



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:05 PM UTC

PDB ID : 7QR8 / pdb_00007qr8
Title : SpCas9 bound to PTPRC off-target1 DNA substrate
Authors : Pacesa, M.; Jinek, M.
Deposited on : 2022-01-10
Resolution : 2.75 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

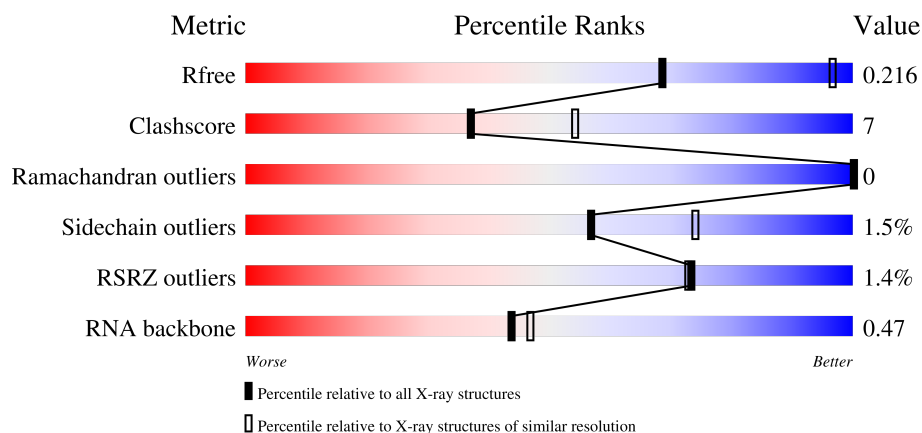
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)
RNA backbone	3983	1179 (3.00-2.52)

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	85	52% 38% 11%
2	B	1368	80% 17% .
3	C	28	68% 32%
4	D	12	8% 42% 42% 17%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 13644 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 PTPRC sgRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	85	Total	C	N	O	P	0	0	1
			1775	794	324	573	84			

- Molecule 2 is a protein called CRISPR-associated endonuclease Cas9/Csn1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1332	Total	C	N	O	S	0	0	0
			10883	6933	1889	2039	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	10	ALA	ASP	engineered mutation	UNP Q99ZW2
B	840	ALA	HIS	engineered mutation	UNP Q99ZW2

- Molecule 3 is a DNA chain called PTPRC off-target1 target strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	28	Total	C	N	O	P	0	0	0
			578	276	105	170	27			

- Molecule 4 is a DNA chain called PTPRC off-target1 non-target strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	10	Total	C	N	O	P	0	0	0
			205	98	34	63	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	3	Total 3	K 3	0	0
6	B	6	Total 6	K 6	0	0
6	C	1	Total 1	K 1	0	0

- Molecule 7 is water.

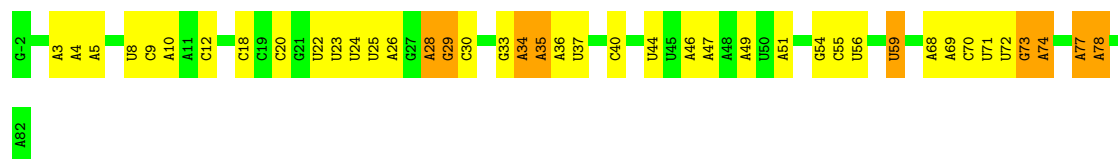
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	50	Total 50	O 50	0	0
7	B	126	Total 126	O 126	0	0
7	C	11	Total 11	O 11	0	0
7	D	4	Total 4	O 4	0	0

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.

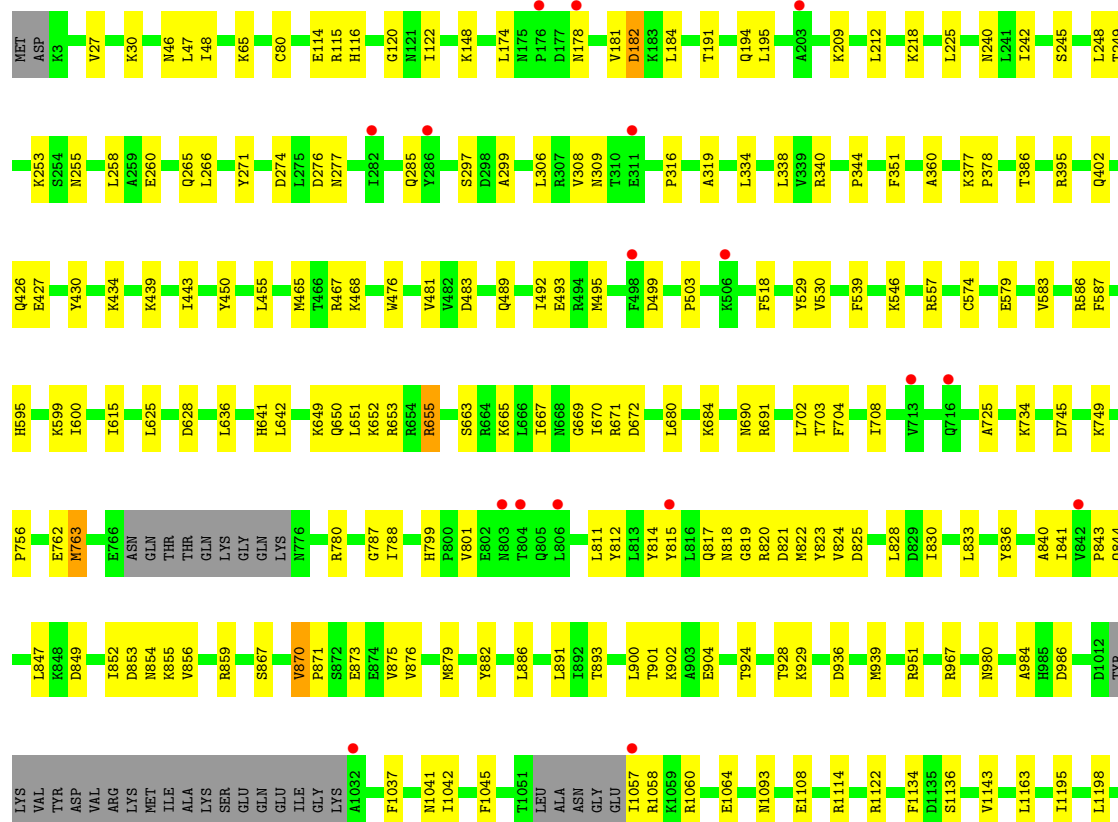
• Molecule 1: PTPRC sgRNA

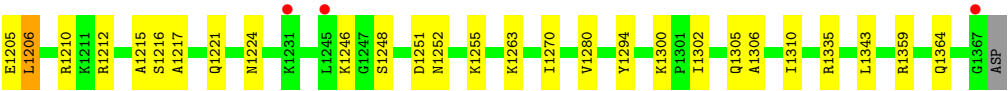
Chain A: 



• Molecule 2: CRISPR-associated endonuclease Cas9/Csn1

Chain B: 

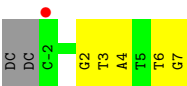




• Molecule 3: PTPRC off-target1 target strand



• Molecule 4: PTPRC off-target1 non-target strand



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.28Å 68.03Å 186.74Å 90.00° 110.74° 90.00°	Depositor
Resolution (Å)	48.12 – 2.75 48.12 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.12-2.75) 98.3 (48.12-2.75)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.222 , 0.250 (Not available) , 0.216	Depositor DCC
R_{free} test set	2898 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.649	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13644	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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	0.34	1/1987 (0.1%)	0.42	0/3095
2	B	0.16	0/11075	0.35	1/14881 (0.0%)
3	C	0.48	0/648	0.63	0/1001
4	D	0.22	0/228	0.48	0/350
All	All	0.22	1/13938 (0.0%)	0.38	1/19327 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	18	C	O3'-P	-5.05	1.53	1.61

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1206	LEU	CA-C-O	5.17	121.59	118.33

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	1775	0	894	31	0
2	B	10883	0	11054	146	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	578	0	319	8	0
4	D	205	0	115	7	0
5	A	2	0	0	0	0
6	A	3	0	0	0	0
6	B	6	0	0	0	0
6	C	1	0	0	0	0
7	A	50	0	0	0	0
7	B	126	0	0	6	0
7	C	11	0	0	0	0
7	D	4	0	0	0	0
All	All	13644	0	12382	181	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 7.

All (181) 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:1057:ILE:HG23	2:B:1058:ARG:H	1.16	1.11
2:B:1057:ILE:HG23	2:B:1058:ARG:N	1.87	0.85
2:B:178:ASN:HB2	2:B:184:LEU:HD11	1.65	0.76
2:B:825:ASP:HB2	2:B:879:MET:HE1	1.67	0.76
1:A:3:A:H2'	1:A:4:A:C8	2.22	0.74
2:B:253:LYS:HA	2:B:258:LEU:HD12	1.70	0.73
2:B:870:VAL:HG22	2:B:871:PRO:HD2	1.71	0.72
1:A:44:U:O2'	2:B:402:GLN:NE2	2.22	0.71
2:B:1057:ILE:CG2	2:B:1058:ARG:H	1.93	0.69
1:A:46:A:H2'	1:A:47:A:H8	1.57	0.69
2:B:495:MET:HE2	3:C:8:DG:H5'	1.73	0.69
1:A:5:A:H61	3:C:15:DA:H2	1.41	0.68
2:B:1215:ALA:HB2	2:B:1221:GLN:HG3	1.76	0.68
2:B:48:ILE:HG22	2:B:984:ALA:HB1	1.74	0.67
1:A:3:A:H2'	1:A:4:A:H8	1.59	0.67
2:B:849:ASP:HB3	2:B:854:ASN:HD22	1.58	0.67
4:D:6:DT:H2''	4:D:7:DG:H5''	1.79	0.65
2:B:80:CYS:SG	7:B:1588:HOH:O	2.54	0.65
1:A:46:A:H2'	1:A:47:A:C8	2.31	0.65
2:B:308:VAL:HG11	2:B:319:ALA:HB3	1.80	0.63
2:B:680:LEU:HG	2:B:684:LYS:HE3	1.81	0.62
2:B:652:LYS:O	2:B:655:ARG:NH2	2.32	0.62
2:B:182:ASP:OD1	2:B:209:LYS:HB2	2.00	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:691:ARG:NH1	2:B:702:LEU:HD23	2.14	0.61
2:B:847:LEU:HD22	2:B:849:ASP:HB2	1.81	0.61
2:B:1163:LEU:HG	2:B:1343:LEU:HD21	1.82	0.61
2:B:1114:ARG:NH1	4:D:4:DA:OP1	2.34	0.60
2:B:184:LEU:HD12	2:B:299:ALA:HB2	1.82	0.60
2:B:1263:LYS:HD2	2:B:1302:ILE:HD11	1.83	0.60
4:D:3:DT:H2''	4:D:4:DA:C8	2.36	0.60
2:B:1212:ARG:NH1	7:B:1508:HOH:O	2.35	0.59
3:C:-1:DC:H2''	3:C:0:DA:C8	2.37	0.59
2:B:649:LYS:HE2	2:B:653:ARG:HH21	1.68	0.58
2:B:114:GLU:HG3	2:B:116:HIS:H	1.67	0.58
3:C:-6:DA:H61	4:D:6:DT:H3	1.48	0.58
2:B:492:ILE:HG12	2:B:625:LEU:HD22	1.84	0.58
2:B:574:CYS:SG	7:B:1614:HOH:O	2.56	0.58
1:A:59:U:OP1	2:B:467:ARG:NH2	2.37	0.57
2:B:600:ILE:HG23	2:B:650:GLN:HB3	1.85	0.57
1:A:49:A:N3	2:B:1122:ARG:NH2	2.52	0.57
2:B:530:VAL:HG12	2:B:579:GLU:HB2	1.87	0.57
2:B:245:SER:HA	2:B:297:SER:HB2	1.87	0.56
2:B:852:ILE:HA	2:B:855:LYS:HB2	1.88	0.56
2:B:338:LEU:HD13	2:B:386:THR:HG22	1.89	0.55
2:B:780:ARG:HB2	2:B:812:TYR:CE1	2.40	0.55
2:B:395:ARG:NH1	7:B:1515:HOH:O	2.39	0.55
1:A:73:G:H5'	1:A:74:A:OP2	2.07	0.54
1:A:22:U:H2'	1:A:23:U:C6	2.43	0.54
1:A:35:A:H2'	1:A:36:A:C8	2.43	0.54
2:B:836:TYR:CD1	2:B:859:ARG:HA	2.43	0.53
1:A:74:A:OP1	1:A:74:A:H8	1.92	0.53
3:C:12:DT:H2'	3:C:13:DG:C8	2.44	0.53
2:B:1045:PHE:HD2	2:B:1064:GLU:HG2	1.74	0.52
1:A:22:U:H2'	1:A:23:U:H6	1.75	0.51
2:B:841:ILE:HD13	2:B:900:LEU:HG	1.91	0.51
2:B:1122:ARG:HG2	2:B:1134:PHE:CE2	2.45	0.51
2:B:1335:ARG:NH2	4:D:2:DG:O6	2.41	0.51
2:B:249:THR:OG1	2:B:265:GLN:NE2	2.44	0.51
2:B:843:PRO:HG2	2:B:867:SER:O	2.11	0.51
3:C:13:DG:C8	3:C:14:DT:H72	2.46	0.51
2:B:240:ASN:ND2	2:B:255:ASN:OD1	2.40	0.50
2:B:819:GLY:O	2:B:828:LEU:N	2.36	0.50
2:B:1205:GLU:OE1	2:B:1359:ARG:NH2	2.44	0.50
2:B:546:LYS:HE2	2:B:684:LYS:HD3	1.93	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1060:ARG:NH2	2:B:1064:GLU:OE2	2.44	0.50
2:B:1300:LYS:O	2:B:1305:GLN:NE2	2.45	0.50
2:B:27:VAL:HG11	2:B:48:ILE:HG12	1.93	0.50
2:B:426:GLN:NE2	7:B:1503:HOH:O	2.42	0.50
2:B:725:ALA:O	2:B:734:LYS:NZ	2.45	0.50
2:B:1057:ILE:HG12	2:B:1058:ARG:N	2.27	0.50
2:B:756:PRO:HD2	2:B:939:MET:HE2	1.94	0.49
3:C:15:DA:H2'	3:C:16:DT:H71	1.92	0.49
1:A:28:A:O4'	2:B:122:ILE:HD11	2.13	0.48
2:B:468:LYS:N	2:B:481:VAL:O	2.41	0.48
2:B:1270:ILE:HG13	2:B:1294:TYR:CE2	2.48	0.48
1:A:29:G:H2'	1:A:30:C:C6	2.49	0.48
2:B:586:ARG:HD2	2:B:587:PHE:O	2.14	0.48
2:B:702:LEU:CD1	2:B:704:PHE:CD2	2.96	0.48
1:A:29:G:H2'	1:A:30:C:H6	1.79	0.48
2:B:499:ASP:HB2	2:B:663:SER:HB3	1.95	0.48
2:B:822:MET:HG3	2:B:856:VAL:HG22	1.94	0.48
1:A:34:A:H8	1:A:34:A:P	2.35	0.48
2:B:27:VAL:HG22	2:B:48:ILE:HD11	1.95	0.48
2:B:882:TYR:CZ	2:B:886:LEU:HD11	2.49	0.48
1:A:35:A:H8	1:A:35:A:OP2	1.97	0.48
2:B:936:ASP:OD2	2:B:951:ARG:NH1	2.47	0.48
2:B:641:HIS:CE1	2:B:642:LEU:HG	2.48	0.47
2:B:274:ASP:O	2:B:277:ASN:HB3	2.15	0.47
2:B:876:VAL:HG13	2:B:901:THR:HG22	1.97	0.47
1:A:54:G:H2'	1:A:55:C:C6	2.49	0.47
2:B:1136:SER:HA	4:D:2:DG:O3'	2.15	0.47
1:A:24:U:H2'	1:A:25:U:C6	2.50	0.47
1:A:70:C:C2	1:A:71:U:C5	3.03	0.46
2:B:114:GLU:HG2	2:B:120:GLY:C	2.40	0.46
2:B:1057:ILE:CG2	2:B:1058:ARG:N	2.59	0.46
2:B:814:TYR:CZ	2:B:819:GLY:HA2	2.50	0.46
2:B:665:LYS:HA	2:B:669:GLY:HA3	1.98	0.46
2:B:1216:SER:OG	2:B:1217:ALA:N	2.48	0.46
2:B:455:LEU:HA	2:B:465:MET:HE2	1.98	0.45
2:B:840:ALA:HB2	2:B:844:GLN:HE21	1.81	0.45
2:B:1224:ASN:HB2	2:B:1280:VAL:HG11	1.99	0.45
2:B:465:MET:HE3	2:B:465:MET:HB3	1.85	0.45
1:A:34:A:P	1:A:34:A:C8	3.09	0.45
2:B:821:ASP:HB3	2:B:824:VAL:O	2.15	0.45
3:C:14:DT:H2'	3:C:15:DA:C8	2.50	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:A:C8	1:A:34:A:O5'	2.69	0.45
2:B:114:GLU:HG2	2:B:120:GLY:HA2	1.99	0.45
2:B:702:LEU:HD13	2:B:704:PHE:CD2	2.52	0.45
1:A:34:A:C8	1:A:34:A:OP1	2.69	0.45
2:B:665:LYS:HG2	2:B:670:ILE:HG23	1.98	0.45
2:B:833:LEU:HA	2:B:836:TYR:HD2	1.82	0.45
2:B:271:TYR:O	2:B:274:ASP:HB2	2.16	0.45
2:B:439:LYS:HG2	2:B:476:TRP:CD1	2.52	0.45
2:B:967:ARG:NH1	2:B:986:ASP:OD1	2.49	0.45
2:B:1206:LEU:HD13	2:B:1210:ARG:NH2	2.31	0.45
2:B:1248:SER:HB3	2:B:1251:ASP:HB2	1.99	0.45
2:B:450:TYR:HE1	2:B:628:ASP:HB2	1.82	0.44
2:B:468:LYS:HG3	2:B:483:ASP:HB2	1.99	0.44
2:B:636:LEU:HD13	2:B:651:LEU:HD22	1.99	0.44
2:B:663:SER:O	2:B:667:ILE:HG12	2.18	0.44
2:B:853:ASP:OD1	2:B:893:THR:OG1	2.32	0.44
2:B:799:HIS:HB3	2:B:815:TYR:OH	2.17	0.44
2:B:1306:ALA:O	2:B:1310:ILE:HG12	2.17	0.44
2:B:306:LEU:HD13	2:B:316:PRO:HB2	1.99	0.44
2:B:817:GLN:HG2	2:B:822:MET:SD	2.58	0.44
2:B:980:ASN:ND2	7:B:1524:HOH:O	2.51	0.44
1:A:77:A:C8	1:A:78:A:C8	3.06	0.44
2:B:489:GLN:HA	2:B:492:ILE:HD12	1.99	0.44
2:B:763:MET:HE3	2:B:928:THR:HG22	2.00	0.43
2:B:583:VAL:HG13	2:B:583:VAL:O	2.17	0.43
2:B:704:PHE:O	2:B:708:ILE:HG12	2.18	0.43
2:B:443:ILE:HD11	2:B:476:TRP:HZ2	1.83	0.43
2:B:1246:LYS:HD3	2:B:1255:LYS:HZ1	1.84	0.43
2:B:1246:LYS:HB2	2:B:1255:LYS:HZ2	1.83	0.43
2:B:148:LYS:HE3	2:B:430:TYR:CZ	2.54	0.43
2:B:691:ARG:HH12	2:B:702:LEU:HD23	1.83	0.43
2:B:218:LYS:HB2	2:B:248:LEU:HD21	2.01	0.43
2:B:1212:ARG:NH2	2:B:1280:VAL:O	2.51	0.43
2:B:340:ARG:O	2:B:344:PRO:HG3	2.18	0.43
2:B:811:LEU:HD11	2:B:815:TYR:HE2	1.84	0.43
2:B:1037:PHE:O	2:B:1041:ASN:ND2	2.51	0.43
2:B:518:PHE:CD1	2:B:667:ILE:HD12	2.53	0.42
2:B:529:TYR:HA	2:B:579:GLU:O	2.19	0.42
1:A:24:U:H2'	1:A:25:U:H6	1.84	0.42
2:B:195:LEU:HD21	2:B:285:GLN:HB3	2.02	0.42
2:B:672:ASP:OD1	2:B:703:THR:N	2.32	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:A:C2'	1:A:35:A:H5'	2.49	0.42
2:B:174:LEU:HD12	2:B:174:LEU:HA	1.78	0.42
2:B:225:LEU:HD23	2:B:242:ILE:HG21	2.02	0.42
1:A:23:U:H2'	1:A:24:U:C6	2.54	0.42
2:B:901:THR:O	2:B:904:GLU:HG2	2.20	0.42
2:B:924:THR:O	2:B:929:LYS:HE3	2.19	0.42
2:B:1246:LYS:N	2:B:1252:ASN:HD21	2.18	0.42
2:B:818:ASN:C	2:B:820:ARG:H	2.28	0.42
2:B:47:LEU:O	2:B:48:ILE:HD13	2.18	0.42
1:A:35:A:OP2	1:A:35:A:C8	2.72	0.42
2:B:212:LEU:HD21	2:B:225:LEU:HD22	2.01	0.42
1:A:26:A:OP1	2:B:115:ARG:NH1	2.53	0.41
2:B:427:GLU:HB2	2:B:434:LYS:HB2	2.02	0.41
2:B:46:ASN:O	2:B:1093:ASN:ND2	2.53	0.41
2:B:266:LEU:HD23	2:B:271:TYR:CZ	2.55	0.41
2:B:276:ASP:HB3	2:B:599:LYS:HE2	2.02	0.41
2:B:191:THR:O	2:B:194:GLN:HB2	2.20	0.41
2:B:65:LYS:NZ	2:B:1108:GLU:OE2	2.35	0.41
2:B:377:LYS:HB3	2:B:378:PRO:HD3	2.02	0.41
2:B:763:MET:HE3	2:B:763:MET:HB2	1.56	0.41
1:A:69:A:H2'	1:A:70:C:H6	1.85	0.41
2:B:557:ARG:HA	2:B:595:HIS:CD2	2.56	0.41
2:B:1246:LYS:H	2:B:1252:ASN:HD21	1.69	0.41
2:B:745:ASP:O	2:B:749:LYS:HG2	2.21	0.41
4:D:7:DG:OP1	4:D:7:DG:H4'	2.20	0.41
2:B:539:PHE:HB3	2:B:690:ASN:ND2	2.36	0.41
2:B:787:GLY:HA3	2:B:891:LEU:HD21	2.04	0.41
2:B:870:VAL:HG11	2:B:902:LYS:HB3	2.03	0.40
2:B:1143:VAL:HG13	2:B:1195:ILE:HG23	2.02	0.40
2:B:1163:LEU:HD23	2:B:1163:LEU:HA	1.94	0.40
2:B:334:LEU:O	2:B:338:LEU:HG	2.21	0.40
2:B:1163:LEU:HD21	2:B:1198:LEU:HD12	2.03	0.40
2:B:499:ASP:O	2:B:503:PRO:HA	2.21	0.40
2:B:351:PHE:C	2:B:360:ALA:HB2	2.46	0.40
2:B:823:TYR:HD1	2:B:875:VAL:HG11	1.85	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	1210/1368 (88%)	1166 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	1181/1225 (96%)	1163 (98%)	18 (2%)	57	74

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	30	LYS
2	B	181	VAL
2	B	182	ASP
2	B	260	GLU
2	B	309	ASN
2	B	493	GLU
2	B	615	ILE
2	B	655	ARG
2	B	671	ARG
2	B	762	GLU
2	B	763	MET
2	B	788	ILE
2	B	801	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	830	ILE
2	B	870	VAL
2	B	873	GLU
2	B	1042	ILE
2	B	1364	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	175	ASN
2	B	265	GLN
2	B	309	ASN
2	B	402	GLN
2	B	459	ASN
2	B	501	ASN
2	B	690	ASN
2	B	698	HIS
2	B	776	ASN
2	B	844	GLN
2	B	980	ASN
2	B	1234	ASN
2	B	1252	ASN
2	B	1254	GLN
2	B	1272	GLN
2	B	1297	HIS
2	B	1350	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	82/85 (96%)	21 (25%)	0

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	8	U
1	A	9	C
1	A	10	A
1	A	12	C
1	A	20	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	28	A
1	A	29	G
1	A	33	G
1	A	34	A
1	A	35	A
1	A	37	U
1	A	40	C
1	A	51	A
1	A	56	U
1	A	59	U
1	A	68	A
1	A	72	U
1	A	73	G
1	A	74	A
1	A	77	A
1	A	78	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 ⓘ

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	85/85 (100%)	-0.36	0 100 100	44, 68, 139, 168	0
2	B	1332/1368 (97%)	0.22	20 (1%) 72 71	40, 77, 124, 146	0
3	C	28/28 (100%)	-0.07	0 100 100	53, 86, 126, 129	0
4	D	10/12 (83%)	0.02	1 (10%) 12 12	58, 80, 116, 140	0
All	All	1455/1493 (97%)	0.18	21 (1%) 73 73	40, 77, 125, 168	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1057	ILE	4.0
2	B	498	PHE	3.4
2	B	803	ASN	3.4
2	B	1245	LEU	3.4
2	B	311	GLU	3.1
2	B	1367	GLY	2.8
2	B	716	GLN	2.8
2	B	282	ILE	2.8
2	B	506	LYS	2.8
2	B	804	THR	2.7
2	B	842	VAL	2.7
2	B	806	LEU	2.7
2	B	1032	ALA	2.6
2	B	178	ASN	2.5
2	B	1231	LYS	2.4
4	D	-2	DC	2.4
2	B	203	ALA	2.4
2	B	713	VAL	2.3
2	B	176	PRO	2.1
2	B	286	TYR	2.1
2	B	815	TYR	2.1

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	1406	1/1	0.57	0.12	119,119,119,119	0
6	K	A	103	1/1	0.77	0.14	123,123,123,123	0
6	K	B	1403	1/1	0.82	0.10	91,91,91,91	0
6	K	B	1404	1/1	0.83	0.10	101,101,101,101	0
6	K	C	101	1/1	0.85	0.15	96,96,96,96	0
6	K	B	1401	1/1	0.88	0.17	103,103,103,103	0
6	K	B	1405	1/1	0.88	0.10	113,113,113,113	0
5	MG	A	101	1/1	0.89	0.11	69,69,69,69	0
6	K	B	1402	1/1	0.96	0.08	70,70,70,70	0
6	K	A	105	1/1	0.96	0.06	67,67,67,67	0
5	MG	A	102	1/1	0.97	0.08	53,53,53,53	0
6	K	A	104	1/1	0.98	0.04	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.