



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:12 PM UTC

PDB ID : 5DLJ / pdb_00005dlj
Title : Crystal Structure of Cas-DNA-N1 complex
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Deposited on : 2015-09-05
Resolution : 2.60 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

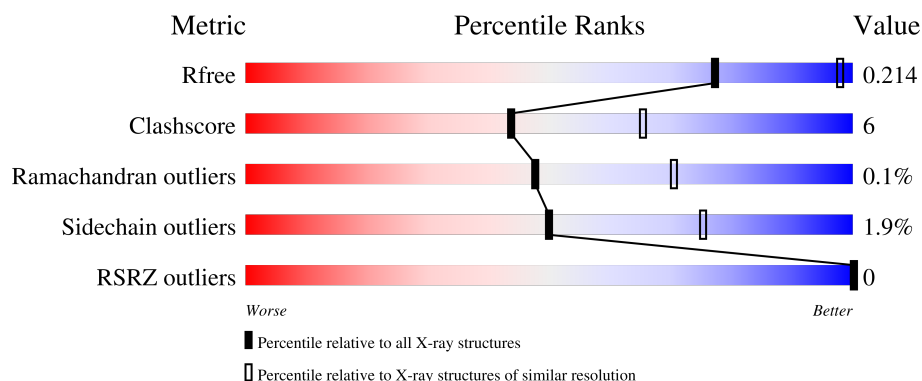
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	280	 85% 10% • 5%
1	B	280	 81% 17% •
1	C	280	 82% 16% •
1	D	280	 82% 12% • 5%
2	G	39	 69% 31%

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Mol	Chain	Length	Quality of chain
3	H	39	 69%31%
4	E	78	 86%14%
4	F	78	 73%27%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11444 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 Cas1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	267	Total	C	N	O	S	0	0	0
			2052	1311	364	370	7			
1	C	280	Total	C	N	O	S	0	0	0
			2162	1385	383	387	7			
1	A	267	Total	C	N	O	S	0	0	0
			2052	1311	364	370	7			
1	B	280	Total	C	N	O	S	0	0	0
			2162	1385	383	387	7			

- Molecule 2 is a DNA chain called 39-mer DNA N1-F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	39	Total	C	N	O	P	0	0	0
			790	383	121	248	38			

- Molecule 3 is a DNA chain called 39-mer DNA N1-R.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	39	Total	C	N	O	P	0	0	0
			787	382	119	248	38			

- Molecule 4 is a protein called CRISPR-associated endoribonuclease Cas2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	78	Total	C	N	O	S	0	0	0
			617	394	109	110	4			
4	E	78	Total	C	N	O	S	0	0	0
			617	394	109	110	4			

- Molecule 5 is water.

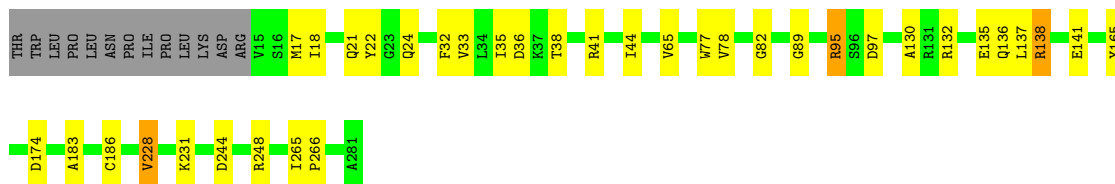
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	2	Total O 2 2	0	0
5	C	30	Total O 30 30	0	0
5	A	53	Total O 53 53	0	0
5	B	64	Total O 64 64	0	0
5	G	15	Total O 15 15	0	0
5	H	18	Total O 18 18	0	0
5	F	16	Total O 16 16	0	0
5	E	7	Total O 7 7	0	0

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.

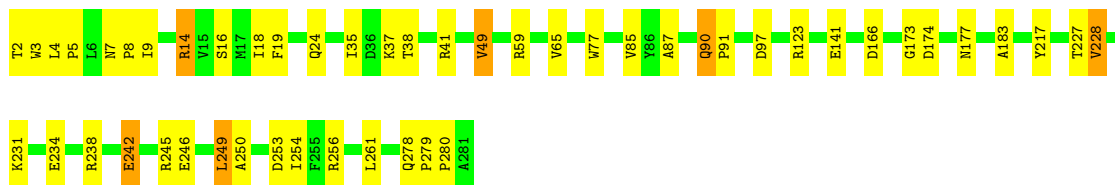
- Molecule 1: CRISPR-associated endonuclease Cas1

Chain D: 



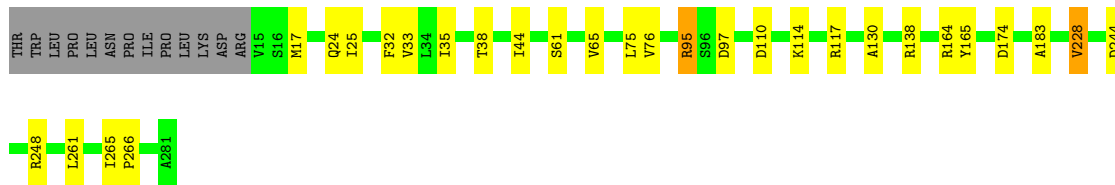
- Molecule 1: CRISPR-associated endonuclease Cas1

Chain C: 



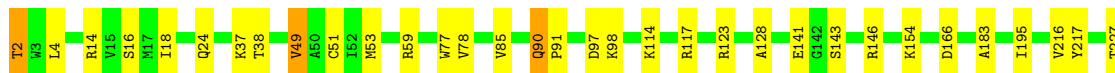
- Molecule 1: CRISPR-associated endonuclease Cas1

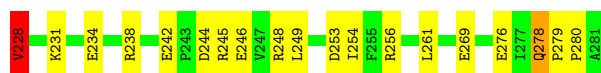
Chain A: 



- Molecule 1: CRISPR-associated endonuclease Cas1

Chain B: 





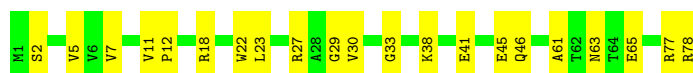
- Molecule 2: 39-mer DNA N1-F



- Molecule 3: 39-mer DNA N1-R



- Molecule 4: CRISPR-associated endoribonuclease Cas2



- Molecule 4: CRISPR-associated endoribonuclease Cas2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.89Å 193.65Å 193.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.84 – 2.60 39.84 – 2.60	Depositor EDS
% Data completeness (in resolution range)	89.1 (39.84-2.60) 89.0 (39.84-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.61Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.179 , 0.207 0.186 , 0.214	Depositor DCC
R_{free} test set	3602 reflections (4.39%)	wwPDB-VP
Wilson B-factor (Å ²)	38.3	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 15.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.477 for -h,l,k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11444	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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

5.1 Standard geometry

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	0.44	0/2092	0.88	3/2837 (0.1%)
1	B	0.46	0/2207	0.88	9/2997 (0.3%)
1	C	0.49	0/2207	0.90	8/2997 (0.3%)
1	D	0.55	1/2092 (0.0%)	0.92	4/2837 (0.1%)
2	G	0.29	0/879	0.53	0/1356
3	H	0.29	0/875	0.50	0/1349
4	E	0.39	0/628	0.78	0/851
4	F	0.51	0/628	0.82	1/851 (0.1%)
All	All	0.46	1/11608 (0.0%)	0.83	25/16075 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	18	ILE	C-O	-5.54	1.18	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	130	ALA	N-CA-C	8.34	122.72	112.54
1	B	228	VAL	N-CA-C	7.78	117.84	110.53
1	C	228	VAL	N-CA-C	6.12	116.90	110.72
1	B	242	GLU	CA-C-N	6.08	125.76	119.56
1	B	242	GLU	C-N-CA	6.08	125.76	119.56
1	C	166	ASP	CA-C-N	5.75	125.42	119.56
1	C	166	ASP	C-N-CA	5.75	125.42	119.56
1	B	166	ASP	CA-C-N	5.52	125.19	119.56
1	B	166	ASP	C-N-CA	5.52	125.19	119.56
1	D	138	ARG	N-CA-C	5.52	117.73	111.11
1	D	95	ARG	N-CA-C	5.51	117.95	109.07
1	C	90	GLN	CA-C-N	5.46	124.98	119.19
1	C	90	GLN	C-N-CA	5.46	124.98	119.19
1	A	35	ILE	N-CA-C	5.45	115.74	108.11
1	B	90	GLN	CA-C-N	5.44	124.78	118.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	90	GLN	C-N-CA	5.44	124.78	118.85
1	C	280	PRO	CB-CA-C	-5.38	102.69	111.56
1	D	35	ILE	N-CA-C	5.37	115.63	108.11
1	B	278	GLN	CA-C-N	5.34	125.89	120.38
1	B	278	GLN	C-N-CA	5.34	125.89	120.38
1	C	242	GLU	CA-C-N	-5.21	114.42	120.04
1	C	242	GLU	C-N-CA	-5.21	114.42	120.04
1	A	138	ARG	N-CA-C	5.12	117.25	111.11
4	F	12	PRO	O-C-N	5.06	123.53	121.15
1	A	95	ARG	N-CA-C	5.04	117.19	109.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	2052	0	2106	24	0
1	B	2162	0	2224	29	1
1	C	2162	0	2224	30	3
1	D	2052	0	2106	25	1
2	G	790	0	451	13	2
3	H	787	0	451	11	1
4	E	617	0	626	12	0
4	F	617	0	626	16	0
5	A	53	0	0	6	0
5	B	64	0	0	6	0
5	C	30	0	0	1	0
5	D	2	0	0	0	0
5	E	7	0	0	0	0
5	F	16	0	0	1	0
5	G	15	0	0	0	0
5	H	18	0	0	0	0
All	All	11444	0	10814	135	4

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 6.

All (135) 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:B:2:THR:CG2	5:B:363:HOH:O	1.84	1.22
4:F:5:VAL:HG21	4:E:5:VAL:HG21	1.38	1.06
1:B:2:THR:HG21	5:B:363:HOH:O	1.51	0.95
1:A:25:ILE:O	5:A:301:HOH:O	1.95	0.84
1:C:238:ARG:NH1	1:C:246:GLU:OE2	2.17	0.78
1:B:183:ALA:HB1	1:B:228:VAL:HG13	1.64	0.78
1:B:51:CYS:SG	5:B:348:HOH:O	2.43	0.75
1:D:183:ALA:HB1	1:D:228:VAL:HG13	1.70	0.73
1:A:183:ALA:HB1	1:A:228:VAL:HG13	1.69	0.73
3:H:21:DT:OP1	4:E:77:ARG:NH1	2.23	0.71
1:D:248:ARG:NH2	2:G:29:DG:OP1	2.26	0.69
1:B:53:MET:SD	5:B:348:HOH:O	2.52	0.68
1:C:2:THR:OG1	1:C:3:TRP:N	2.27	0.68
2:G:35:DT:H5'	2:G:35:DT:H6	1.59	0.67
4:F:77:ARG:NH2	5:F:101:HOH:O	2.29	0.65
1:A:75:LEU:O	5:A:303:HOH:O	2.15	0.64
2:G:21:DT:OP1	4:F:77:ARG:NH1	2.29	0.64
2:G:6:DT:H72	2:G:7:DC:H41	1.64	0.63
1:A:165:TYR:OH	5:A:302:HOH:O	2.12	0.62
1:B:253:ASP:OD1	1:B:256:ARG:NH2	2.32	0.62
1:A:248:ARG:NH2	3:H:29:DG:OP1	2.33	0.61
1:B:16:SER:HA	1:B:261:LEU:HD11	1.82	0.60
4:F:30:VAL:HG11	4:E:7:VAL:HG21	1.82	0.60
1:B:114:LYS:NZ	1:B:276:GLU:OE1	2.29	0.60
1:B:238:ARG:NH1	1:B:246:GLU:OE2	2.35	0.60
1:A:24:GLN:HB2	5:A:353:HOH:O	2.03	0.59
4:F:7:VAL:HG21	4:E:30:VAL:HG11	1.85	0.58
1:D:36:ASP:HB3	1:D:38:THR:H	1.68	0.58
1:D:132:ARG:HG2	1:D:136:GLN:HB3	1.85	0.58
1:D:135:GLU:O	1:D:138:ARG:HB2	2.05	0.56
1:C:16:SER:HA	1:C:261:LEU:HD11	1.88	0.55
1:B:123:ARG:NH1	1:B:141:GLU:OE2	2.40	0.55
4:E:2:SER:HB2	4:E:61:ALA:HB3	1.88	0.55
1:B:234:GLU:HG2	1:B:238:ARG:HH21	1.72	0.55
1:A:61:SER:N	5:A:301:HOH:O	2.22	0.55
1:B:24:GLN:OE1	1:B:59:ARG:NH2	2.41	0.54
1:D:32:PHE:HB2	1:D:44:ILE:HB	1.90	0.54
1:A:32:PHE:HB2	1:A:44:ILE:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:41:GLU:O	4:F:45:GLU:HG3	2.08	0.53
3:H:6:DT:H73	3:H:7:DC:H41	1.74	0.53
5:B:319:HOH:O	4:F:22:TRP:HD1	1.91	0.52
1:D:138:ARG:NH2	2:G:35:DT:C4	2.76	0.52
4:E:41:GLU:O	4:E:45:GLU:HG3	2.10	0.51
1:C:234:GLU:HG2	1:C:238:ARG:HH21	1.75	0.51
2:G:21:DT:P	4:F:77:ARG:HH12	2.34	0.50
1:C:24:GLN:OE1	1:C:59:ARG:NH2	2.45	0.50
4:F:2:SER:HB2	4:F:61:ALA:HB3	1.91	0.50
1:D:95:ARG:NH2	1:D:97:ASP:OD2	2.42	0.50
1:A:24:GLN:OE1	1:B:59:ARG:NH1	2.44	0.49
1:A:117:ARG:NH2	1:A:130:ALA:O	2.46	0.49
1:A:38:THR:HG23	2:G:6:DT:H5''	1.94	0.49
1:B:244:ASP:O	1:B:248:ARG:HG3	2.12	0.49
3:H:21:DT:OP1	4:E:78:ARG:NH2	2.45	0.49
2:G:22:DC:H2''	2:G:23:DA:C8	2.47	0.48
1:C:245:ARG:O	1:C:249:LEU:HD13	2.13	0.48
4:F:27:ARG:HB2	4:F:30:VAL:HB	1.96	0.48
1:D:165:TYR:HB2	2:G:33:DT:O2	2.14	0.48
1:B:2:THR:HG23	5:B:363:HOH:O	1.82	0.48
3:H:22:DC:H2''	3:H:23:DA:C8	2.48	0.48
1:D:137:LEU:O	1:D:141:GLU:HG3	2.13	0.47
1:D:244:ASP:OD1	1:D:244:ASP:N	2.47	0.47
1:C:9:ILE:O	1:C:14:ARG:NH1	2.47	0.47
3:H:21:DT:H5''	4:E:77:ARG:CZ	2.45	0.47
1:C:37:LYS:HG3	1:C:38:THR:HG23	1.96	0.47
1:B:4:LEU:HD23	1:B:4:LEU:HA	1.77	0.46
2:G:31:DT:H4'	2:G:32:DT:C6	2.51	0.46
1:B:18:ILE:HD12	1:B:49:VAL:HG21	1.97	0.46
1:A:110:ASP:O	1:A:114:LYS:HG3	2.15	0.46
3:H:21:DT:H5''	4:E:77:ARG:NH1	2.31	0.46
4:E:27:ARG:HB2	4:E:30:VAL:HB	1.96	0.46
1:A:38:THR:CG2	2:G:6:DT:H5''	2.44	0.46
1:C:123:ARG:NH1	1:C:141:GLU:OE2	2.48	0.46
1:C:90:GLN:HA	1:C:91:PRO:HD3	1.76	0.45
3:H:21:DT:P	4:E:77:ARG:HH12	2.37	0.45
1:B:97:ASP:OD1	1:B:97:ASP:N	2.49	0.45
1:C:174:ASP:HB2	1:C:177:ASN:HD22	1.79	0.45
1:A:244:ASP:OD1	1:A:244:ASP:N	2.50	0.45
1:D:21:GLN:HG2	1:D:22:TYR:CG	2.52	0.45
1:D:24:GLN:OE1	1:C:59:ARG:NH1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:GLN:HA	1:B:91:PRO:HD3	1.75	0.45
4:F:38:LYS:HE2	4:F:38:LYS:HB2	1.67	0.45
1:A:265:ILE:HB	1:A:266:PRO:HD3	1.99	0.44
1:C:19:PHE:CB	1:C:245:ARG:HH12	2.30	0.44
2:G:21:DT:OP1	4:F:78:ARG:NH2	2.51	0.44
1:D:265:ILE:HB	1:D:266:PRO:HD3	2.00	0.44
1:B:245:ARG:O	1:B:249:LEU:HG	2.17	0.44
4:F:7:VAL:HA	4:F:29:GLY:O	2.18	0.44
1:D:77:TRP:CD2	1:C:65:VAL:HG11	2.53	0.44
1:D:78:VAL:HB	1:D:82:GLY:HA2	2.00	0.44
1:A:95:ARG:HB3	1:A:97:ASP:OD1	2.17	0.44
1:C:18:ILE:HD12	1:C:49:VAL:HG21	1.99	0.44
1:A:76:VAL:HG22	5:A:303:HOH:O	2.18	0.44
1:A:174:ASP:N	1:A:174:ASP:OD1	2.51	0.44
1:C:24:GLN:HB3	1:C:59:ARG:HH21	1.83	0.43
1:C:227:THR:HB	1:C:254:ILE:HG21	2.00	0.43
1:A:65:VAL:HG11	1:B:77:TRP:CG	2.53	0.43
2:G:35:DT:H5'	2:G:35:DT:C6	2.47	0.43
1:B:278:GLN:HA	1:B:279:PRO:HD3	1.75	0.43
4:E:7:VAL:HA	4:E:29:GLY:O	2.18	0.43
1:C:278:GLN:HA	1:C:279:PRO:HD3	1.89	0.43
1:A:164:ARG:HA	3:H:36:DT:O2	2.17	0.43
1:C:7:ASN:HA	1:C:8:PRO:HD3	1.92	0.43
1:C:256:ARG:HE	4:F:65:GLU:HB2	1.82	0.43
1:D:183:ALA:HB1	1:D:228:VAL:CG1	2.45	0.43
1:B:37:LYS:HG3	1:B:38:THR:HG23	2.00	0.43
1:B:98:LYS:NZ	1:B:269:GLU:OE1	2.52	0.43
3:H:13:DT:H2''	3:H:14:DG:C8	2.54	0.43
1:A:183:ALA:HB1	1:A:228:VAL:CG1	2.45	0.43
1:D:17:MET:HB3	1:D:17:MET:HE2	1.80	0.43
4:F:23:LEU:HD22	4:F:33:GLY:HA3	2.01	0.43
1:C:231:LYS:HD3	1:C:254:ILE:HD11	2.00	0.43
1:C:4:LEU:HD23	1:C:5:PRO:HD2	2.01	0.42
1:B:227:THR:HB	1:B:254:ILE:HG21	2.00	0.42
1:C:35:ILE:HG12	1:C:41:ARG:HG2	2.02	0.42
1:B:195:ILE:HD12	1:B:216:VAL:HG22	2.02	0.42
1:A:17:MET:HE2	1:A:17:MET:HB3	1.86	0.42
1:B:143:SER:HA	1:B:146:ARG:HH21	1.83	0.42
1:D:65:VAL:HG11	1:C:77:TRP:CG	2.55	0.42
1:D:95:ARG:HB3	1:D:97:ASP:OD1	2.19	0.42
1:C:173:GLY:N	5:C:306:HOH:O	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:31:DT:H4'	3:H:32:DT:C6	2.55	0.41
1:D:174:ASP:OD1	1:D:174:ASP:N	2.51	0.41
1:C:253:ASP:OD1	1:C:256:ARG:NH1	2.48	0.41
1:A:17:MET:SD	1:A:261:LEU:HD22	2.60	0.41
1:C:183:ALA:HB1	1:C:228:VAL:HB	2.02	0.41
1:D:97:ASP:OD1	1:D:97:ASP:N	2.54	0.41
1:D:186:CYS:SG	1:D:248:ARG:HG3	2.61	0.41
1:C:97:ASP:OD1	1:C:97:ASP:N	2.52	0.41
1:C:231:LYS:HG2	1:C:250:ALA:HB1	2.02	0.41
1:B:117:ARG:NH2	1:B:128:ALA:O	2.49	0.41
4:F:18:ARG:NH1	4:F:46:GLN:OE1	2.52	0.41
1:D:36:ASP:OD2	1:D:41:ARG:NE	2.44	0.40
1:A:97:ASP:OD1	1:A:97:ASP:N	2.54	0.40
1:B:53:MET:HE3	1:B:78:VAL:HG11	2.03	0.40
1:D:89:GLY:HA2	1:C:87:ALA:O	2.22	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:132:ARG:NH1	1:C:173:GLY:O[1_455]	1.94	0.26
1:C:217:TYR:OH	2:G:3:DT:OP1[3_445]	2.16	0.04
1:C:37:LYS:NZ	2:G:38:DT:O4[1_655]	2.18	0.02
1:B:217:TYR:OH	3:H:3:DT:OP1[2_344]	2.18	0.02

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	265/280 (95%)	262 (99%)	3 (1%)	0	100	100
1	B	278/280 (99%)	272 (98%)	5 (2%)	1 (0%)	30	51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	278/280 (99%)	271 (98%)	7 (2%)	0	100	100
1	D	265/280 (95%)	261 (98%)	4 (2%)	0	100	100
4	E	76/78 (97%)	74 (97%)	2 (3%)	0	100	100
4	F	76/78 (97%)	74 (97%)	2 (3%)	0	100	100
All	All	1238/1276 (97%)	1214 (98%)	23 (2%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	280	PRO

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	
1	A	212/225 (94%)	210 (99%)	2 (1%)	70	87
1	B	225/225 (100%)	218 (97%)	7 (3%)	35	63
1	C	225/225 (100%)	220 (98%)	5 (2%)	45	72
1	D	212/225 (94%)	209 (99%)	3 (1%)	59	81
4	E	64/64 (100%)	64 (100%)	0	100	100
4	F	64/64 (100%)	62 (97%)	2 (3%)	35	63
All	All	1002/1028 (98%)	983 (98%)	19 (2%)	50	75

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	33	VAL
1	D	228	VAL
1	D	231	LYS
1	C	14	ARG
1	C	49	VAL
1	C	85	VAL

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Mol	Chain	Res	Type
1	C	242	GLU
1	C	249	LEU
1	A	33	VAL
1	A	228	VAL
1	B	2	THR
1	B	14	ARG
1	B	49	VAL
1	B	85	VAL
1	B	154	LYS
1	B	228	VAL
1	B	231	LYS
4	F	11	VAL
4	F	63	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	43	HIS
1	D	278	GLN
1	A	278	GLN
4	E	71	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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 ⓘ

There are no ligands in this entry.

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	267/280 (95%)	-1.62	0 100 100	17, 42, 75, 102	0
1	B	280/280 (100%)	-1.66	0 100 100	19, 41, 78, 116	0
1	C	280/280 (100%)	-1.66	0 100 100	18, 40, 74, 101	0
1	D	267/280 (95%)	-1.64	0 100 100	17, 40, 76, 98	0
2	G	39/39 (100%)	-1.90	0 100 100	30, 61, 129, 141	0
3	H	39/39 (100%)	-1.88	0 100 100	28, 61, 133, 148	0
4	E	78/78 (100%)	-1.65	0 100 100	24, 43, 84, 109	0
4	F	78/78 (100%)	-1.64	0 100 100	24, 42, 80, 107	0
All	All	1328/1354 (98%)	-1.66	0 100 100	17, 42, 80, 148	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.