



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2026 – 08:40 PM UTC

PDB ID : 6LWQ / pdb_00006lwq
Title : Crystal structure of human NEIL1(R242) bound to duplex DNA containing a C:T mismatch
Authors : Liu, M.H.; Zhang, J.; Zhu, C.X.; Zhang, X.X.; Gao, Y.Q.; Yi, C.Q.
Deposited on : 2020-02-07
Resolution : 2.89 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

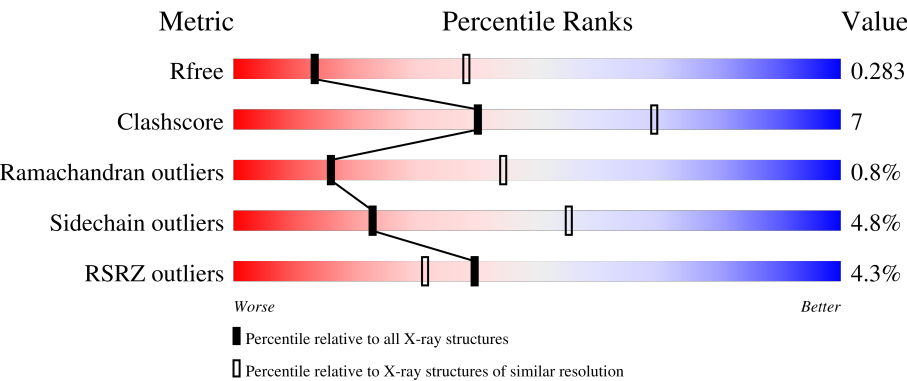
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	295	% 75% 14% .. 8%
1	D	295	2% 74% 13% .. 12%
1	G	295	9% 69% 7% . 22%
2	B	13	77% 23%
2	E	13	69% 23% 8%

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Mol	Chain	Length	Quality of chain
2	H	13	8% 69% 23% 8%
3	C	13	54% 46%
3	F	13	62% 31% 8%
3	I	13	38% 38% 23%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7521 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 Endonuclease 8-like 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2142	1365	390	377	10			
1	D	260	Total	C	N	O	S	0	0	0
			2047	1306	374	358	9			
1	G	229	Total	C	N	O	S	0	0	0
			1659	1028	322	302	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	242	ARG	LYS	variant	UNP Q96FI4
D	242	ARG	LYS	variant	UNP Q96FI4
G	242	ARG	LYS	variant	UNP Q96FI4

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*GP*TP*CP*CP*AP*TP*GP*TP*CP*TP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	13	Total	C	N	O	P	0	0	0
			258	125	43	78	12			
2	E	13	Total	C	N	O	P	0	0	0
			258	125	43	78	12			
2	H	13	Total	C	N	O	P	0	0	0
			258	125	43	78	12			

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	P	0	0	0
			267	127	53	75	12			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	13	Total 267	C 127	N 53	O 75	P 12	0	0	0
3	I	13	Total 267	C 127	N 53	O 75	P 12	0	0	0

- Molecule 4 is water.

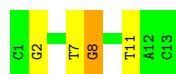
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total 49	O 49	0	0
4	B	5	Total 5	O 5	0	0
4	C	1	Total 1	O 1	0	0
4	D	39	Total 39	O 39	0	0
4	E	4	Total 4	O 4	0	0

Chain B:  77% 23%



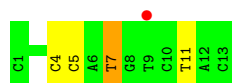
- Molecule 2: DNA (5'-D(*CP*GP*TP*CP*CP*AP*TP*GP*TP*CP*TP*AP*C)-3')

Chain E:  69% 23% 8%



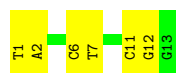
- Molecule 2: DNA (5'-D(*CP*GP*TP*CP*CP*AP*TP*GP*TP*CP*TP*AP*C)-3')

Chain H:  8% 69% 23% 8%



- Molecule 3: DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3')

Chain C:  54% 46%

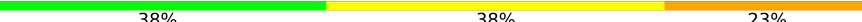


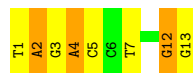
- Molecule 3: DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3')

Chain F:  62% 31% 8%



- Molecule 3: DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3')

Chain I:  38% 38% 23%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.08Å 109.23Å 171.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.03 – 2.89 43.03 – 2.89	Depositor EDS
% Data completeness (in resolution range)	97.5 (43.03-2.89) 97.5 (43.03-2.89)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.237 , 0.283 0.237 , 0.283	Depositor DCC
R_{free} test set	1529 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	68.0	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 71.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7521	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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	1.10	3/2201 (0.1%)	1.44	6/2977 (0.2%)
1	D	1.12	2/2102 (0.1%)	1.51	12/2843 (0.4%)
1	G	1.05	0/1695	1.41	2/2267 (0.1%)
2	B	0.49	0/287	1.03	1/440 (0.2%)
2	E	0.52	0/287	1.12	4/440 (0.9%)
2	H	0.45	0/287	1.01	2/440 (0.5%)
3	C	0.56	0/300	0.89	0/462
3	F	0.51	0/300	0.88	1/462 (0.2%)
3	I	0.47	0/300	1.17	5/462 (1.1%)
All	All	0.99	5/7759 (0.1%)	1.36	33/10793 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	150	ARG	CD-NE	12.43	1.63	1.46
1	D	150	ARG	NE-CZ	11.27	1.45	1.33
1	A	95	ARG	C-O	6.20	1.31	1.23
1	A	83	GLY	C-O	5.55	1.31	1.23
1	A	91	GLU	CG-CD	5.48	1.65	1.52

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	105	PRO	CA-C-N	17.60	141.84	119.84
1	D	105	PRO	C-N-CA	17.60	141.84	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	3	DG	C4'-C3'-O3'	9.44	124.15	110.00
1	D	104	ALA	CA-C-N	8.15	128.77	120.38
1	D	104	ALA	C-N-CA	8.15	128.77	120.38
2	E	11	DT	C2'-C3'-O3'	-7.33	100.50	111.50
1	D	105	PRO	C-N-CD	-6.99	96.34	125.00
2	E	8	DG	O5'-P-OP1	-6.89	88.32	109.00
1	A	106	PRO	N-CA-C	-6.88	98.31	112.47
3	I	4	DA	C2'-C3'-O3'	6.75	121.62	111.50
1	A	105	PRO	N-CA-C	6.43	118.55	110.70
3	I	3	DG	C2'-C3'-O3'	-6.43	101.86	111.50
2	H	7	DT	C2'-C3'-O3'	-6.21	102.18	111.50
1	G	199	GLU	CA-C-N	6.14	128.79	120.38
1	G	199	GLU	C-N-CA	6.14	128.79	120.38
1	D	107	GLY	CA-C-N	-6.03	113.31	120.13
1	D	107	GLY	C-N-CA	-6.03	113.31	120.13
3	I	12	DG	C2'-C3'-O3'	5.86	120.29	111.50
1	D	18	CYS	CA-C-O	-5.86	112.81	119.78
2	B	11	DT	C2'-C3'-O3'	-5.83	102.76	111.50
3	I	2	DA	C2'-C3'-O3'	5.75	120.12	111.50
2	H	11	DT	C2'-C3'-O3'	-5.46	103.30	111.50
1	D	110	LEU	O-C-N	5.46	129.96	123.24
2	E	8	DG	O5'-P-OP2	5.31	123.92	108.00
1	A	273	ASP	CA-CB-CG	5.27	117.87	112.60
1	D	23	PHE	CB-CA-C	5.20	118.39	109.80
1	A	110	LEU	O-C-N	5.19	129.63	123.24
1	D	252	ASP	CA-CB-CG	5.12	117.72	112.60
1	A	18	CYS	CA-C-O	-5.12	113.57	119.56
1	D	131	PRO	N-CA-CB	5.06	107.82	103.27
3	F	3	DG	C2'-C3'-O3'	5.05	119.08	111.50
2	E	2	DG	C4'-C3'-O3'	5.02	117.53	110.00
1	A	19	ARG	CA-C-O	-5.02	115.53	120.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	PRO	Peptide

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	2142	0	2110	28	1
1	D	2047	0	2005	34	1
1	G	1659	0	1473	13	0
2	B	258	0	149	2	0
2	E	258	0	149	4	0
2	H	258	0	149	3	0
3	C	267	0	147	7	0
3	F	267	0	147	4	0
3	I	267	0	147	8	0
4	A	49	0	0	4	0
4	B	5	0	0	1	0
4	C	1	0	0	0	0
4	D	39	0	0	4	0
4	E	4	0	0	0	0
All	All	7521	0	6476	90	1

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 (90) 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:D:105:PRO:CB	1:D:106:PRO:HD2	1.65	1.15
1:D:105:PRO:HB3	1:D:106:PRO:HD2	1.27	1.10
1:D:104:ALA:HB1	1:D:105:PRO:CD	1.81	1.09
1:D:104:ALA:CB	1:D:105:PRO:HD2	1.85	1.06
1:A:47:ILE:O	1:D:150:ARG:NH1	1.92	1.02
1:G:89:PRO:HG2	1:G:92:GLU:HB2	1.40	1.00
1:D:104:ALA:HB1	1:D:105:PRO:HD2	0.97	0.95
1:A:159:ARG:NH2	1:A:274:ARG:HE	1.78	0.80
1:A:41:GLU:O	1:A:70:GLN:NE2	2.17	0.76
1:G:55:GLU:OE2	1:G:130:GLN:HG3	1.87	0.74
1:A:250:GLU:OE1	1:A:253:PHE:HB3	1.89	0.71
1:D:189:PRO:HB2	1:D:192:GLU:HG2	1.72	0.70
2:H:4:DC:H2'	2:H:5:DC:C6	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:ARG:NH1	4:A:301:HOH:O	2.26	0.67
1:D:2:PRO:O	4:D:301:HOH:O	2.15	0.64
3:I:4:DA:H2'	3:I:5:DC:O4'	2.00	0.62
1:D:81:MET:HE2	2:E:8:DG:C8	2.36	0.60
1:G:103:THR:HG1	1:G:110:LEU:N	2.00	0.60
1:A:26:CYS:HB2	4:A:341:HOH:O	2.03	0.59
3:C:1:DT:O4'	3:I:13:DG:H2''	2.03	0.58
1:D:167:ASP:OD1	1:D:169:ARG:HB2	2.04	0.58
3:F:2:DA:H2''	3:F:3:DG:OP2	2.03	0.58
3:C:11:DC:H2''	3:C:12:DG:O5'	2.04	0.57
1:A:249:GLY:C	1:A:251:GLU:H	2.13	0.57
1:G:2:PRO:N	2:H:7:DT:H2''	2.20	0.56
1:A:159:ARG:CZ	1:A:274:ARG:HE	2.19	0.56
1:A:47:ILE:C	1:D:150:ARG:NH1	2.64	0.56
1:D:81:MET:CE	2:E:8:DG:C8	2.90	0.54
1:A:167:ASP:OD1	1:A:169:ARG:HB2	2.06	0.54
1:D:99:LEU:C	1:D:100:ARG:HG3	2.33	0.54
1:D:99:LEU:HD13	1:D:123:TRP:CZ2	2.43	0.54
1:G:167:ASP:OD1	1:G:169:ARG:HB2	2.07	0.54
1:A:59:ILE:HD11	4:A:345:HOH:O	2.08	0.53
3:C:1:DT:H2''	3:C:2:DA:C8	2.44	0.52
1:D:146:GLU:HG3	1:D:150:ARG:CD	2.40	0.52
3:I:12:DG:H1'	3:I:13:DG:OP1	2.10	0.52
3:I:1:DT:H2'	3:I:2:DA:N7	2.26	0.51
1:A:159:ARG:NH2	1:A:274:ARG:NE	2.54	0.51
3:F:2:DA:H1'	3:F:3:DG:H5'	1.92	0.51
1:D:93:LEU:HD13	1:D:100:ARG:HD2	1.92	0.50
1:G:99:LEU:HD13	1:G:123:TRP:CZ2	2.47	0.50
1:A:99:LEU:HD13	1:A:123:TRP:CZ2	2.47	0.49
1:D:137:VAL:HG12	1:D:144:PHE:CE1	2.47	0.49
1:A:59:ILE:CD1	4:A:345:HOH:O	2.60	0.49
1:D:146:GLU:OE2	1:D:150:ARG:HD2	2.12	0.49
3:F:1:DT:H2''	3:F:2:DA:C8	2.48	0.49
3:C:1:DT:C5'	3:I:13:DG:H2''	2.43	0.49
1:D:176:ASN:O	4:D:302:HOH:O	2.20	0.48
2:B:12:DA:H61	3:C:1:DT:H3	1.61	0.48
1:D:25:GLY:O	1:D:41:GLU:HA	2.14	0.47
1:D:132:GLY:O	1:D:169:ARG:HG2	2.14	0.47
1:G:239:LEU:HD23	1:G:239:LEU:N	2.30	0.47
3:C:1:DT:C2'	3:C:2:DA:C8	2.99	0.46
3:F:6:DC:H2'	3:F:7:DT:C6	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:MET:HE1	1:A:277:ARG:NH2	2.30	0.46
2:B:9:DT:H1'	4:B:102:HOH:O	2.15	0.46
1:A:242:ARG:HG3	1:A:247:GLU:HG2	1.99	0.46
1:D:198:LEU:O	1:D:201:LEU:HD12	2.16	0.45
1:D:17:ALA:HB1	1:D:87:LEU:HD22	1.98	0.45
1:A:137:VAL:HG12	1:A:144:PHE:CE1	2.51	0.45
1:A:242:ARG:HG3	1:A:247:GLU:CG	2.46	0.45
1:A:25:GLY:O	1:A:41:GLU:HA	2.17	0.45
1:A:93:LEU:HD13	1:A:100:ARG:HD2	1.98	0.45
1:D:104:ALA:CB	1:D:105:PRO:CD	2.55	0.45
1:A:99:LEU:C	1:A:100:ARG:HG3	2.42	0.44
1:D:2:PRO:HA	2:E:7:DT:O2	2.17	0.44
1:D:120:PHE:HB2	4:D:309:HOH:O	2.18	0.44
1:A:198:LEU:O	1:A:201:LEU:HD12	2.17	0.44
1:D:187:LYS:NZ	1:D:282:GLN:OE1	2.43	0.44
1:D:277:ARG:NH2	2:E:7:DT:OP2	2.48	0.43
1:A:132:GLY:O	1:A:169:ARG:HG2	2.19	0.43
3:C:6:DC:H2'	3:C:7:DT:C6	2.54	0.42
1:G:130:GLN:HB2	1:G:133:ARG:HD2	2.00	0.42
1:G:117:ILE:O	3:I:7:DT:H4'	2.18	0.42
1:A:155:LYS:HD3	3:I:12:DG:OP1	2.19	0.42
1:D:150:ARG:NH1	4:D:304:HOH:O	2.52	0.42
3:I:4:DA:C2'	3:I:5:DC:O4'	2.66	0.41
1:G:132:GLY:O	1:G:169:ARG:HG2	2.21	0.41
1:D:81:MET:HE1	1:D:277:ARG:NH2	2.36	0.41
1:D:137:VAL:HG12	1:D:144:PHE:CD1	2.56	0.41
1:D:155:LYS:HE2	1:D:155:LYS:HB2	1.83	0.41
1:G:277:ARG:NH2	2:H:7:DT:OP2	2.54	0.41
1:A:195:ARG:HG2	1:A:199:GLU:CD	2.46	0.40
1:A:249:GLY:C	1:A:251:GLU:N	2.78	0.40
1:A:105:PRO:HA	1:A:106:PRO:O	2.22	0.40
1:G:137:VAL:HG12	1:G:144:PHE:CE1	2.56	0.40
1:G:99:LEU:C	1:G:100:ARG:HG3	2.47	0.40
1:A:145:ARG:NH1	1:A:226:LEU:HB2	2.37	0.40
1:D:105:PRO:HB3	1:D:106:PRO:CD	2.19	0.40
1:D:146:GLU:OE1	1:D:146:GLU:HA	2.21	0.40

All (1) 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:A:199:GLU:OE1	1:D:34:ARG:NH2[3_445]	2.16	0.04

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	266/295 (90%)	247 (93%)	17 (6%)	2 (1%)	16	44
1	D	254/295 (86%)	235 (92%)	16 (6%)	3 (1%)	10	34
1	G	215/295 (73%)	198 (92%)	16 (7%)	1 (0%)	24	54
All	All	735/885 (83%)	680 (92%)	49 (7%)	6 (1%)	16	44

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	106	PRO
1	A	244	TYR
1	A	106	PRO
1	D	104	ALA
1	G	90	ARG
1	D	83	GLY

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	
1	A	224/248 (90%)	210 (94%)	14 (6%)	16	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	213/248 (86%)	206 (97%)	7 (3%)	33	67
1	G	151/248 (61%)	144 (95%)	7 (5%)	24	57
All	All	588/744 (79%)	560 (95%)	28 (5%)	23	55

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	SER
1	A	46	ARG
1	A	70	GLN
1	A	128	LYS
1	A	142	GLN
1	A	155	LYS
1	A	177	TYR
1	A	187	LYS
1	A	198	LEU
1	A	201	LEU
1	A	247	GLU
1	A	250	GLU
1	A	251	GLU
1	A	269	SER
1	D	46	ARG
1	D	61	SER
1	D	70	GLN
1	D	150	ARG
1	D	155	LYS
1	D	177	TYR
1	D	234	LYS
1	G	2	PRO
1	G	86	GLN
1	G	95	ARG
1	G	130	GLN
1	G	177	TYR
1	G	239	LEU
1	G	252	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	GLN

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Mol	Chain	Res	Type
1	A	139	GLN
1	A	143	GLN
1	A	282	GLN
1	D	70	GLN
1	D	130	GLN
1	G	130	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

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 ⓘ

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	270/295 (91%)	0.07	2 (0%) 84 79	39, 61, 95, 131	0
1	D	260/295 (88%)	0.29	5 (1%) 66 58	41, 75, 122, 142	0
1	G	229/295 (77%)	1.02	28 (12%) 8 7	88, 115, 151, 174	0
2	B	13/13 (100%)	0.32	0 100 100	64, 95, 139, 148	0
2	E	13/13 (100%)	-0.02	0 100 100	68, 90, 111, 122	0
2	H	13/13 (100%)	1.00	1 (7%) 19 16	124, 137, 143, 146	0
3	C	13/13 (100%)	0.14	0 100 100	72, 95, 123, 126	0
3	F	13/13 (100%)	0.28	0 100 100	57, 79, 109, 114	0
3	I	13/13 (100%)	0.42	0 100 100	101, 127, 146, 150	0
All	All	837/963 (86%)	0.42	36 (4%) 40 31	39, 85, 140, 174	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	289	ALA	4.1
1	G	81	MET	3.7
1	G	113	CYS	3.6
1	G	167	ASP	3.4
1	G	114	PHE	3.4
1	A	200	ALA	3.3
2	H	9	DT	3.2
1	G	53	GLY	3.1
1	G	184	TYR	3.1
1	G	86	GLN	3.0
1	G	58	LEU	3.0
1	A	290	PRO	3.0
1	D	275	HIS	2.9
1	D	290	PRO	2.9
1	G	166	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	G	155	LYS	2.8
1	G	271	LEU	2.8
1	D	263	TYR	2.8
1	G	2	PRO	2.8
1	G	240	GLY	2.7
1	G	170	PHE	2.7
1	D	2	PRO	2.6
1	G	87	LEU	2.6
1	G	263	TYR	2.6
1	G	171	PHE	2.5
1	G	49	ALA	2.3
1	G	99	LEU	2.3
1	D	277	ARG	2.1
1	G	120	PHE	2.1
1	G	253	PHE	2.1
1	G	80	GLY	2.1
1	G	75	LEU	2.1
1	G	266	PRO	2.1
1	G	159	ARG	2.1
1	G	27	VAL	2.0
1	G	127	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 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.