A:O2'	1:A:43:G:OP1	2.27	0.49
2:B:22:THR:HG22	2:B:23:ASP:H	1.77	0.49
2:B:1207:GLU:OE2	2:B:1210:ARG:NH1	2.46	0.49
2:B:179:SER:O	2:B:209:LYS:NZ	2.36	0.49
2:B:193:ASN:ND2	2:B:201:ILE:O	2.40	0.48
1:A:33:G:O3'	7:A:2034:HOH:O	2.19	0.48
2:B:1221:GLN:NE2	4:D:6:DG:OP2	2.46	0.48
2:B:139:ARG:NH1	2:B:418:GLU:OE1	2.46	0.48
3:C:27:DA:OP2	7:C:2010:HOH:O	2.20	0.48
2:B:866:LYS:N	2:B:866:LYS:HD3	2.28	0.48
2:B:1283:ALA:C	7:B:2097:HOH:O	2.52	0.48
4:D:7:DA:H2''	4:D:8:DG:H8	1.79	0.47
4:D:7:DA:H2''	4:D:8:DG:C8	2.49	0.47
2:B:173:ASP:OD1	2:B:173:ASP:N	2.46	0.47
2:B:497:ASN:HD21	3:C:19:DA:P	2.37	0.47
2:B:165:ARG:NH2	2:B:446:PHE:O	2.46	0.46
2:B:1286:ASN:CA	7:B:2097:HOH:O	2.60	0.46
2:B:217:SER:OG	2:B:220:ARG:HD3	2.16	0.45
2:B:623:LEU:HG	2:B:654:ARG:O	2.16	0.45
2:B:114:GLU:OE2	2:B:116:HIS:ND1	2.35	0.45
2:B:323:LYS:O	2:B:327:GLU:HG3	2.17	0.45
2:B:1122:ARG:HG3	2:B:1134:PHE:CE2	2.52	0.45
2:B:756:PRO:HD2	2:B:939:MET:HE2	1.99	0.44
2:B:450:TYR:OH	2:B:627:GLU:HG2	2.17	0.44
2:B:675:SER:HB3	2:B:677:LYS:HD3	1.98	0.44
2:B:1215:ALA:HB2	2:B:1221:GLN:HG3	1.99	0.44
2:B:870:VAL:HG13	2:B:871:PRO:HD2	1.99	0.44
2:B:447:ARG:NH2	7:B:2034:HOH:O	2.50	0.44
2:B:1333:ARG:NH2	4:D:6:DG:O6	2.44	0.44
2:B:1193:ASP:N	2:B:1193:ASP:OD1	2.51	0.43
2:B:338:LEU:HB3	2:B:383:MET:HE2	2.00	0.43
2:B:814:TYR:CZ	2:B:830:ILE:HG23	2.53	0.43
1:A:42:A:HO2'	1:A:43:G:P	2.39	0.43
2:B:215:ARG:HG3	2:B:307:ARG:CZ	2.48	0.43
2:B:703:THR:O	2:B:703:THR:HG23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:597:LEU:O	2:B:601:ILE:HG12	2.19	0.43
1:A:37:U:H2'	1:A:38:A:H5''	2.01	0.42
2:B:910:GLU:HG2	2:B:1033:THR:HG23	2.02	0.42
3:C:19:DA:H5''	3:C:19:DA:H8	1.84	0.42
2:B:35:LEU:HB2	2:B:1358:THR:HB	2.02	0.42
2:B:338:LEU:HD13	2:B:386:THR:HG22	2.02	0.42
2:B:1270:ILE:HG13	2:B:1294:TYR:CE2	2.55	0.42
2:B:861:ASP:O	2:B:864:ARG:HG2	2.19	0.42
2:B:1295:ASN:HA	2:B:1298:ARG:HD2	2.02	0.42
2:B:526:LYS:HD2	3:C:27:DA:OP1	2.20	0.41
2:B:999:LYS:HB3	2:B:1073:VAL:HG12	2.01	0.41
2:B:962:LEU:HD22	2:B:1043:MET:SD	2.60	0.41
1:A:73:G:H3'	1:A:74:A:C5'	2.50	0.41
2:B:165:ARG:O	2:B:412:HIS:HA	2.21	0.41
2:B:306:LEU:O	2:B:320:SER:HB2	2.20	0.41
2:B:507:VAL:HG11	2:B:660:GLY:O	2.21	0.41
2:B:864:ARG:NH1	2:B:869:ASN:O	2.44	0.41
2:B:1216:SER:OG	2:B:1217:ALA:N	2.53	0.41
2:B:54:ASP:O	2:B:735:LYS:NZ	2.45	0.41
2:B:321:MET:HE2	2:B:402:GLN:HA	2.03	0.41
2:B:1179:ILE:O	2:B:1183:GLU:HG3	2.21	0.40
2:B:136:TYR:CD2	2:B:321:MET:HG2	2.56	0.40
2:B:1041:ASN:HB3	2:B:1044:ASN:OD1	2.22	0.40
2:B:69:ARG:O	2:B:73:THR:HG23	2.21	0.40
2:B:204:SER:O	2:B:204:SER:OG	2.29	0.40
2:B:540:LEU:O	2:B:690:ASN:ND2	2.44	0.40
2:B:823:TYR:CE2	2:B:864:ARG:HB3	2.56	0.40
1:A:73:G:H3'	1:A:74:A:H5''	2.04	0.40
2:B:1091:GLN:HG3	2:B:1091:GLN:O	2.22	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	1304/1372 (95%)	1275 (98%)	28 (2%)	1 (0%)	48 71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	398	LEU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	B	1181/1226 (96%)	1028 (87%)	153 (13%)	4 9

All (153) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	4	LYS
2	B	22	THR
2	B	27	VAL
2	B	31	LYS
2	B	38	THR
2	B	39	ASP
2	B	44	LYS
2	B	45	LYS
2	B	64	LEU
2	B	65	LYS
2	B	73	THR
2	B	82	LEU
2	B	102	GLU
2	B	106	LEU
2	B	114	GLU
2	B	122	ILE
2	B	123	VAL
2	B	139	ARG
2	B	141	LYS
2	B	146	THR

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Mol	Chain	Res	Type
2	B	161	MET
2	B	165	ARG
2	B	173	ASP
2	B	174	LEU
2	B	178	ASN
2	B	188	LEU
2	B	198	GLU
2	B	199	ASN
2	B	212	LEU
2	B	215	ARG
2	B	219	SER
2	B	226	ILE
2	B	244	LEU
2	B	248	LEU
2	B	263	LYS
2	B	300	ILE
2	B	304	ASP
2	B	307	ARG
2	B	311	GLU
2	B	313	THR
2	B	321	MET
2	B	345	GLU
2	B	346	LYS
2	B	376	ILE
2	B	377	LYS
2	B	382	LYS
2	B	386	THR
2	B	403	ARG
2	B	409	SER
2	B	419	LEU
2	B	425	ARG
2	B	438	GLU
2	B	445	THR
2	B	455	LEU
2	B	466	THR
2	B	470	GLU
2	B	502	LEU
2	B	512	SER
2	B	514	LEU
2	B	536	LYS
2	B	540	LEU
2	B	543	GLU

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Mol	Chain	Res	Type
2	B	551	LEU
2	B	557	ARG
2	B	598	LEU
2	B	599	LYS
2	B	602	LYS
2	B	610	GLU
2	B	623	LEU
2	B	630	GLU
2	B	635	ARG
2	B	637	LYS
2	B	638	THR
2	B	649	LYS
2	B	653	ARG
2	B	661	ARG
2	B	666	LEU
2	B	671	ARG
2	B	677	LYS
2	B	692	ASN
2	B	694	MET
2	B	696	LEU
2	B	697	ILE
2	B	701	SER
2	B	703	THR
2	B	718	ASP
2	B	738	LEU
2	B	750	VAL
2	B	751	MET
2	B	762	GLU
2	B	775	LYS
2	B	782	LYS
2	B	785	GLU
2	B	794	GLN
2	B	801	VAL
2	B	803	ASN
2	B	811	LEU
2	B	830	ILE
2	B	833	LEU
2	B	842	VAL
2	B	853	ASP
2	B	856	VAL
2	B	866	LYS
2	B	878	LYS

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Mol	Chain	Res	Type
2	B	884	ARG
2	B	887	LEU
2	B	918	LYS
2	B	919	ARG
2	B	921	LEU
2	B	935	LEU
2	B	945	GLU
2	B	957	THR
2	B	959	LYS
2	B	968	LYS
2	B	977	GLU
2	B	998	ILE
2	B	1003	LYS
2	B	1007	GLU
2	B	1012	ASP
2	B	1035	LYS
2	B	1051	THR
2	B	1071	GLU
2	B	1072	ILE
2	B	1085	LYS
2	B	1087	LEU
2	B	1091	GLN
2	B	1097	LYS
2	B	1111	LEU
2	B	1115	ASN
2	B	1119	LEU
2	B	1123	LYS
2	B	1144	LEU
2	B	1146	VAL
2	B	1151	LYS
2	B	1156	LYS
2	B	1158	LYS
2	B	1176	LYS
2	B	1188	LYS
2	B	1207	GLU
2	B	1219	GLU
2	B	1222	LYS
2	B	1240	SER
2	B	1241	HIS
2	B	1263	LYS
2	B	1266	LEU
2	B	1275	GLU

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Mol	Chain	Res	Type
2	B	1296	LYS
2	B	1310	ILE
2	B	1314	THR
2	B	1325	LYS
2	B	1338	SER
2	B	1347	LEU
2	B	1358	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	129	HIS
2	B	415	HIS
2	B	497	ASN
2	B	501	ASN
2	B	674	GLN
2	B	698	HIS
2	B	803	ASN
2	B	807	GLN
2	B	980	ASN
2	B	1044	ASN
2	B	1101	GLN
2	B	1177	ASN
2	B	1264	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	80/83 (96%)	18 (22%)	5 (6%)

All (18) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	9	A
1	A	20	A
1	A	24	U
1	A	28	A
1	A	29	G
1	A	31	U
1	A	32	A

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Mol	Chain	Res	Type
1	A	35	A
1	A	38	A
1	A	40	C
1	A	42	A
1	A	43	G
1	A	51	A
1	A	56	U
1	A	59	U
1	A	68	A
1	A	69	A
1	A	74	A

All (5) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	8	A
1	A	27	G
1	A	28	A
1	A	42	A
1	A	68	A

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 12 are monoatomic - leaving 0 for Mogul analysis.

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	81/83 (97%)	-0.44	0 100 100	21, 41, 100, 108	0
2	B	1316/1372 (95%)	0.17	40 (3%) 52 48	18, 53, 92, 126	0
3	C	28/28 (100%)	-0.33	0 100 100	27, 46, 78, 92	0
4	D	11/12 (91%)	0.61	1 (9%) 15 12	36, 58, 106, 113	0
All	All	1436/1495 (96%)	0.13	41 (2%) 53 49	18, 53, 93, 126	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1365	LEU	5.7
2	B	713	VAL	5.6
2	B	306	LEU	4.6
2	B	178	ASN	4.5
2	B	1030	GLY	4.2
2	B	1241	HIS	4.0
4	D	2	DA	3.8
2	B	204	SER	3.8
2	B	3	LYS	3.6
2	B	199	ASN	3.4
2	B	721	HIS	3.4
2	B	718	ASP	3.3
2	B	312	ILE	3.2
2	B	216	LEU	3.1
2	B	868	ASP	3.0
2	B	688	PHE	3.0
2	B	269	ASP	2.9
2	B	203	ALA	2.8
2	B	775	LYS	2.8
2	B	304	ASP	2.8
2	B	1252	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	644	ASP	2.6
2	B	765	ARG	2.5
2	B	1051	THR	2.4
2	B	377	LYS	2.4
2	B	1307	GLU	2.3
2	B	248	LEU	2.3
2	B	24	GLU	2.3
2	B	661	ARG	2.3
2	B	862	LYS	2.3
2	B	198	GLU	2.3
2	B	1253	GLU	2.3
2	B	776	ASN	2.2
2	B	825	ASP	2.2
2	B	1077	GLY	2.1
2	B	356	LYS	2.1
2	B	257	ASP	2.1
2	B	1363	SER	2.1
2	B	703	THR	2.1
2	B	258	LEU	2.1
2	B	1299	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
6	K	B	2370	1/1	0.58	0.24	50,50,50,50	0
6	K	B	2372	1/1	0.88	0.12	50,50,50,50	0
5	MG	A	1083	1/1	0.89	0.20	63,63,63,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	K	A	1183	1/1	0.90	0.10	50,50,50,50	0
6	K	A	1182	1/1	0.93	0.13	79,79,79,79	0
6	K	B	2371	1/1	0.96	0.20	50,50,50,50	0
6	K	B	2366	1/1	0.97	0.05	39,39,39,39	0
6	K	B	2367	1/1	0.97	0.05	57,57,57,57	0
6	K	B	2369	1/1	0.97	0.04	43,43,43,43	0
6	K	B	2373	1/1	0.97	0.07	50,50,50,50	0
6	K	B	2368	1/1	0.98	0.04	32,32,32,32	0
5	MG	A	1084	1/1	0.99	0.11	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.