



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2026 – 11:12 AM UTC

PDB ID : 2GB7 / pdb_00002gb7
Title : Metal-depleted Ecl18kI in complex with uncleaved, modified DNA
Authors : Bochtler, M.; Szczepanowski, R.H.; Tamulaitis, G.; Grazulis, S.; Czapinska, H.; Manakova, E.; Siksny, V.
Deposited on : 2006-03-10
Resolution : 1.70 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

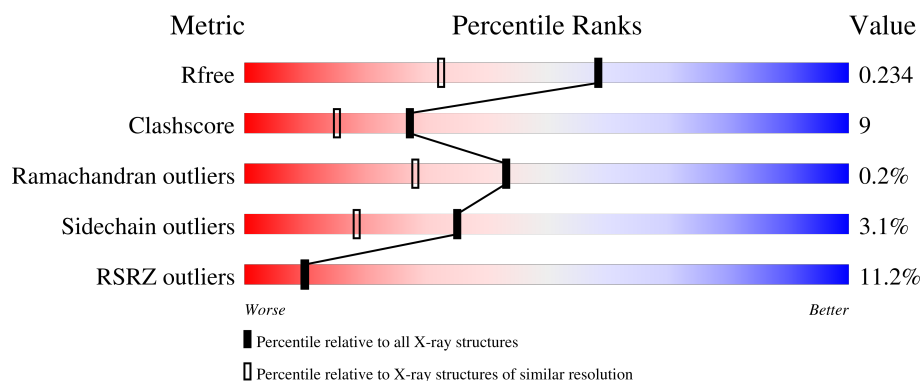
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	E	9	67% 33%
1	G	9	56% 44%
2	F	9	33% 67%
2	H	9	89% 11%
3	A	305	19% 80% 15% .

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Mol	Chain	Length	Quality of chain
3	B	305	8% 80% 16%
3	C	305	8% 78% 15%
3	D	305	10% 80% 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11749 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 DNA chain called DNA STRAND 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	9	Total	C	N	O	P	0	9	0
			364	172	74	102	16			
1	G	9	Total	C	N	O	P	0	9	0
			364	172	74	102	16			

- Molecule 2 is a DNA chain called DNA STRAND 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	9	Total	C	N	O	P	0	9	0
			362	172	68	106	16			
2	H	9	Total	C	N	O	P	0	9	0
			362	172	68	106	16			

- Molecule 3 is a protein called R.Ecl18kI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	293	Total	C	N	O	S	53	0	0
			2421	1537	404	467	13			
3	B	295	Total	C	N	O	S	26	0	0
			2441	1552	405	471	13			
3	C	293	Total	C	N	O	S	30	0	0
			2421	1542	400	466	13			
3	D	292	Total	C	N	O	S	10	0	0
			2412	1537	398	464	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	GLN	ARG	engineered mutation	UNP O87963
B	277	GLN	ARG	engineered mutation	UNP O87963
C	277	GLN	ARG	engineered mutation	UNP O87963
D	277	GLN	ARG	engineered mutation	UNP O87963

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	20	Total 20	O 20	0	0
4	F	26	Total 26	O 26	0	0
4	G	24	Total 24	O 24	0	0
4	H	29	Total 29	O 29	0	0
4	A	83	Total 83	O 83	0	0
4	B	144	Total 144	O 144	0	0
4	C	128	Total 128	O 128	0	0
4	D	148	Total 148	O 148	0	0

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.

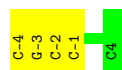
• Molecule 1: DNA STRAND 1

Chain E: 

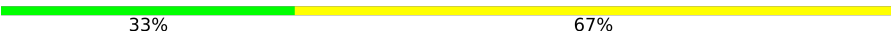


• Molecule 1: DNA STRAND 1

Chain G: 

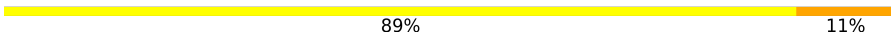


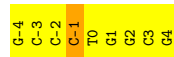
• Molecule 2: DNA STRAND 2

Chain F: 



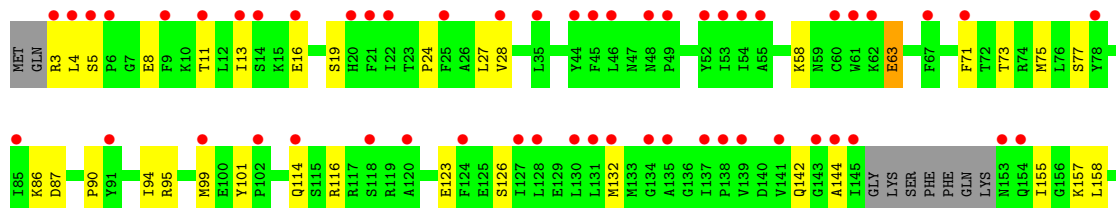
• Molecule 2: DNA STRAND 2

Chain H: 



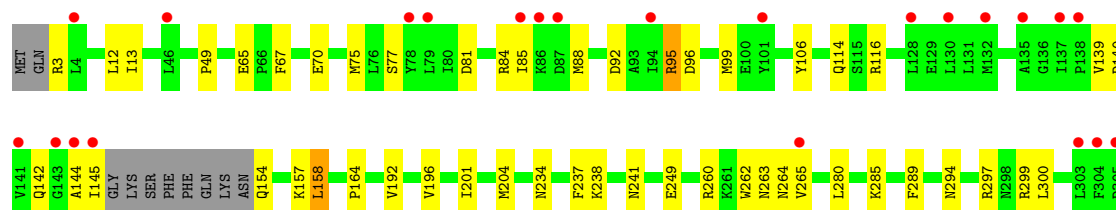
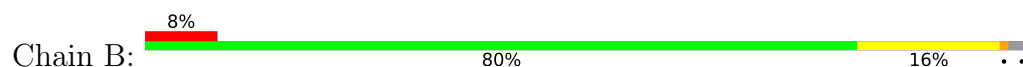
• Molecule 3: R.Ecl18kI

Chain A: 

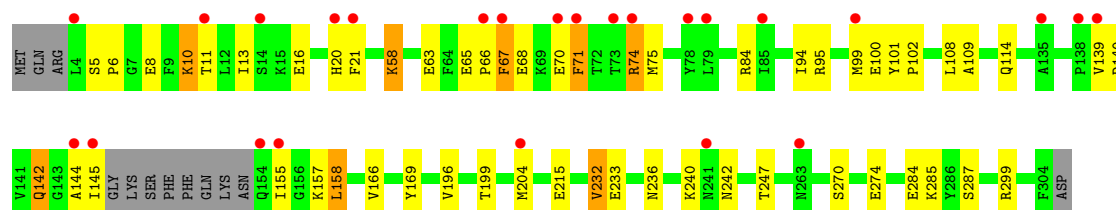
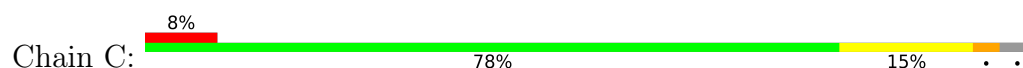




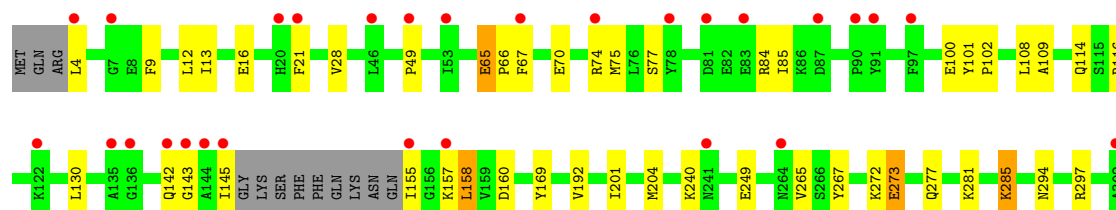
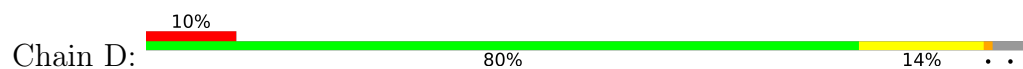
• Molecule 3: R.Ecl18kI



• Molecule 3: R.Ecl18kI



• Molecule 3: R.Ecl18kI



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.70Å 96.74Å 192.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.70 20.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	97.5 (20.00-1.70) 97.5 (20.00-1.70)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.51 (at 1.68Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.207 , 0.236 0.206 , 0.234	Depositor DCC
R_{free} test set	7898 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11749	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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](#)

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	E	0.23	0/408	0.60	0/626
1	G	0.23	0/408	0.69	0/626
2	F	0.24	0/404	0.74	0/620
2	H	0.24	0/404	0.65	0/620
3	A	0.66	0/2468	0.91	2/3330 (0.1%)
3	B	0.70	0/2489	0.88	2/3357 (0.1%)
3	C	0.68	0/2469	0.90	2/3332 (0.1%)
3	D	0.69	0/2460	0.93	0/3320
All	All	0.64	0/11510	0.87	6/15831 (0.0%)

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
2	H	0	1
3	B	0	1
All	All	0	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	289	PHE	CA-C-N	5.94	125.42	119.24
3	A	289	PHE	C-N-CA	5.94	125.42	119.24
3	B	289	PHE	CA-C-N	5.19	125.24	119.32
3	B	289	PHE	C-N-CA	5.19	125.24	119.32
3	C	232	VAL	CA-C-N	5.18	127.53	120.54
3	C	232	VAL	C-N-CA	5.18	127.53	120.54

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	241	ASN	Peptide
2	H	-1[A]	DC	Sidechain

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	E	364	0	185	8	0
1	G	364	0	189	8	0
2	F	362	0	194	15	0
2	H	362	0	195	13	0
3	A	2421	0	2387	32	0
3	B	2441	0	2405	51	0
3	C	2421	0	2388	57	0
3	D	2412	0	2380	37	0
4	A	83	0	0	2	0
4	B	144	0	0	10	0
4	C	128	0	0	10	0
4	D	148	0	0	6	0
4	E	20	0	0	3	0
4	F	26	0	0	1	0
4	G	24	0	0	3	0
4	H	29	0	0	1	0
All	All	11749	0	10323	193	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 9.

All (193) 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)
3:C:67:PHE:HD1	3:C:67:PHE:C	1.48	1.21
1:G:-3[A]:DG:H5''	3:C:157:LYS:HB2	1.32	1.05
3:C:63:GLU:HG3	4:C:418:HOH:O	1.57	1.04
3:C:67:PHE:C	3:C:67:PHE:CD1	2.27	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:77:SER:HB2	4:B:363:HOH:O	1.61	0.97
3:A:27:LEU:HD11	3:A:63:GLU:HG2	1.50	0.91
3:A:13:ILE:HA	3:A:75:MET:HE2	1.52	0.91
1:G:-3[A]:DG:C5'	3:C:157:LYS:HB2	2.00	0.90
3:C:16:GLU:HB2	4:C:403:HOH:O	1.72	0.88
2:H:-3[B]:DC:H5''	3:C:157:LYS:HB2	1.56	0.85
3:D:21:PHE:CE1	3:D:67:PHE:HD2	1.94	0.85
3:D:160:ASP:HB3	4:D:401:HOH:O	1.77	0.84
3:C:21:PHE:CE1	3:C:67:PHE:CZ	2.66	0.83
3:B:140:ASP:HB3	3:B:145:ILE:HD11	1.59	0.82
3:D:240:LYS:HD3	4:D:410:HOH:O	1.78	0.81
2:F:-3[B]:DC:H5''	3:A:157:LYS:HB2	1.62	0.81
3:C:21:PHE:CE1	3:C:67:PHE:CE1	2.70	0.80
3:C:21:PHE:CD1	3:C:67:PHE:CZ	2.70	0.79
3:A:3:ARG:HG2	3:A:4:LEU:H	1.50	0.77
3:C:67:PHE:HD1	3:C:68:GLU:N	1.80	0.77
3:A:27:LEU:CD1	3:A:63:GLU:HG2	2.15	0.77
3:B:70:GLU:HG3	4:B:374:HOH:O	1.87	0.74
1:E:-4[A]:DC:H2''	1:E:-3[A]:DG:OP2	1.86	0.74
3:C:67:PHE:CD1	3:C:68:GLU:N	2.54	0.74
4:G:19:HOH:O	3:D:157:LYS:HG3	1.88	0.73
3:C:13:ILE:HD13	3:C:75:MET:CE	2.19	0.72
3:B:196:VAL:CG2	3:B:204:MET:HE3	2.20	0.72
3:A:197:ASN:HB3	4:B:387:HOH:O	1.90	0.71
3:B:12:LEU:HG	3:B:75:MET:HE1	1.73	0.70
3:A:63:GLU:O	3:A:63:GLU:HG3	1.90	0.70
3:C:13:ILE:HD13	3:C:75:MET:HE3	1.73	0.70
3:C:67:PHE:HD1	3:C:67:PHE:O	1.74	0.68
3:A:270:SER:O	3:A:274:GLU:HG2	1.94	0.67
1:E:-4[B]:DC:H2''	1:E:-3[B]:DG:OP2	1.93	0.67
3:B:13:ILE:CG1	3:B:75:MET:HE2	2.24	0.67
3:D:21:PHE:CE1	3:D:67:PHE:CD2	2.82	0.66
3:D:158:LEU:HD11	3:D:192:VAL:HG22	1.77	0.66
3:C:95:ARG:O	3:C:99:MET:HG3	1.96	0.66
1:E:0[A]:DA:C2	4:E:162:HOH:O	2.48	0.65
3:A:94:ILE:HD12	3:B:85:ILE:HD13	1.77	0.65
2:F:-3[A]:DC:H5''	3:B:157:LYS:HB2	1.79	0.65
2:H:0[A]:DT:H72	4:D:348:HOH:O	1.97	0.64
3:B:294:ASN:HA	3:B:297:ARG:HG2	1.79	0.64
3:B:237:PHE:CD2	3:B:238:LYS:HG2	2.33	0.63
3:D:273:GLU:O	3:D:277:GLN:HG3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:144:ALA:O	3:A:285:LYS:HD3	1.99	0.63
3:B:196:VAL:HG23	3:B:204:MET:HE3	1.81	0.63
3:B:13:ILE:HG12	3:B:75:MET:HE2	1.80	0.62
3:A:77:SER:HB2	4:A:339:HOH:O	2.00	0.62
3:C:75:MET:HE1	4:C:403:HOH:O	2.00	0.62
3:D:84:ARG:NH2	3:D:100:GLU:HG3	2.16	0.61
3:D:77:SER:HB2	4:D:358:HOH:O	2.01	0.60
3:A:90:PRO:HB2	3:B:85:ILE:HG22	1.82	0.60
3:B:85:ILE:HG22	3:B:85:ILE:O	2.00	0.60
3:D:12:LEU:HG	3:D:75:MET:HE1	1.84	0.60
3:B:158:LEU:C	3:B:158:LEU:HD23	2.26	0.60
3:C:144:ALA:HA	3:C:285:LYS:HE2	1.83	0.59
1:G:-4[B]:DC:H4'	4:G:19:HOH:O	2.01	0.59
3:C:11:THR:C	4:C:427:HOH:O	2.45	0.59
3:C:21:PHE:CZ	3:C:67:PHE:CE1	2.90	0.58
3:C:20:HIS:O	3:C:67:PHE:CZ	2.55	0.58
3:A:94:ILE:CD1	3:B:85:ILE:HD13	2.34	0.56
3:B:234:ASN:HB2	4:B:448:HOH:O	2.05	0.56
3:B:88:MET:HE1	3:B:96:ASP:HB2	1.87	0.56
3:B:88:MET:HE2	3:B:92:ASP:HB3	1.87	0.56
3:D:265:VAL:HB	4:D:434:HOH:O	2.05	0.55
3:A:3:ARG:O	3:A:4:LEU:HG	2.06	0.55
3:B:234:ASN:CG	4:B:448:HOH:O	2.49	0.55
3:C:155:ILE:HD11	3:C:169:TYR:CE2	2.41	0.55
3:A:3:ARG:CG	3:A:4:LEU:H	2.18	0.55
3:B:95:ARG:O	3:B:99:MET:HG3	2.06	0.55
3:D:142:GLN:HG3	3:D:143:GLY:N	2.21	0.55
3:D:142:GLN:HG2	3:D:160:ASP:HA	1.88	0.54
3:B:85:ILE:O	3:B:85:ILE:CG2	2.55	0.54
3:A:132:MET:HE1	3:A:252:LEU:HD21	1.90	0.54
2:H:0[B]:DT:H72	4:C:412:HOH:O	2.07	0.53
3:D:84:ARG:NH2	3:D:100:GLU:CG	2.71	0.53
3:C:75:MET:HA	4:C:366:HOH:O	2.08	0.53
3:C:13:ILE:HD13	3:C:75:MET:HE2	1.90	0.53
3:C:139:VAL:O	3:C:299:ARG:NH1	2.42	0.53
3:C:71:PHE:CD2	3:C:71:PHE:C	2.87	0.53
3:C:70:GLU:HG3	4:C:415:HOH:O	2.08	0.52
3:A:86:LYS:O	3:A:87:ASP:HB2	2.08	0.52
3:C:67:PHE:CD1	3:C:67:PHE:O	2.54	0.52
3:B:67:PHE:HA	3:B:70:GLU:HG2	1.91	0.51
2:F:2[A]:DG:O4'	3:A:114:GLN:HG3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:196:VAL:HG22	3:C:204:MET:HE3	1.92	0.51
1:G:-3[B]:DG:H5''	3:D:157:LYS:HB2	1.92	0.50
3:D:142:GLN:HG3	3:D:143:GLY:H	1.76	0.50
1:E:0[A]:DA:H2	4:E:162:HOH:O	1.90	0.50
3:B:13:ILE:HA	3:B:75:MET:HE2	1.93	0.50
3:D:70:GLU:HG3	3:D:74:ARG:HH12	1.76	0.50
1:E:0[B]:DA:H2'	1:E:0[B]:DA:N3	2.27	0.50
2:F:-3[A]:DC:H5''	3:B:157:LYS:CB	2.41	0.50
3:C:236:ASN:O	3:C:240:LYS:HE2	2.11	0.50
3:D:155:ILE:HD11	3:D:169:TYR:CE2	2.47	0.50
3:B:234:ASN:HB2	4:B:382:HOH:O	2.12	0.50
3:A:3:ARG:HG2	3:A:4:LEU:N	2.24	0.50
3:B:234:ASN:CB	4:B:448:HOH:O	2.60	0.49
3:C:20:HIS:O	3:C:67:PHE:CE2	2.65	0.49
3:C:232:VAL:HG22	3:C:247:THR:HG21	1.94	0.49
3:D:84:ARG:CZ	3:D:100:GLU:HG3	2.42	0.49
3:B:237:PHE:HD2	3:B:238:LYS:HG2	1.77	0.49
3:B:280:LEU:HD21	3:B:300:LEU:HD21	1.95	0.48
2:H:-2[A]:DC:H2'	2:H:-1[A]:DC:C6	2.47	0.48
3:C:84:ARG:CZ	3:C:100:GLU:OE2	2.62	0.48
3:C:196:VAL:CG2	3:C:204:MET:HE3	2.44	0.48
3:B:144:ALA:HA	3:B:285:LYS:HG2	1.97	0.47
3:C:13:ILE:HD12	3:C:108:LEU:HD13	1.96	0.47
2:F:2[B]:DG:O4'	3:B:114:GLN:HG3	2.13	0.47
3:C:199:THR:HG21	4:C:426:HOH:O	2.13	0.47
2:F:3[A]:DC:H2''	2:F:4[A]:DG:C8	2.49	0.47
3:B:294:ASN:OD1	4:B:412:HOH:O	2.20	0.47
2:F:0[A]:DT:H3'	3:B:116:ARG:HB3	1.97	0.47
2:H:2[B]:DG:N3	3:D:114:GLN:HB3	2.30	0.47
3:C:158:LEU:C	3:C:158:LEU:HD23	2.39	0.47
3:D:201:ILE:HD13	3:D:204:MET:CE	2.45	0.47
3:B:234:ASN:ND2	4:B:448:HOH:O	2.47	0.47
1:G:-4[A]:DC:H1'	1:G:-3[A]:DG:C8	2.49	0.46
3:A:16:GLU:HB3	3:A:71:PHE:HE1	1.80	0.46
3:C:270:SER:O	3:C:274:GLU:HG3	2.14	0.46
2:F:-3[B]:DC:C5'	3:A:157:LYS:HB2	2.41	0.46
3:D:70:GLU:HB3	4:D:435:HOH:O	2.15	0.45
3:C:101:TYR:N	3:C:102:PRO:CD	2.79	0.45
1:E:0[A]:DA:OP1	3:A:184:THR:HG23	2.16	0.45
3:C:74:ARG:C	4:C:366:HOH:O	2.58	0.45
3:C:94:ILE:HD12	3:D:85:ILE:HD13	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:294:ASN:HA	3:D:297:ARG:HG2	1.98	0.45
2:F:0[B]:DT:OP2	2:F:0[B]:DT:H4'	2.16	0.45
3:D:9:PHE:O	3:D:13:ILE:HG13	2.16	0.45
1:G:-2[B]:DC:H2'	1:G:-1[B]:DC:C6	2.52	0.45
3:A:95:ARG:O	3:A:99:MET:HG3	2.17	0.45
4:E:307:HOH:O	2:F:3[B]:DC:H5	1.99	0.45
3:B:237:PHE:CE2	3:B:238:LYS:HE3	2.52	0.44
2:H:-4[A]:DG:H2''	2:H:-3[A]:DC:OP2	2.16	0.44
3:B:139:VAL:O	3:B:299:ARG:NH1	2.49	0.44
1:G:-4[B]:DC:H1'	1:G:-3[B]:DG:C8	2.53	0.44
4:G:8:HOH:O	2:H:0[A]:DT:H6	1.99	0.44
3:C:140:ASP:OD2	3:C:166:VAL:HG22	2.17	0.44
2:H:3[A]:DC:H2''	2:H:4[A]:DG:C8	2.52	0.44
3:B:81:ASP:CG	3:B:84:ARG:HB2	2.43	0.44
3:D:101:TYR:N	3:D:102:PRO:CD	2.80	0.44
1:E:4[B]:DC:O2	2:F:4[B]:DG:N2	2.43	0.44
4:H:14:HOH:O	3:C:10:LYS:HE3	2.16	0.44
3:A:24:PRO:HB3	3:A:123:GLU:HG3	1.98	0.44
3:C:65:GLU:O	3:C:66:PRO:C	2.59	0.44
3:B:164:PRO:HG3	3:B:262:TRP:CD2	2.52	0.43
1:E:0[A]:DA:N3	1:E:0[A]:DA:H2'	2.34	0.43
2:H:2[A]:DG:O4'	3:C:114:GLN:HG3	2.18	0.43
3:C:21:PHE:CZ	3:C:67:PHE:CZ	3.07	0.43
3:D:155:ILE:HD11	3:D:169:TYR:CD2	2.54	0.43
3:B:201:ILE:HD13	3:B:204:MET:HE2	1.99	0.43
3:B:260:ARG:HA	3:B:263:ASN:HD22	1.82	0.43
3:C:84:ARG:NH2	3:C:100:GLU:OE2	2.52	0.43
2:H:0[A]:DT:H3'	3:D:116:ARG:HB3	2.00	0.43
3:C:5:SER:OG	3:C:8:GLU:HG3	2.19	0.43
3:C:75:MET:CE	4:C:403:HOH:O	2.63	0.43
3:A:90:PRO:HG3	3:B:88:MET:O	2.19	0.42
3:C:58:LYS:HB3	3:C:58:LYS:HE3	1.66	0.42
3:C:142:GLN:NE2	3:C:155:ILE:HG22	2.34	0.42
3:D:13:ILE:HA	3:D:75:MET:HE2	2.01	0.42
2:H:2[B]:DG:O4'	3:D:114:GLN:HG3	2.20	0.42
3:D:49:PRO:HD2	3:D:249:GLU:HG3	2.02	0.42
3:A:27:LEU:HD11	3:A:63:GLU:CG	2.36	0.42
3:D:28:VAL:HG12	3:D:130:LEU:HD12	2.02	0.42
3:A:73:THR:HG21	3:B:99:MET:HG2	2.02	0.42
3:A:294:ASN:HB2	4:A:369:HOH:O	2.18	0.42
3:B:196:VAL:HG23	3:B:204:MET:CE	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:5:SER:OG	3:A:8:GLU:HB2	2.19	0.42
3:A:28:VAL:HG21	3:A:126:SER:HB2	2.02	0.42
3:C:215:GLU:OE1	3:C:242:ASN:ND2	2.52	0.42
3:B:65:GLU:HG3	4:B:442:HOH:O	2.19	0.42
2:F:1[B]:DG:OP2	3:B:106:TYR:OH	2.26	0.42
3:B:262:TRP:O	3:B:265:VAL:HG12	2.19	0.42
2:H:0[A]:DT:OP2	2:H:1[A]:DG:H2'	2.20	0.42
2:H:3[B]:DC:H2''	2:H:4[B]:DG:C8	2.54	0.42
2:F:2[B]:DG:N3	3:B:114:GLN:HB3	2.35	0.41
3:D:267:TYR:O	3:D:272:LYS:HE3	2.21	0.41
3:D:281:LYS:HG3	3:D:285:LYS:NZ	2.35	0.41
3:B:196:VAL:CG2	3:B:204:MET:CE	2.96	0.41
3:C:5:SER:HB2	3:C:6:PRO:HD2	2.02	0.41
3:C:67:PHE:HD1	3:C:68:GLU:CA	2.33	0.41
2:F:-3[B]:DC:N4	4:F:218:HOH:O	2.53	0.41
3:B:158:LEU:HD11	3:B:192:VAL:HG22	2.02	0.41
3:C:67:PHE:CD1	3:C:68:GLU:HA	2.56	0.41
3:D:13:ILE:HD13	3:D:108:LEU:HA	2.03	0.41
1:G:-3[A]:DG:H5'	3:C:157:LYS:HB2	1.95	0.40
3:A:192:VAL:HB	3:A:193:PRO:HD3	2.04	0.40
3:B:49:PRO:HD2	3:B:249:GLU:HG3	2.04	0.40
3:B:280:LEU:CD2	3:B:300:LEU:HD21	2.51	0.40
2:F:0[B]:DT:H3'	3:A:116:ARG:HB3	2.03	0.40
3:C:109:ALA:CB	3:D:109:ALA:HB3	2.52	0.40
3:D:65:GLU:N	3:D:66:PRO:HD2	2.37	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
3	A	289/305 (95%)	283 (98%)	5 (2%)	1 (0%)	36 22

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	B	291/305 (95%)	285 (98%)	6 (2%)	0	100	100
3	C	289/305 (95%)	284 (98%)	4 (1%)	1 (0%)	36	22
3	D	288/305 (94%)	285 (99%)	3 (1%)	0	100	100
All	All	1157/1220 (95%)	1137 (98%)	18 (2%)	2 (0%)	43	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	74	ARG
3	A	101	TYR

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	
3	A	272/283 (96%)	261 (96%)	11 (4%)	28	12
3	B	274/283 (97%)	268 (98%)	6 (2%)	45	29
3	C	272/283 (96%)	262 (96%)	10 (4%)	30	14
3	D	271/283 (96%)	264 (97%)	7 (3%)	40	23
All	All	1089/1132 (96%)	1055 (97%)	34 (3%)	35	18

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	11	THR
3	A	19	SER
3	A	58	LYS
3	A	63	GLU
3	A	142	GLN
3	A	155	ILE
3	A	158	LEU
3	A	263	ASN
3	A	265	VAL

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Mol	Chain	Res	Type
3	A	280	LEU
3	A	281	LYS
3	B	3	ARG
3	B	95	ARG
3	B	142	GLN
3	B	154	GLN
3	B	158	LEU
3	B	264	ASN
3	C	10	LYS
3	C	58	LYS
3	C	67	PHE
3	C	71	PHE
3	C	142	GLN
3	C	145	ILE
3	C	158	LEU
3	C	233	GLU
3	C	284	GLU
3	C	287	SER
3	D	4	LEU
3	D	16	GLU
3	D	65	GLU
3	D	145	ILE
3	D	158	LEU
3	D	273	GLU
3	D	285	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	190	GLN
3	A	253	GLN
3	A	263	ASN
3	A	276	GLN
3	B	39	GLN
3	B	59	ASN
3	B	197	ASN
3	B	263	ASN
3	B	264	ASN
3	C	142	GLN
3	C	276	GLN
3	D	253	GLN
3	D	263	ASN

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Mol	Chain	Res	Type
3	D	276	GLN
3	D	277	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(Å²)	Q<0.9	
1	E	9/9 (100%)	0.21	0	100	100	6, 7, 12, 17	9 (100%)
1	G	9/9 (100%)	0.14	0	100	100	3, 5, 9, 12	9 (100%)
2	F	9/9 (100%)	0.23	0	100	100	4, 7, 14, 16	9 (100%)
2	H	9/9 (100%)	-0.07	0	100	100	3, 5, 10, 11	9 (100%)
3	A	293/305 (96%)	1.07	57 (19%)	3	2	8, 23, 41, 51	19 (6%)
3	B	295/305 (96%)	0.43	23 (7%)	19	20	7, 15, 29, 43	19 (6%)
3	C	293/305 (96%)	0.55	25 (8%)	16	17	6, 14, 31, 40	26 (8%)
3	D	292/305 (95%)	0.57	30 (10%)	12	12	5, 15, 31, 51	17 (5%)
All	All	1209/1256 (96%)	0.64	135 (11%)	10	10	3, 16, 34, 51	117 (9%)

All (135) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	145	ILE	6.1
3	C	67	PHE	6.1
3	A	67	PHE	5.2
3	A	4	LEU	4.8
3	C	78	TYR	4.5
3	A	20	HIS	4.5
3	D	67	PHE	4.5
3	C	66	PRO	4.3
3	A	78	TYR	4.3
3	C	154	GLN	3.8
3	A	135	ALA	3.7
3	B	144	ALA	3.5
3	C	85	ILE	3.5
3	C	79	LEU	3.4
3	A	138	PRO	3.4
3	C	241	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
3	C	144	ALA	3.4
3	B	85	ILE	3.3
3	A	53	ILE	3.2
3	C	145	ILE	3.2
3	A	21	PHE	3.2
3	C	155	ILE	3.2
3	A	3	ARG	3.2
3	B	305	ASP	3.2
3	D	145	ILE	3.1
3	B	143	GLY	3.1
3	D	144	ALA	3.1
3	A	49	PRO	3.1
3	A	46	LEU	3.1
3	D	155	ILE	3.0
3	A	9	PHE	3.0
3	A	61	TRP	3.0
3	B	46	LEU	2.9
3	A	6	PRO	2.9
3	D	83	GLU	2.9
3	B	141	VAL	2.9
3	A	128	LEU	2.9
3	D	46	LEU	2.9
3	C	138	PRO	2.9
3	A	153	ASN	2.9
3	D	4	LEU	2.9
3	A	13	ILE	2.8
3	D	20	HIS	2.8
3	C	139	VAL	2.8
3	D	97	PHE	2.7
3	A	132	MET	2.7
3	B	304	PHE	2.6
3	D	21	PHE	2.6
3	B	94	ILE	2.6
3	A	270	SER	2.6
3	A	131	LEU	2.6
3	A	139	VAL	2.6
3	C	11	THR	2.6
3	C	74	ARG	2.6
3	B	303	LEU	2.6
3	A	5	SER	2.5
3	B	101	TYR	2.5
3	A	120	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
3	C	20	HIS	2.5
3	A	44	TYR	2.5
3	C	73	THR	2.5
3	C	4	LEU	2.5
3	D	81	ASP	2.5
3	A	91	TYR	2.5
3	D	78	TYR	2.5
3	D	87	ASP	2.4
3	A	143	GLY	2.4
3	C	99	MET	2.4
3	D	90	PRO	2.4
3	B	78	TYR	2.4
3	A	85	ILE	2.4
3	B	86	LYS	2.4
3	D	157	LYS	2.4
3	D	143	GLY	2.4
3	A	102	PRO	2.4
3	A	130	LEU	2.4
3	B	135	ALA	2.4
3	D	91	TYR	2.4
3	C	14	SER	2.4
3	C	71	PHE	2.4
3	B	87	ASP	2.4
3	A	141	VAL	2.4
3	B	265	VAL	2.4
3	A	154	GLN	2.3
3	D	264	ASN	2.3
3	A	144	ALA	2.3
3	B	145	ILE	2.3
3	D	7	GLY	2.3
3	A	55	ALA	2.3
3	A	35	LEU	2.3
3	D	53	ILE	2.3
3	B	138	PRO	2.3
3	B	4	LEU	2.3
3	A	16	GLU	2.3
3	A	118	SER	2.3
3	A	62	LYS	2.3
3	A	134	GLY	2.3
3	A	60	CYS	2.2
3	C	135	ALA	2.2
3	D	302	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
3	A	25	PHE	2.2
3	A	45	PHE	2.2
3	A	273	GLU	2.2
3	C	70	GLU	2.2
3	A	71	PHE	2.2
3	A	124	PHE	2.2
3	D	74	ARG	2.2
3	D	136	GLY	2.2
3	A	22	ILE	2.2
3	A	137	ILE	2.2
3	A	11	THR	2.2
3	A	14	SER	2.2
3	B	132	MET	2.1
3	D	142	GLN	2.1
3	D	135	ALA	2.1
3	C	263	ASN	2.1
3	A	114	GLN	2.1
3	B	128	LEU	2.1
3	B	130	LEU	2.1
3	C	21	PHE	2.1
3	A	28	VAL	2.1
3	A	48	ASN	2.1
3	C	204	MET	2.1
3	B	79	LEU	2.1
3	D	303	LEU	2.1
3	D	49	PRO	2.1
3	D	122	LYS	2.1
3	D	304	PHE	2.1
3	A	54	ILE	2.1
3	A	127	ILE	2.1
3	B	137	ILE	2.1
3	A	99	MET	2.0
3	A	264	ASN	2.0
3	D	241	ASN	2.0
3	A	52	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.