



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:01 AM UTC

PDB ID : 1LQG / pdb_00001lqg
Title : ESCHERICHIA COLI URACIL-DNA GLYCOSYLASE COMPLEX WITH
URACIL-DNA GLYCOSYLASE INHIBITOR PROTEIN
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Deposited on : 2002-05-10
Resolution : 2.90 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

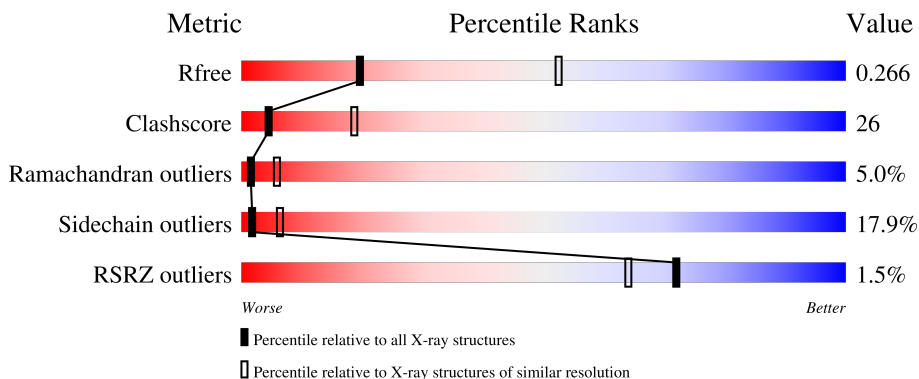
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	229	% 45% 40% 13% ..
1	B	229	3% 47% 37% 11% ..
2	C	84	% 42% 39% 6% 13%
2	D	84	43% 39% . 17%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4765 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 URACIL-DNA GLYCOSYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	0	0
			1755	1128	312	312	3			
1	B	224	Total	C	N	O	S	0	0	0
			1749	1127	311	308	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	cloning artifact	UNP P12295
B	1	MET	-	cloning artifact	UNP P12295

- Molecule 2 is a protein called URACIL-DNA GLYCOSYLASE INHIBITOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	73	Total	C	N	O	S	0	0	0
			567	358	88	118	3			
2	D	70	Total	C	N	O	S	0	0	0
			538	341	84	110	3			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	71	Total	O	0	0
			71	71		
3	B	51	Total	O	0	0
			51	51		
3	C	23	Total	O	0	0
			23	23		
3	D	11	Total	O	0	0
			11	11		

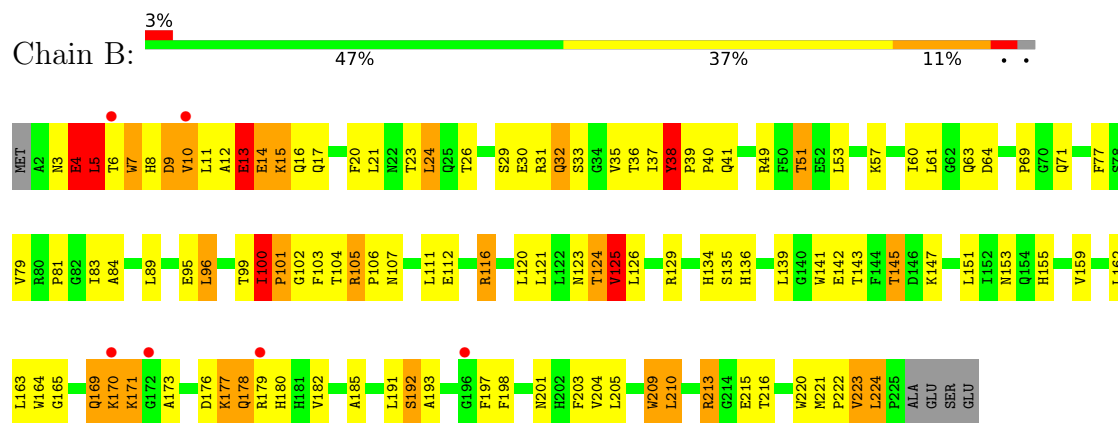
3 Residue-property plots [i](#)

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.

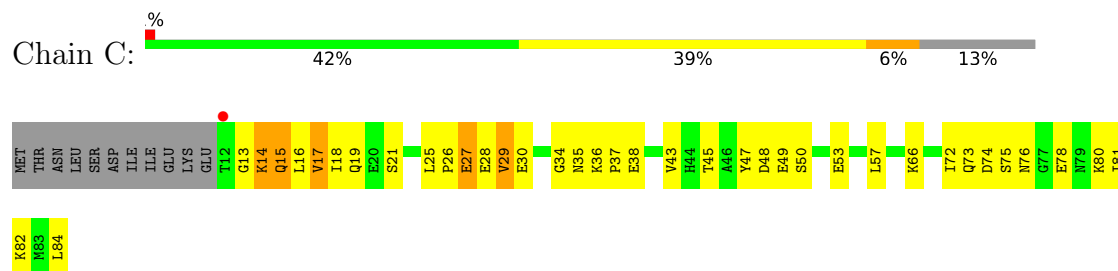
• Molecule 1: URACIL-DNA GLYCOSYLASE



• Molecule 1: URACIL-DNA GLYCOSYLASE

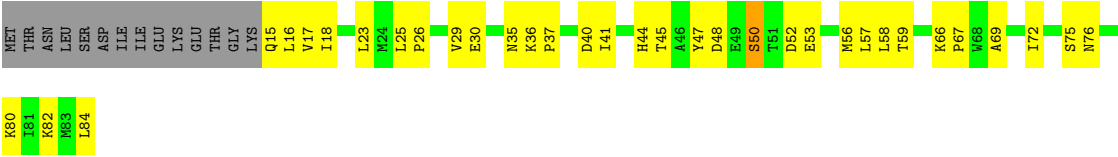


• Molecule 2: URACIL-DNA GLYCOSYLASE INHIBITOR



• Molecule 2: URACIL-DNA GLYCOSYLASE INHIBITOR

Chain D: 43% 39% • 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.16Å 89.90Å 141.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 20.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	93.6 (20.00-2.90) 93.2 (20.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.22 (at 2.79Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.183 , 0.277 0.177 , 0.266	Depositor DCC
R_{free} test set	1080 reflections (6.99%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 99.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4765	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% 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	0.74	1/1811 (0.1%)	1.23	27/2475 (1.1%)
1	B	0.67	0/1805	1.23	17/2466 (0.7%)
2	C	0.74	0/575	1.27	6/780 (0.8%)
2	D	0.69	0/546	1.08	4/743 (0.5%)
All	All	0.71	1/4737 (0.0%)	1.22	54/6464 (0.8%)

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
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	125	VAL	CA-CB	5.00	1.60	1.54

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	ILE	CA-C-N	11.09	133.71	119.84
1	B	100	ILE	C-N-CA	11.09	133.71	119.84
1	B	8	HIS	N-CA-C	-10.66	100.23	113.02
2	C	50	SER	N-CA-C	-9.58	99.01	112.13
1	B	38	TYR	N-CA-C	9.43	130.65	109.81
1	B	14	GLU	N-CA-C	-9.21	101.19	111.14
1	B	38	TYR	CA-C-N	-8.83	111.28	120.38
1	B	38	TYR	C-N-CA	-8.83	111.28	120.38
1	A	38	TYR	CA-C-N	-8.78	111.34	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	38	TYR	C-N-CA	-8.78	111.34	120.38
1	B	7	TRP	N-CA-C	-8.66	99.04	112.99
1	A	38	TYR	N-CA-C	8.46	128.50	109.81
2	C	14	LYS	N-CA-C	8.34	122.07	108.55
1	A	37	ILE	N-CA-C	-8.02	101.31	110.05
1	B	23	THR	N-CA-C	-7.84	102.73	111.82
1	A	105	ARG	N-CA-C	7.75	119.54	109.93
1	B	126	LEU	N-CA-C	7.68	122.47	113.18
1	A	85	ILE	CA-C-N	7.58	125.12	119.66
1	A	85	ILE	C-N-CA	7.58	125.12	119.66
1	B	182	VAL	N-CA-C	7.11	116.56	106.53
1	B	125	VAL	N-CA-C	-7.04	97.99	108.12
1	A	5	LEU	N-CA-C	7.01	121.24	111.56
1	B	155	HIS	N-CA-C	6.93	121.34	113.02
1	A	126	LEU	N-CA-C	6.76	121.28	112.89
1	A	65	PRO	N-CA-C	6.68	120.57	111.22
1	A	15	LYS	N-CA-C	-6.67	104.50	114.64
1	A	12	ALA	N-CA-C	-6.36	105.67	113.50
1	B	111	LEU	N-CA-C	6.33	119.94	112.72
1	A	80	ARG	CA-C-N	6.29	127.70	119.84
1	A	80	ARG	C-N-CA	6.29	127.70	119.84
1	A	8	HIS	N-CA-C	-6.23	97.54	110.80
1	B	123	ASN	N-CA-C	-6.14	100.56	110.32
2	C	17	VAL	N-CA-C	6.13	116.33	108.12
2	C	34	GLY	N-CA-C	-6.13	106.62	113.79
2	C	16	LEU	N-CA-C	-5.99	95.31	107.41
1	A	17	GLN	CA-C-N	5.80	125.93	119.32
1	A	17	GLN	C-N-CA	5.80	125.93	119.32
1	A	179	ARG	N-CA-C	-5.75	105.15	111.82
1	A	197	PHE	N-CA-C	5.69	118.58	111.24
2	C	45	THR	N-CA-C	5.57	118.03	109.07
1	A	117	GLN	N-CA-C	-5.55	105.86	113.30
1	B	13	GLU	N-CA-C	-5.48	99.12	110.80
1	A	182	VAL	N-CA-C	5.48	115.48	107.37
1	A	16	GLN	N-CA-C	-5.36	107.11	113.97
1	A	125	VAL	N-CA-C	-5.33	100.65	108.11
1	A	13	GLU	N-CA-C	-5.31	106.86	113.55
2	D	75	SER	N-CA-C	5.24	116.99	111.28
1	A	89	LEU	N-CA-C	-5.22	105.66	112.23
2	D	66	LYS	CA-C-N	5.20	125.50	119.93
2	D	66	LYS	C-N-CA	5.20	125.50	119.93
1	B	89	LEU	N-CA-C	-5.14	105.59	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	17	VAL	N-CA-C	5.06	115.05	107.51
1	A	221	MET	CA-C-N	5.06	124.97	119.76
1	A	221	MET	C-N-CA	5.06	124.97	119.76

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	38	TYR	Sidechain
1	B	38	TYR	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	A	1755	0	1671	111	0
1	B	1749	0	1682	101	0
2	C	567	0	555	18	0
2	D	538	0	528	14	0
3	A	71	0	0	1	0
3	B	51	0	0	2	0
3	C	23	0	0	1	0
3	D	11	0	0	1	0
All	All	4765	0	4436	238	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 26.

All (238) 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:A:201:ASN:HB3	1:A:204:VAL:HG12	1.43	1.01
1:B:191:LEU:HG	1:B:191:LEU:O	1.68	0.90
1:A:201:ASN:HB3	1:A:204:VAL:CG1	2.03	0.87
1:A:227:GLU:O	1:A:228:SER:HB2	1.75	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:THR:HG23	1:B:105:ARG:H	1.39	0.87
1:B:165:GLY:O	1:B:169:GLN:HG2	1.77	0.85
1:B:11:LEU:O	1:B:15:LYS:HD3	1.78	0.83
1:A:173:ALA:O	1:A:174:ILE:HD13	1.80	0.81
1:A:157:GLU:OE1	1:A:213:ARG:NH2	2.14	0.81
1:B:163:LEU:HB3	1:B:169:GLN:HA	1.63	0.81
1:A:194:HIS:HB2	3:A:282:HOH:O	1.82	0.79
1:A:124:THR:HA	1:A:145:THR:HG21	1.65	0.77
1:B:223:VAL:HG13	1:B:224:LEU:N	1.99	0.77
1:B:38:TYR:O	1:B:39:PRO:C	2.27	0.76
1:A:64:ASP:HB3	1:A:125:VAL:HG13	1.66	0.76
1:A:173:ALA:C	1:A:174:ILE:HD13	2.11	0.75
1:A:6:THR:HG23	1:A:52:GLU:OE1	1.87	0.74
1:A:6:THR:HB	1:A:8:HIS:CD2	2.21	0.74
1:B:3:ASN:O	1:B:4:GLU:HB2	1.85	0.74
1:A:30:GLU:O	1:A:35:VAL:HG13	1.87	0.73
1:A:9:ASP:C	1:A:11:LEU:H	1.97	0.73
1:A:201:ASN:CB	1:A:204:VAL:HG12	2.20	0.71
1:A:67:HIS:HB2	1:A:131:GLY:H	1.56	0.70
1:B:7:TRP:O	1:B:11:LEU:HB2	1.93	0.69
1:B:159:VAL:H	1:B:180:HIS:HD2	1.39	0.69
1:B:4:GLU:O	1:B:5:LEU:HB2	1.91	0.68
1:B:191:LEU:HB2	2:D:56:MET:SD	2.33	0.68
1:B:153:ASN:HA	1:B:180:HIS:CE1	2.29	0.68
1:A:8:HIS:ND1	1:A:8:HIS:C	2.52	0.68
1:A:153:ASN:ND2	1:A:175:ILE:HG23	2.10	0.67
1:A:163:LEU:HB3	1:A:169:GLN:HA	1.78	0.65
1:A:119:VAL:HG22	1:A:218:ILE:HD13	1.77	0.65
1:A:100:ILE:HG23	1:A:103:PHE:HB2	1.79	0.65
1:A:186:PRO:HG2	1:A:196:GLY:HA3	1.79	0.64
1:B:112:GLU:OE2	1:B:116:ARG:HD2	1.97	0.64
1:A:170:LYS:O	1:A:171:LYS:C	2.37	0.63
1:B:11:LEU:O	1:B:15:LYS:HB2	1.98	0.63
2:C:30:GLU:HA	2:C:35:ASN:O	1.98	0.63
1:A:201:ASN:O	1:A:204:VAL:HG12	1.99	0.63
1:A:99:THR:O	1:A:101:PRO:HD3	1.99	0.63
1:B:223:VAL:CG1	1:B:224:LEU:N	2.62	0.63
1:B:5:LEU:HD23	1:B:53:LEU:HG	1.80	0.62
1:B:210:LEU:O	1:B:213:ARG:HG3	1.99	0.62
1:A:10:VAL:C	1:A:12:ALA:H	2.06	0.62
1:B:40:PRO:O	1:B:41:GLN:C	2.41	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:GLU:HB3	1:A:37:ILE:HD11	1.81	0.62
1:A:6:THR:CG2	1:A:7:TRP:N	2.63	0.62
1:B:124:THR:HG22	1:B:145:THR:HG21	1.82	0.62
1:B:213:ARG:HG2	1:B:213:ARG:HH11	1.64	0.62
1:B:99:THR:O	1:B:101:PRO:HD3	2.00	0.61
1:B:107:ASN:O	1:B:107:ASN:ND2	2.33	0.61
1:B:38:TYR:O	1:B:40:PRO:N	2.34	0.61
1:A:95:GLU:OE2	1:A:200:CYS:O	2.19	0.61
1:A:176:ASP:HB3	1:A:179:ARG:HB2	1.82	0.61
1:A:30:GLU:CB	1:A:37:ILE:HD11	2.31	0.61
1:B:9:ASP:C	1:B:11:LEU:H	2.09	0.60
2:D:26:PRO:HA	2:D:37:PRO:HG2	1.82	0.60
1:A:153:ASN:HD22	1:A:175:ILE:HG23	1.65	0.60
1:A:201:ASN:HB2	1:A:205:LEU:HD13	1.82	0.60
1:A:6:THR:HG22	1:A:7:TRP:N	2.14	0.60
1:B:96:LEU:HD13	1:B:203:PHE:CD2	2.36	0.60
1:B:37:ILE:HD13	1:B:136:HIS:HE1	1.66	0.60
1:A:96:LEU:HA	1:A:99:THR:HG22	1.85	0.59
1:A:99:THR:HG23	1:A:100:ILE:N	2.18	0.59
1:A:67:HIS:HB2	1:A:131:GLY:N	2.17	0.59
1:B:14:GLU:HG3	1:B:147:LYS:HG2	1.83	0.59
1:B:95:GLU:O	1:B:99:THR:HG22	2.03	0.59
1:A:170:LYS:O	1:A:172:GLY:N	2.37	0.58
1:A:9:ASP:C	1:A:11:LEU:N	2.62	0.58
1:B:163:LEU:HB3	1:B:169:GLN:CA	2.33	0.57
1:A:153:ASN:HD21	1:A:176:ASP:H	1.52	0.57
1:A:9:ASP:O	1:A:11:LEU:N	2.36	0.57
1:B:95:GLU:OE2	1:B:198:PHE:O	2.22	0.57
2:D:18:ILE:HD12	2:D:44:HIS:HB3	1.87	0.57
1:A:154:GLN:O	1:A:179:ARG:NH1	2.36	0.57
2:D:23:LEU:HD11	2:D:40:ASP:HB3	1.86	0.57
1:A:159:VAL:H	1:A:180:HIS:HD2	1.51	0.56
1:A:41:GLN:NE2	1:A:44:VAL:HG21	2.20	0.56
1:A:153:ASN:HA	1:A:180:HIS:CE1	2.40	0.56
1:B:104:THR:HG23	1:B:105:ARG:N	2.14	0.56
1:A:187:HIS:CG	1:A:188:PRO:HD2	2.41	0.56
1:B:49:ARG:HG2	1:B:49:ARG:HH11	1.71	0.56
1:A:14:GLU:HG3	1:A:147:LYS:HG2	1.87	0.55
1:A:23:THR:O	1:A:26:THR:HG22	2.07	0.55
1:A:175:ILE:HG22	1:A:176:ASP:N	2.22	0.55
1:A:169:GLN:O	1:A:170:LYS:C	2.50	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:ASP:HB3	1:A:125:VAL:CG1	2.37	0.54
2:C:84:LEU:OXT	3:C:94:HOH:O	2.18	0.54
1:B:147:LYS:HE3	1:B:151:LEU:HD11	1.89	0.54
1:A:187:HIS:ND1	1:A:188:PRO:HD2	2.23	0.54
1:B:223:VAL:HG13	1:B:224:LEU:H	1.69	0.54
1:B:147:LYS:HG3	1:B:151:LEU:HD11	1.90	0.54
1:B:165:GLY:C	1:B:169:GLN:HG2	2.33	0.53
1:B:69:PRO:HD2	3:B:262:HOH:O	2.08	0.53
1:B:176:ASP:C	1:B:178:GLN:H	2.17	0.53
2:C:73:GLN:HA	2:C:78:GLU:O	2.09	0.53
1:A:223:VAL:HG13	1:A:224:LEU:N	2.23	0.52
1:B:10:VAL:C	1:B:12:ALA:H	2.17	0.52
1:A:6:THR:CG2	1:A:7:TRP:H	2.21	0.52
1:B:32:GLN:HB2	3:B:243:HOH:O	2.09	0.52
1:A:38:TYR:O	1:A:39:PRO:C	2.46	0.52
1:A:58:VAL:HG12	1:A:59:VAL:N	2.24	0.52
1:A:64:ASP:HB2	1:A:65:PRO:HD2	1.92	0.52
1:B:3:ASN:O	1:B:4:GLU:CB	2.55	0.52
1:B:147:LYS:O	1:B:151:LEU:CD1	2.58	0.52
1:B:147:LYS:O	1:B:151:LEU:HD12	2.10	0.52
1:B:10:VAL:CB	1:B:151:LEU:HD21	2.40	0.52
1:A:179:ARG:HH22	1:B:14:GLU:CD	2.16	0.52
2:C:49:GLU:CG	2:C:49:GLU:O	2.57	0.51
1:A:10:VAL:C	1:A:12:ALA:N	2.67	0.51
1:A:5:LEU:H	1:A:5:LEU:HD23	1.76	0.51
1:B:176:ASP:O	1:B:178:GLN:N	2.44	0.51
1:B:106:PRO:HG3	1:B:222:PRO:HG2	1.93	0.51
1:B:9:ASP:O	1:B:11:LEU:N	2.41	0.50
1:B:60:ILE:HB	1:B:121:LEU:HD23	1.93	0.50
1:A:153:ASN:ND2	1:A:176:ASP:H	2.08	0.50
1:B:83:ILE:CG2	1:B:84:ALA:N	2.74	0.50
1:A:201:ASN:HB3	1:A:204:VAL:HG11	1.91	0.50
1:B:104:THR:HG23	1:B:221:MET:HE1	1.94	0.50
1:A:177:LYS:O	1:A:180:HIS:O	2.30	0.50
1:A:179:ARG:NH2	1:B:14:GLU:OE1	2.45	0.50
1:A:3:ASN:O	1:A:4:GLU:C	2.54	0.50
1:A:99:THR:CG2	1:A:100:ILE:N	2.74	0.49
1:B:10:VAL:C	1:B:12:ALA:N	2.66	0.49
1:A:58:VAL:CG1	1:A:59:VAL:N	2.74	0.49
1:B:30:GLU:HB2	1:B:37:ILE:HD11	1.94	0.49
1:B:163:LEU:HD13	1:B:169:GLN:HA	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:25:LEU:O	2:D:29:VAL:HG23	2.12	0.49
1:B:125:VAL:HA	1:B:141:TRP:HB3	1.94	0.49
2:D:47:TYR:HE2	3:D:94:HOH:O	1.95	0.49
2:C:74:ASP:OD2	2:C:78:GLU:HB3	2.11	0.49
1:B:201:ASN:O	1:B:204:VAL:HG12	2.13	0.48
2:C:49:GLU:O	2:C:49:GLU:HG3	2.14	0.48
1:A:30:GLU:CD	1:A:129:ARG:NH1	2.71	0.48
1:B:147:LYS:HG3	1:B:151:LEU:CD1	2.42	0.48
1:B:193:ALA:HA	1:B:197:PHE:HB3	1.94	0.48
1:B:191:LEU:O	1:B:191:LEU:CG	2.50	0.48
1:B:170:LYS:O	1:B:171:LYS:C	2.56	0.48
1:B:49:ARG:HG2	1:B:49:ARG:NH1	2.29	0.48
1:B:57:LYS:NZ	1:B:216:THR:HG23	2.29	0.48
1:A:13:GLU:C	1:A:15:LYS:H	2.22	0.48
1:B:209:TRP:O	1:B:210:LEU:C	2.57	0.47
2:C:15:GLN:HG3	2:C:17:VAL:HG23	1.96	0.47
1:B:29:SER:O	1:B:33:SER:HB2	2.15	0.47
1:A:200:CYS:HB2	1:A:202:HIS:CD2	2.49	0.47
1:A:179:ARG:NH2	1:B:14:GLU:CD	2.73	0.47
1:B:14:GLU:HA	1:B:17:GLN:HG2	1.96	0.47
1:B:64:ASP:HB3	1:B:125:VAL:CG1	2.45	0.47
1:A:9:ASP:O	1:A:9:ASP:CG	2.57	0.46
1:B:176:ASP:HB3	1:B:179:ARG:HB2	1.97	0.46
1:B:213:ARG:HH11	1:B:213:ARG:CG	2.28	0.46
1:A:166:SER:HB2	2:C:28:GLU:CD	2.40	0.46
1:B:213:ARG:HD3	1:B:215:GLU:OE2	2.16	0.46
1:A:204:VAL:HG13	1:A:205:LEU:N	2.30	0.46
1:B:30:GLU:CB	1:B:37:ILE:HD11	2.45	0.46
1:B:63:GLN:OE1	2:D:23:LEU:HB3	2.16	0.46
1:B:153:ASN:HA	1:B:180:HIS:HE1	1.76	0.46
1:B:104:THR:O	1:B:105:ARG:HB2	2.15	0.45
1:A:38:TYR:O	1:A:40:PRO:N	2.49	0.45
1:A:95:GLU:O	1:A:99:THR:HG22	2.16	0.45
1:B:153:ASN:ND2	1:B:176:ASP:H	2.15	0.45
1:B:129:ARG:HB2	1:B:135:SER:OG	2.15	0.45
1:B:24:LEU:HD12	1:B:24:LEU:HA	1.67	0.45
2:D:30:GLU:HA	2:D:35:ASN:O	2.17	0.45
1:A:175:ILE:CG2	1:A:176:ASP:N	2.79	0.45
1:A:106:PRO:HG3	1:A:222:PRO:HG2	1.99	0.45
1:A:162:LEU:HD11	1:A:183:LEU:HD12	1.99	0.45
2:C:36:LYS:HA	2:C:37:PRO:HD3	1.87	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ASP:OD1	1:A:11:LEU:HD13	2.17	0.44
1:A:112:GLU:O	1:A:116:ARG:HG3	2.17	0.44
1:A:204:VAL:CG1	1:A:205:LEU:N	2.80	0.44
1:B:53:LEU:HD12	1:B:53:LEU:O	2.17	0.44
1:A:201:ASN:CB	1:A:204:VAL:CG1	2.86	0.44
2:C:30:GLU:OE1	2:C:36:LYS:HE3	2.17	0.44
1:B:103:PHE:HE2	1:B:220:TRP:O	2.01	0.44
2:C:49:GLU:O	2:C:49:GLU:OE2	2.34	0.44
1:A:40:PRO:O	1:A:41:GLN:C	2.59	0.44
1:B:163:LEU:HD11	1:B:173:ALA:HB1	1.98	0.44
1:A:30:GLU:OE1	1:A:129:ARG:NH1	2.51	0.43
1:A:114:TRP:O	1:A:119:VAL:HG23	2.18	0.43
1:B:169:GLN:O	1:B:171:LYS:N	2.52	0.43
1:A:159:VAL:O	1:A:180:HIS:HB3	2.18	0.43
1:A:181:HIS:HB2	1:A:209:TRP:CZ2	2.53	0.43
1:A:30:GLU:HB2	1:A:37:ILE:HD11	1.99	0.43
2:C:47:TYR:CG	2:C:48:ASP:N	2.87	0.43
2:C:25:LEU:HB3	2:C:26:PRO:HD2	2.00	0.43
1:A:191:LEU:O	1:A:191:LEU:HG	2.18	0.43
1:A:43:ASP:O	1:A:44:VAL:C	2.60	0.43
1:B:51:THR:HG23	1:B:120:LEU:HD12	2.00	0.43
1:B:57:LYS:HZ1	1:B:216:THR:HG23	1.84	0.43
1:B:71:GLN:O	1:B:79:VAL:HA	2.19	0.42
2:C:81:ILE:HG22	2:C:82:LYS:N	2.34	0.42
1:B:9:ASP:C	1:B:11:LEU:N	2.74	0.42
1:A:187:HIS:CE1	1:A:188:PRO:HD2	2.54	0.42
1:B:64:ASP:HB3	1:B:125:VAL:HG13	2.02	0.42
2:D:48:ASP:OD1	2:D:50:SER:OG	2.36	0.42
1:A:41:GLN:HE22	1:A:44:VAL:HG21	1.84	0.42
1:A:6:THR:HG23	1:A:7:TRP:H	1.85	0.42
1:A:57:LYS:HE3	1:A:117:GLN:O	2.19	0.42
1:A:163:LEU:HB3	1:A:169:GLN:CA	2.48	0.42
1:B:7:TRP:C	1:B:9:ASP:H	2.21	0.42
1:A:24:LEU:HD12	1:A:24:LEU:HA	1.83	0.42
1:A:152:ILE:O	1:A:156:ARG:HB2	2.19	0.42
1:B:151:LEU:HD12	1:B:151:LEU:N	2.34	0.42
1:A:189:SER:O	1:A:190:PRO:C	2.61	0.42
1:B:64:ASP:CB	1:B:125:VAL:HG13	2.50	0.42
1:B:139:LEU:HD12	1:B:139:LEU:N	2.35	0.42
1:A:223:VAL:CG1	1:A:224:LEU:N	2.83	0.42
1:B:15:LYS:HG3	1:B:20:PHE:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ILE:HA	1:A:101:PRO:HD2	1.83	0.41
1:B:164:TRP:HA	1:B:185:ALA:O	2.20	0.41
1:A:96:LEU:HD12	1:A:100:ILE:HG22	2.02	0.41
1:B:153:ASN:HD21	1:B:176:ASP:H	1.68	0.41
2:C:38:GLU:O	2:C:38:GLU:HG3	2.20	0.41
1:A:11:LEU:O	1:A:15:LYS:HG3	2.21	0.41
1:A:64:ASP:CB	1:A:125:VAL:HG13	2.44	0.41
1:A:154:GLN:HB3	1:A:155:HIS:ND1	2.35	0.41
1:B:32:GLN:HG2	1:B:33:SER:N	2.35	0.41
1:A:31:ARG:HG2	1:A:37:ILE:HD12	2.02	0.41
1:B:176:ASP:C	1:B:178:GLN:N	2.78	0.41
2:D:44:HIS:HE1	2:D:59:THR:OG1	2.04	0.41
1:A:29:SER:O	1:A:30:GLU:C	2.62	0.41
1:B:96:LEU:O	1:B:100:ILE:HG23	2.21	0.41
2:C:21:SER:HA	2:C:43:VAL:O	2.21	0.41
2:C:27:GLU:H	2:C:27:GLU:HG2	1.54	0.41
2:D:67:PRO:HB3	2:D:84:LEU:HD13	2.03	0.41
1:A:19:TYR:CG	1:A:20:PHE:N	2.89	0.41
1:B:20:PHE:CE2	1:B:24:LEU:HD22	2.56	0.41
1:B:112:GLU:OE2	1:B:116:ARG:NH1	2.35	0.41
2:C:29:VAL:O	2:C:30:GLU:C	2.62	0.41
2:D:41:ILE:HD13	2:D:41:ILE:HA	1.92	0.41
2:D:50:SER:C	2:D:52:ASP:H	2.29	0.41
2:D:58:LEU:HD12	2:D:69:ALA:HB3	2.02	0.41
1:B:14:GLU:C	1:B:16:GLN:H	2.29	0.40
1:A:95:GLU:CD	1:A:200:CYS:O	2.64	0.40
1:A:112:GLU:O	1:A:112:GLU:HG2	2.21	0.40
1:A:15:LYS:HA	1:A:20:PHE:CD1	2.57	0.40
1:A:19:TYR:CD2	1:A:19:TYR:C	2.99	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	
1	A	224/229 (98%)	177 (79%)	36 (16%)	11 (5%)	1	6
1	B	222/229 (97%)	174 (78%)	33 (15%)	15 (7%)	1	2
2	C	71/84 (84%)	62 (87%)	7 (10%)	2 (3%)	4	15
2	D	68/84 (81%)	62 (91%)	5 (7%)	1 (2%)	8	28
All	All	585/626 (94%)	475 (81%)	81 (14%)	29 (5%)	1	6

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	225	PRO
1	A	226	ALA
1	A	227	GLU
1	B	4	GLU
1	B	5	LEU
1	B	6	THR
1	B	171	LYS
2	C	15	GLN
1	A	4	GLU
1	A	103	PHE
1	A	106	PRO
1	B	10	VAL
1	B	13	GLU
1	B	38	TYR
1	B	101	PRO
1	B	177	LYS
2	D	50	SER
1	A	44	VAL
1	A	81	PRO
1	B	77	PHE
1	A	10	VAL
1	B	105	ARG
1	B	170	LYS
1	B	192	SER
1	A	38	TYR
1	A	137	ALA
1	B	81	PRO
1	B	102	GLY
2	C	13	GLY

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	181/194 (93%)	150 (83%)	31 (17%)	2	7
1	B	181/194 (93%)	147 (81%)	34 (19%)	1	5
2	C	64/78 (82%)	52 (81%)	12 (19%)	1	5
2	D	61/78 (78%)	51 (84%)	10 (16%)	2	8
All	All	487/544 (90%)	400 (82%)	87 (18%)	2	6

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLU
1	A	16	GLN
1	A	17	GLN
1	A	21	LEU
1	A	24	LEU
1	A	35	VAL
1	A	36	THR
1	A	41	GLN
1	A	51	THR
1	A	61	LEU
1	A	83	ILE
1	A	92	MET
1	A	96	LEU
1	A	97	GLU
1	A	116	ARG
1	A	125	VAL
1	A	134	HIS
1	A	142	GLU
1	A	143	THR
1	A	145	THR
1	A	157	GLU
1	A	162	LEU
1	A	166	SER
1	A	169	GLN

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Mol	Chain	Res	Type
1	A	178	GLN
1	A	179	ARG
1	A	209	TRP
1	A	212	GLN
1	A	221	MET
1	A	223	VAL
1	A	228	SER
1	B	4	GLU
1	B	5	LEU
1	B	9	ASP
1	B	13	GLU
1	B	15	LYS
1	B	21	LEU
1	B	24	LEU
1	B	26	THR
1	B	31	ARG
1	B	32	GLN
1	B	35	VAL
1	B	36	THR
1	B	51	THR
1	B	61	LEU
1	B	96	LEU
1	B	100	ILE
1	B	116	ARG
1	B	124	THR
1	B	125	VAL
1	B	134	HIS
1	B	142	GLU
1	B	143	THR
1	B	145	THR
1	B	162	LEU
1	B	169	GLN
1	B	177	LYS
1	B	178	GLN
1	B	192	SER
1	B	205	LEU
1	B	209	TRP
1	B	210	LEU
1	B	213	ARG
1	B	223	VAL
1	B	224	LEU
2	C	14	LYS

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Mol	Chain	Res	Type
2	C	18	ILE
2	C	19	GLN
2	C	27	GLU
2	C	29	VAL
2	C	53	GLU
2	C	57	LEU
2	C	66	LYS
2	C	72	ILE
2	C	75	SER
2	C	76	ASN
2	C	80	LYS
2	D	15	GLN
2	D	16	LEU
2	D	36	LYS
2	D	45	THR
2	D	53	GLU
2	D	57	LEU
2	D	72	ILE
2	D	76	ASN
2	D	80	LYS
2	D	82	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	22	ASN
1	A	25	GLN
1	A	46	ASN
1	A	153	ASN
1	A	169	GLN
1	A	178	GLN
1	A	180	HIS
1	A	202	HIS
1	A	212	GLN
1	B	22	ASN
1	B	46	ASN
1	B	91	ASN
1	B	134	HIS
1	B	153	ASN
1	B	169	GLN
1	B	180	HIS

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Mol	Chain	Res	Type
2	C	44	HIS
2	D	15	GLN
2	D	35	ASN
2	D	44	HIS
2	D	73	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 [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	226/229 (98%)	-0.27	2 (0%) 81 75	3, 24, 67, 99	0
1	B	224/229 (97%)	-0.12	6 (2%) 56 47	7, 31, 73, 100	0
2	C	73/84 (86%)	-0.35	1 (1%) 73 65	4, 22, 81, 99	0
2	D	70/84 (83%)	-0.20	0 100 100	11, 33, 72, 88	0
All	All	593/626 (94%)	-0.21	9 (1%) 72 64	3, 28, 72, 100	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	196	GLY	4.4
2	C	12	THR	4.3
1	B	172	GLY	4.2
1	B	10	VAL	2.9
1	A	228	SER	2.8
1	B	170	LYS	2.4
1	B	6	THR	2.2
1	B	179	ARG	2.1
1	A	172	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.