



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 07:31 PM UTC

PDB ID : 4YRV / pdb_00004yrv
Title : Crystal structure of Anabaena transcription factor HetR complexed with 21-bp DNA from hetP promoter
Authors : Hu, H.X.; Jiang, Y.L.; Zhao, M.X.; Zhang, C.C.; Chen, Y.; Zhou, C.Z.
Deposited on : 2015-03-16
Resolution : 2.80 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

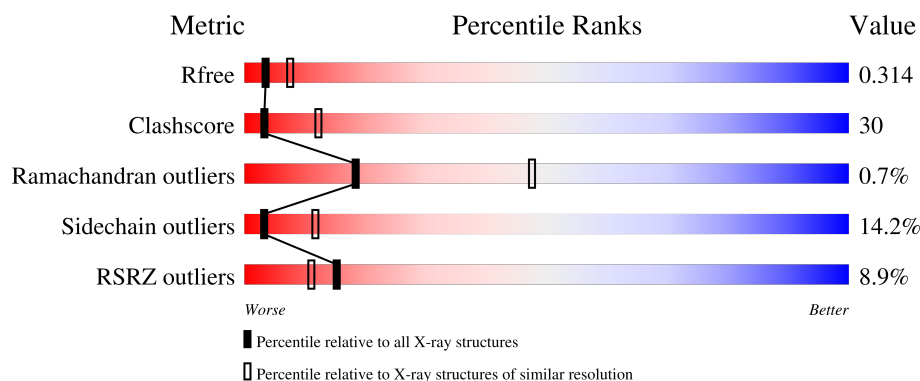
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	307	9% 46% 36% 11% 6%
1	B	307	7% 45% 40% 8% 7%
2	C	21	14% 71% 24% 5%
3	D	21	10% 76% 14% 10%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5648 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 Heterocyst differentiation control protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	286	Total	C	N	O	S	0	0	0
			2368	1514	414	423	17			
1	A	288	Total	C	N	O	S	0	0	0
			2384	1524	418	425	17			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	MET	-	expression tag	UNP P27709
B	-6	GLY	-	expression tag	UNP P27709
B	-5	HIS	-	expression tag	UNP P27709
B	-4	HIS	-	expression tag	UNP P27709
B	-3	HIS	-	expression tag	UNP P27709
B	-2	HIS	-	expression tag	UNP P27709
B	-1	HIS	-	expression tag	UNP P27709
B	0	HIS	-	expression tag	UNP P27709
A	-7	MET	-	expression tag	UNP P27709
A	-6	GLY	-	expression tag	UNP P27709
A	-5	HIS	-	expression tag	UNP P27709
A	-4	HIS	-	expression tag	UNP P27709
A	-3	HIS	-	expression tag	UNP P27709
A	-2	HIS	-	expression tag	UNP P27709
A	-1	HIS	-	expression tag	UNP P27709
A	0	HIS	-	expression tag	UNP P27709

- Molecule 2 is a DNA chain called DNA (5'-D(P*GP*CP*GP*AP*GP*GP*GP*TP*C P*TP*AP*AP*CP*CP*CP*CP*TP*CP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	21	Total	C	N	O	P	0	0	0
			429	203	79	126	21			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	21	Total	C	N	O	P	0	0	0
			432	204	81	126	21			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		
4	C	2	Total	Ca	0	0
			2	2		
4	D	1	Total	Ca	0	0
			1	1		

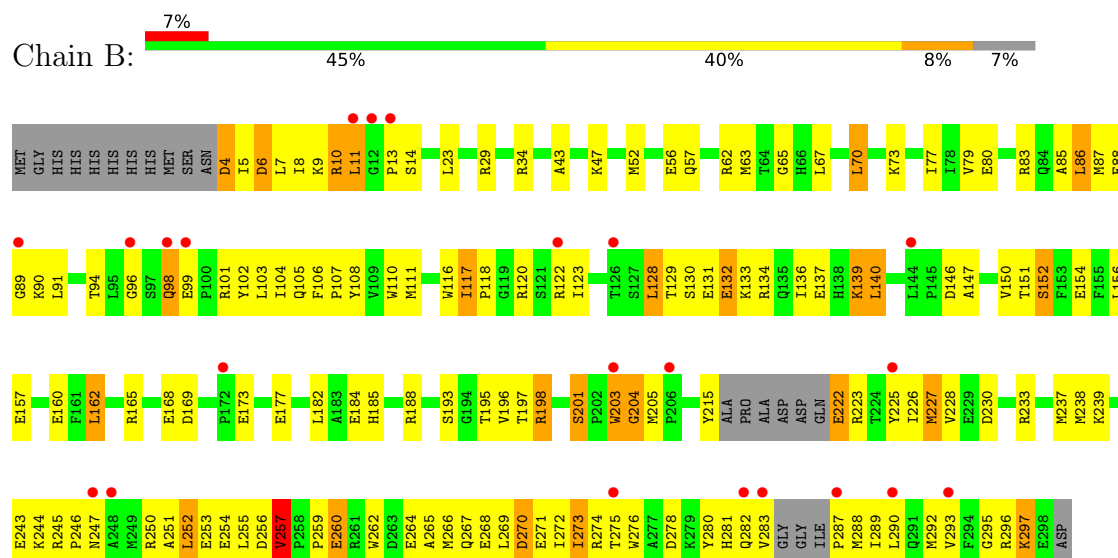
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	13	Total	O	0	0
			13	13		
5	A	13	Total	O	0	0
			13	13		
5	C	3	Total	O	0	0
			3	3		
5	D	2	Total	O	0	0
			2	2		

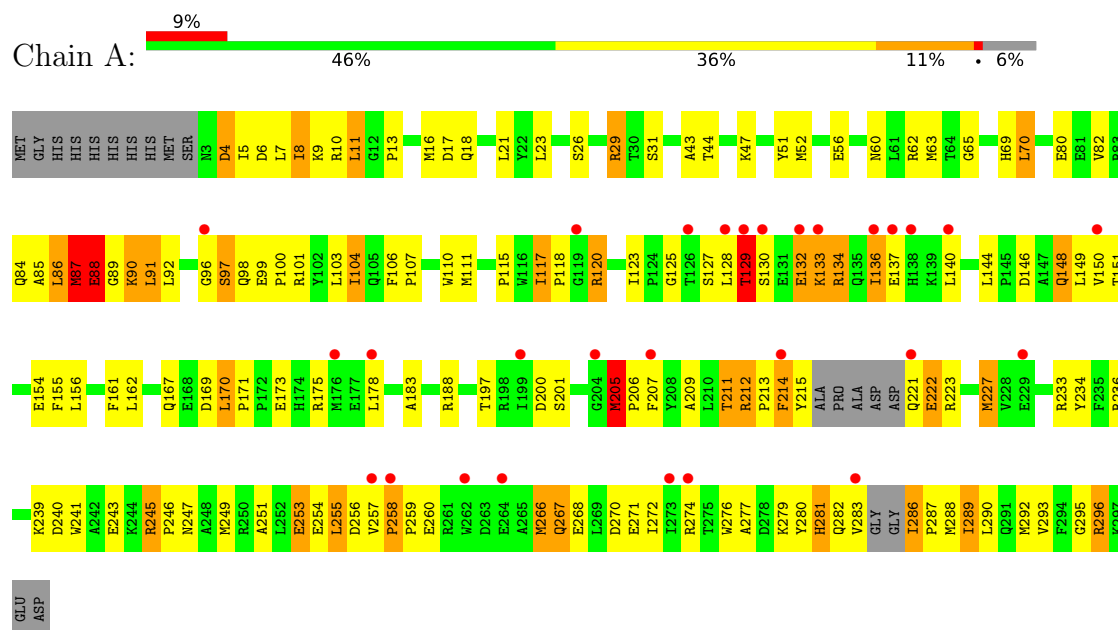
3 Residue-property plots

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: Heterocyst differentiation control protein



• Molecule 1: Heterocyst differentiation control protein



- Molecule 2: DNA (5'-D(P*GP*CP*GP*AP*GP*GP*GP*GP*TP*CP*TP*AP*AP*CP*CP*CP*CP*TP*CP*AP*T)-3')

Chain C: 



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

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	90.56Å 90.56Å 242.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.95 – 2.80 39.95 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.2 (39.95-2.80) 99.1 (39.95-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.58 (at 2.81Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.1_743)	Depositor
R, R_{free}	0.203 , 0.264 0.270 , 0.314	Depositor DCC
R_{free} test set	1298 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	75.4	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 74.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5648	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0.52	0/2443	0.94	11/3298 (0.3%)
1	B	0.48	0/2427	0.90	8/3275 (0.2%)
2	C	0.57	0/480	0.95	1/738 (0.1%)
3	D	0.54	1/484 (0.2%)	1.00	2/745 (0.3%)
All	All	0.51	1/5834 (0.0%)	0.93	22/8056 (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	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	17	DC	O3'-P	-5.37	1.53	1.61

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	MET	CA-C-N	6.76	126.94	120.31
1	A	205	MET	C-N-CA	6.76	126.94	120.31
1	A	281	HIS	N-CA-C	6.31	116.57	108.24
1	A	286	ILE	CA-C-N	6.14	127.52	119.84
1	A	286	ILE	C-N-CA	6.14	127.52	119.84
1	B	123	ILE	N-CA-C	6.06	114.09	107.60
1	A	247	ASN	N-CA-C	6.03	120.33	112.92
2	C	5	DG	C1'-O4'-C4'	-5.61	101.29	109.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	SER	N-CA-C	5.54	119.15	112.38
3	D	18	DT	P-O3'-C3'	-5.54	111.89	120.20
1	B	140	LEU	CA-C-N	5.52	125.28	119.76
1	B	140	LEU	C-N-CA	5.52	125.28	119.76
1	B	204	GLY	N-CA-C	5.48	121.37	114.25
1	A	87	MET	CA-C-N	-5.42	115.14	122.95
1	A	87	MET	C-N-CA	-5.42	115.14	122.95
1	B	205	MET	CA-C-N	5.37	125.29	119.76
1	B	205	MET	C-N-CA	5.37	125.29	119.76
1	B	257	VAL	CA-C-N	5.19	125.73	120.38
1	B	257	VAL	C-N-CA	5.19	125.73	120.38
1	A	125	GLY	N-CA-C	5.17	117.56	110.58
1	A	88	GLU	N-CA-C	5.08	118.31	107.70
3	D	18	DT	C2'-C3'-O3'	-5.07	103.89	111.50

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	129	THR	Peptide
1	A	88	GLU	Peptide
1	A	89	GLY	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2384	0	2377	195	0
1	B	2368	0	2359	182	0
2	C	429	0	236	9	0
3	D	432	0	236	9	0
4	B	1	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
5	A	13	0	0	3	0
5	B	13	0	0	2	0
5	C	3	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	2	0	0	0	0
All	All	5648	0	5208	329	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 30.

All (329) 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:150:VAL:CG1	1:A:154:GLU:HB3	1.59	1.30
1:A:150:VAL:HG13	1:A:154:GLU:CB	1.62	1.28
1:B:150:VAL:CG1	1:B:154:GLU:HB2	1.65	1.26
1:A:213:PRO:O	1:A:214:PHE:HD1	1.16	1.24
1:A:150:VAL:CG1	1:A:154:GLU:CB	2.14	1.19
1:B:150:VAL:CG1	1:B:154:GLU:CB	2.26	1.13
1:A:213:PRO:O	1:A:214:PHE:CD1	2.02	1.13
1:A:128:LEU:HD21	1:A:136:ILE:HG13	1.31	1.07
1:B:150:VAL:HG13	1:B:154:GLU:CB	1.88	1.02
1:B:150:VAL:HG13	1:B:154:GLU:HB2	1.44	0.99
1:A:150:VAL:HG11	1:A:154:GLU:HB3	1.45	0.99
1:A:213:PRO:C	1:A:214:PHE:CD1	2.40	0.98
1:A:150:VAL:CG1	1:A:154:GLU:HB2	1.92	0.97
1:A:150:VAL:HG13	1:A:154:GLU:HB3	1.19	0.97
1:B:150:VAL:HG13	1:B:154:GLU:OE1	1.66	0.96
1:A:117:ILE:O	1:A:120:ARG:NH1	1.98	0.95
1:B:239:LYS:NZ	1:B:243:GLU:OE1	2.01	0.94
1:A:8:ILE:HG13	1:A:21:LEU:HD13	1.50	0.92
1:B:150:VAL:HG12	1:B:154:GLU:HB2	1.51	0.92
1:B:237:MET:HE1	1:B:289:ILE:HB	1.52	0.91
1:A:150:VAL:HG13	1:A:154:GLU:HB2	1.50	0.90
2:C:11:DT:O4	3:D:11:DA:N1	2.06	0.89
1:B:247:ASN:HB2	1:A:296:ARG:HH11	1.38	0.88
1:A:234:TYR:OH	5:A:301:HOH:O	1.90	0.87
1:A:84:GLN:HB3	1:A:90:LYS:HG2	1.55	0.87
1:B:251:ALA:HB3	1:A:292:MET:HE2	1.55	0.86
1:A:137:GLU:HA	1:A:140:LEU:HB2	1.55	0.85
1:B:292:MET:HB3	1:A:292:MET:HG3	1.58	0.85
1:B:227:MET:HE3	1:A:239:LYS:HA	1.58	0.84
1:B:239:LYS:HA	1:A:227:MET:HE3	1.57	0.83
1:B:150:VAL:CG1	1:B:154:GLU:HB3	2.07	0.82
1:B:222:GLU:HB3	1:B:225:TYR:HD1	1.43	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:PRO:C	1:A:214:PHE:HD1	1.85	0.82
1:A:65:GLY:HA2	1:A:70:LEU:HD13	1.60	0.81
1:A:222:GLU:CD	1:A:222:GLU:H	1.87	0.81
1:B:281:HIS:CG	1:B:282:GLN:H	1.99	0.81
1:A:215:TYR:HE1	5:A:302:HOH:O	1.62	0.81
1:B:117:ILE:HB	1:B:120:ARG:HH11	1.46	0.81
1:B:198:ARG:HH21	1:B:198:ARG:HG3	1.46	0.80
1:B:4:ASP:OD2	1:B:4:ASP:N	2.15	0.80
1:A:215:TYR:CE1	5:A:302:HOH:O	2.35	0.77
1:A:128:LEU:O	1:A:133:LYS:N	2.19	0.76
1:B:150:VAL:HG11	1:B:154:GLU:HB3	1.67	0.75
1:B:274:ARG:NH1	1:A:253:GLU:HG2	2.02	0.75
1:A:97:SER:O	1:A:98:GLN:HB3	1.86	0.75
1:B:150:VAL:HG11	1:B:154:GLU:CB	2.17	0.74
1:B:287:PRO:N	1:A:257:VAL:H	1.83	0.74
1:A:281:HIS:CG	1:A:282:GLN:H	2.05	0.74
1:A:85:ALA:HA	1:A:90:LYS:HB3	1.70	0.73
1:B:296:ARG:HH21	1:A:246:PRO:HG2	1.54	0.73
1:A:281:HIS:CD2	1:A:282:GLN:H	2.07	0.73
1:B:129:THR:O	1:B:133:LYS:HG3	1.89	0.72
1:B:11:LEU:HD12	1:B:13:PRO:HB3	1.72	0.71
3:D:17:DC:C2'	3:D:18:DT:H5'	2.21	0.71
2:C:3:DG:H2''	2:C:4:DA:O5'	1.91	0.71
1:B:193:SER:CB	1:B:215:TYR:HE1	2.04	0.71
1:B:132:GLU:HG3	1:B:203:TRP:CE2	2.25	0.70
1:B:116:TRP:CD1	1:B:122:ARG:HG2	2.26	0.70
1:B:150:VAL:HG12	1:B:151:THR:N	2.05	0.70
1:A:170:LEU:O	1:A:175:ARG:NE	2.23	0.70
1:B:238:MET:HE1	1:B:250:ARG:HD2	1.73	0.69
1:B:101:ARG:HH22	1:B:169:ASP:HB2	1.58	0.69
1:B:165:ARG:O	1:B:168:GLU:HB2	1.92	0.69
1:B:297:LYS:HE3	1:A:289:ILE:HD11	1.73	0.69
1:A:150:VAL:HG12	1:A:154:GLU:HB2	1.75	0.68
1:B:256:ASP:O	1:A:286:ILE:N	2.27	0.68
1:B:266:MET:HG2	1:A:249:MET:SD	2.34	0.68
1:A:243:GLU:OE2	1:A:245:ARG:NH1	2.26	0.68
1:B:116:TRP:O	1:B:117:ILE:HD12	1.94	0.68
1:A:128:LEU:HD12	1:A:132:GLU:HB3	1.76	0.68
1:A:87:MET:HE3	1:A:87:MET:HA	1.77	0.67
1:A:62:ARG:HD3	5:C:201:HOH:O	1.95	0.66
1:B:201:SER:O	1:B:204:GLY:N	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:LEU:CD2	1:A:136:ILE:HG13	2.17	0.65
1:B:289:ILE:HD13	1:A:254:GLU:HG2	1.78	0.64
1:B:63:MET:SD	1:A:188:ARG:NH1	2.71	0.64
1:B:281:HIS:CG	1:B:282:GLN:N	2.66	0.64
1:B:5:ILE:HD12	1:B:5:ILE:H	1.64	0.63
1:A:150:VAL:HG12	1:A:151:THR:O	1.98	0.63
1:B:117:ILE:HB	1:B:120:ARG:NH1	2.13	0.63
1:A:99:GLU:HB3	1:A:104:ILE:HD11	1.80	0.63
1:B:150:VAL:HG13	1:B:154:GLU:CD	2.23	0.63
1:A:13:PRO:HG2	1:A:18:GLN:HG2	1.81	0.63
1:B:254:GLU:OE1	1:B:297:LYS:HE2	1.99	0.62
1:B:274:ARG:HH11	1:A:253:GLU:HG2	1.64	0.62
1:A:200:ASP:OD2	1:A:201:SER:N	2.29	0.61
1:B:227:MET:HE1	1:A:239:LYS:HG2	1.82	0.61
1:B:122:ARG:NH2	1:B:147:ALA:O	2.32	0.61
1:A:257:VAL:HG12	1:A:259:PRO:HD3	1.80	0.61
1:A:106:PHE:HB3	1:A:107:PRO:HD3	1.83	0.61
1:A:255:LEU:N	1:A:255:LEU:HD23	2.16	0.61
1:B:5:ILE:HD13	1:A:245:ARG:HH22	1.65	0.61
1:B:150:VAL:CG1	1:B:151:THR:N	2.65	0.60
1:A:6:ASP:HA	1:A:9:LYS:HD3	1.83	0.60
1:B:222:GLU:HB3	1:B:225:TYR:CD1	2.32	0.60
1:A:129:THR:OG1	1:A:130:SER:N	2.35	0.60
1:B:287:PRO:HD3	1:A:257:VAL:HG22	1.83	0.60
1:B:252:LEU:HD12	1:B:252:LEU:C	2.27	0.60
1:A:128:LEU:C	1:A:133:LYS:HD2	2.27	0.60
1:B:260:GLU:OE1	1:B:260:GLU:N	2.32	0.59
1:A:99:GLU:HB3	1:A:104:ILE:CD1	2.32	0.59
1:B:10:ARG:HH22	1:A:86:LEU:C	2.09	0.59
1:A:120:ARG:HD3	1:A:120:ARG:N	2.18	0.59
1:A:170:LEU:HD23	1:A:171:PRO:HD2	1.83	0.59
1:B:96:GLY:HA3	5:B:407:HOH:O	2.03	0.58
1:B:247:ASN:HB2	1:A:296:ARG:NH1	2.13	0.58
1:B:230:ASP:OD1	1:B:233:ARG:NH2	2.37	0.58
1:B:98:GLN:NE2	1:B:225:TYR:HA	2.18	0.58
1:B:227:MET:CE	1:A:239:LYS:HG2	2.34	0.58
1:B:65:GLY:HA2	1:B:70:LEU:HD13	1.87	0.57
1:A:288:MET:C	1:A:289:ILE:HG12	2.29	0.57
1:B:8:ILE:H	1:B:8:ILE:HD12	1.68	0.57
1:B:256:ASP:HB3	1:A:283:VAL:H	1.70	0.57
1:B:278:ASP:OD2	1:B:278:ASP:N	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:ARG:HH21	1:B:198:ARG:CG	2.17	0.56
1:A:222:GLU:CD	1:A:222:GLU:N	2.61	0.56
1:B:83:ARG:HD2	1:B:87:MET:HG3	1.88	0.56
1:B:252:LEU:HD12	1:B:253:GLU:N	2.20	0.56
1:A:31:SER:OG	1:A:98:GLN:O	2.22	0.56
1:B:150:VAL:CG1	1:B:154:GLU:OE1	2.49	0.56
1:A:60:ASN:ND2	1:A:63:MET:HB2	2.20	0.56
1:A:283:VAL:O	1:A:286:ILE:N	2.38	0.56
2:C:5:DG:H2'	2:C:6:DG:H8	1.70	0.56
2:C:11:DT:H3	3:D:11:DA:H2	1.54	0.56
1:B:73:LYS:O	1:B:77:ILE:HD12	2.06	0.56
1:A:129:THR:HA	1:A:133:LYS:HD3	1.88	0.55
1:B:105:GLN:OE1	1:B:105:GLN:HA	2.06	0.55
1:B:86:LEU:HD21	1:A:7:LEU:HD13	1.87	0.55
1:B:102:TYR:HA	1:B:162:LEU:HD23	1.88	0.55
1:B:129:THR:HG22	1:B:131:GLU:H	1.71	0.55
1:A:266:MET:HE2	1:A:270:ASP:OD2	2.06	0.54
1:B:4:ASP:HB2	1:B:5:ILE:HD12	1.88	0.54
1:B:188:ARG:HD2	1:A:63:MET:CE	2.37	0.54
1:B:188:ARG:HD2	1:A:63:MET:HE1	1.90	0.54
1:A:240:ASP:HA	1:A:245:ARG:NH1	2.23	0.54
1:B:14:SER:HB3	1:A:96:GLY:O	2.08	0.54
1:A:8:ILE:HG13	1:A:21:LEU:CD1	2.31	0.54
1:A:274:ARG:HD2	1:A:277:ALA:HB3	1.88	0.54
1:B:130:SER:O	1:B:133:LYS:HB2	2.07	0.54
1:A:268:GLU:O	1:A:272:ILE:HG12	2.08	0.54
1:A:133:LYS:O	1:A:136:ILE:HB	2.07	0.54
1:B:297:LYS:HG3	1:A:289:ILE:HD11	1.89	0.54
3:D:17:DC:H2''	3:D:18:DT:H5'	1.90	0.54
1:A:85:ALA:HA	1:A:90:LYS:O	2.08	0.54
1:B:259:PRO:HG2	1:B:260:GLU:OE1	2.07	0.53
1:B:11:LEU:CD1	1:B:13:PRO:HB3	2.38	0.53
1:B:150:VAL:CG1	1:B:151:THR:H	2.21	0.53
1:B:154:GLU:O	1:B:157:GLU:HB2	2.08	0.53
1:A:26:SER:O	1:A:31:SER:HB3	2.08	0.53
1:A:281:HIS:CG	1:A:282:GLN:N	2.74	0.53
1:B:99:GLU:HB2	1:B:104:ILE:HD11	1.91	0.53
1:A:267:GLN:HG3	1:A:268:GLU:N	2.23	0.53
1:B:11:LEU:HB3	1:B:13:PRO:HD3	1.91	0.53
1:A:110:TRP:CD1	1:A:110:TRP:C	2.87	0.53
3:D:17:DC:H2'	3:D:18:DT:H5'	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:MET:HG3	1:A:251:ALA:HB3	1.91	0.52
1:B:5:ILE:HD12	1:B:5:ILE:N	2.25	0.52
1:A:233:ARG:HG3	1:A:236:ARG:HH12	1.75	0.52
1:B:116:TRP:CH2	1:B:118:PRO:HA	2.45	0.52
1:A:6:ASP:O	1:A:9:LYS:HB2	2.10	0.52
3:D:16:DC:H2''	3:D:17:DC:O5'	2.09	0.52
1:B:62:ARG:NH1	3:D:3:DG:N7	2.57	0.51
1:A:173:GLU:HA	1:A:173:GLU:OE2	2.10	0.51
1:A:136:ILE:O	1:A:140:LEU:HD13	2.10	0.51
1:A:132:GLU:O	1:A:136:ILE:HG12	2.10	0.51
1:B:296:ARG:NH2	1:A:246:PRO:HG2	2.23	0.51
1:B:287:PRO:N	1:A:257:VAL:N	2.57	0.51
1:A:133:LYS:C	1:A:136:ILE:HB	2.35	0.51
1:B:8:ILE:HD12	1:B:8:ILE:N	2.26	0.51
1:B:251:ALA:CB	1:A:292:MET:HE2	2.34	0.51
1:B:260:GLU:H	1:B:260:GLU:CD	2.18	0.51
1:B:270:ASP:O	1:B:274:ARG:HG2	2.11	0.51
1:B:259:PRO:HD3	1:A:286:ILE:HG13	1.92	0.51
1:B:264:GLU:O	1:B:267:GLN:HB3	2.11	0.51
2:C:2:DC:H2'	2:C:3:DG:C8	2.46	0.51
1:B:156:LEU:O	1:B:160:GLU:HG3	2.11	0.50
1:B:222:GLU:OE2	1:B:223:ARG:N	2.44	0.50
1:A:43:ALA:O	1:A:47:LYS:HG3	2.11	0.50
1:A:62:ARG:NH1	2:C:3:DG:N7	2.59	0.50
1:B:280:TYR:O	1:B:281:HIS:HB2	2.10	0.50
1:A:85:ALA:CA	1:A:90:LYS:HB3	2.38	0.50
1:B:106:PHE:HB3	1:B:107:PRO:HD3	1.91	0.50
1:B:106:PHE:HA	1:B:162:LEU:HD21	1.94	0.50
1:B:137:GLU:O	1:B:140:LEU:HB2	2.12	0.50
1:B:223:ARG:NH2	1:A:241:TRP:O	2.43	0.50
1:B:122:ARG:HH22	1:B:147:ALA:H	1.60	0.50
1:B:128:LEU:HD23	1:B:129:THR:N	2.26	0.50
1:B:257:VAL:HG13	1:A:288:MET:HG3	1.93	0.50
1:B:111:MET:SD	1:B:146:ASP:O	2.69	0.50
1:B:182:LEU:O	1:B:185:HIS:HB3	2.12	0.50
1:B:239:LYS:HG3	1:A:227:MET:HE1	1.94	0.50
1:A:13:PRO:HG2	1:A:18:GLN:CG	2.42	0.49
1:B:108:TYR:N	1:B:108:TYR:CD1	2.78	0.49
2:C:11:DT:C4	3:D:11:DA:N1	2.80	0.49
1:A:87:MET:C	1:A:88:GLU:HG2	2.37	0.49
1:A:233:ARG:HG3	1:A:236:ARG:NH1	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:ILE:O	1:A:8:ILE:N	2.46	0.49
1:B:34:ARG:HD2	1:A:69:HIS:ND1	2.28	0.49
1:B:288:MET:HE3	1:A:295:GLY:O	2.12	0.49
1:A:80:GLU:OE1	1:A:80:GLU:HA	2.12	0.49
1:B:254:GLU:HG2	1:A:289:ILE:HD13	1.94	0.49
1:B:293:VAL:O	1:A:290:LEU:HD12	2.13	0.48
1:A:127:SER:C	1:A:129:THR:H	2.20	0.48
1:A:129:THR:HA	1:A:133:LYS:CD	2.44	0.48
1:A:197:THR:OG1	1:A:211:THR:HG21	2.14	0.48
1:B:88:GLU:O	1:B:90:LYS:HG3	2.14	0.48
1:A:84:GLN:C	1:A:90:LYS:HB3	2.39	0.47
1:B:132:GLU:HG3	1:B:203:TRP:CD2	2.50	0.47
1:B:94:THR:C	1:B:96:GLY:N	2.69	0.47
1:B:255:LEU:N	1:B:255:LEU:HD23	2.29	0.47
1:B:269:LEU:O	1:B:272:ILE:HG12	2.14	0.47
1:B:296:ARG:O	1:A:236:ARG:NH2	2.48	0.47
1:B:5:ILE:HG22	1:B:9:LYS:HE3	1.97	0.47
1:B:98:GLN:O	1:B:99:GLU:HG3	2.14	0.47
1:B:116:TRP:CG	1:B:117:ILE:N	2.83	0.47
1:B:182:LEU:HD13	1:A:69:HIS:CE1	2.49	0.47
1:A:128:LEU:HD12	1:A:132:GLU:CB	2.43	0.47
1:B:57:GLN:OE1	1:B:57:GLN:HA	2.14	0.47
1:B:6:ASP:HA	1:B:9:LYS:HB2	1.98	0.46
1:A:123:ILE:HD12	1:A:123:ILE:N	2.31	0.46
1:B:5:ILE:HA	1:B:8:ILE:HD13	1.98	0.46
1:A:8:ILE:CG1	1:A:21:LEU:HD13	2.32	0.46
1:A:133:LYS:HA	1:A:136:ILE:HB	1.97	0.46
1:B:110:TRP:CD1	1:B:110:TRP:C	2.93	0.46
1:B:262:TRP:HZ2	1:A:249:MET:HB2	1.81	0.46
1:A:4:ASP:OD2	1:A:4:ASP:C	2.59	0.46
1:B:273:ILE:O	1:B:276:TRP:HB3	2.16	0.46
1:B:246:PRO:HG2	1:A:296:ARG:HH12	1.80	0.46
1:A:211:THR:O	1:A:212:ARG:HD2	2.16	0.46
1:B:11:LEU:HD12	1:B:13:PRO:CB	2.42	0.45
1:B:268:GLU:O	1:B:269:LEU:C	2.60	0.45
1:A:128:LEU:HD12	1:A:132:GLU:CG	2.46	0.45
1:B:7:LEU:HD13	1:A:52:MET:SD	2.57	0.45
1:B:63:MET:HE3	1:B:67:LEU:HG	1.98	0.45
1:B:150:VAL:HG13	1:B:154:GLU:CG	2.46	0.45
1:B:88:GLU:O	1:B:90:LYS:N	2.50	0.45
1:B:295:GLY:O	1:A:289:ILE:N	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:PRO:O	1:A:148:GLN:NE2	2.44	0.45
2:C:5:DG:H2'	2:C:6:DG:C8	2.49	0.45
1:B:198:ARG:CG	1:B:198:ARG:NH2	2.78	0.45
1:A:286:ILE:HA	1:A:287:PRO:HD3	1.62	0.45
1:B:290:LEU:HD11	1:A:292:MET:SD	2.57	0.45
1:B:136:ILE:O	1:B:139:LYS:HB3	2.16	0.45
1:A:205:MET:HA	1:A:206:PRO:HD3	1.67	0.45
1:A:282:GLN:C	1:A:283:VAL:HG23	2.42	0.45
1:A:256:ASP:C	1:A:258:PRO:HD3	2.41	0.44
1:B:151:THR:O	1:B:152:SER:C	2.60	0.44
1:B:10:ARG:NH2	1:A:86:LEU:C	2.76	0.44
1:B:227:MET:HE3	1:A:239:LYS:CA	2.40	0.44
1:B:227:MET:CE	1:A:239:LYS:HA	2.38	0.44
1:A:123:ILE:HB	1:A:133:LYS:HE2	1.99	0.44
1:B:85:ALA:HA	1:B:90:LYS:O	2.16	0.44
1:A:111:MET:HG2	1:A:146:ASP:O	2.18	0.44
1:A:214:PHE:CD1	1:A:214:PHE:N	2.80	0.44
1:B:5:ILE:O	1:B:9:LYS:HG3	2.18	0.44
1:B:265:ALA:HA	1:A:280:TYR:CZ	2.52	0.44
1:B:268:GLU:O	1:B:271:GLU:HB3	2.17	0.43
1:A:97:SER:O	1:A:98:GLN:CB	2.62	0.43
1:A:178:LEU:HD11	1:A:183:ALA:HB2	1.99	0.43
1:A:292:MET:HE2	1:A:292:MET:HB3	1.63	0.43
1:A:282:GLN:O	1:A:283:VAL:HB	2.18	0.43
1:B:52:MET:HE2	1:A:4:ASP:HB2	1.99	0.43
1:B:228:VAL:HG11	1:A:16:MET:HG3	2.01	0.43
1:A:134:ARG:O	1:A:137:GLU:N	2.52	0.43
1:A:144:LEU:HD23	1:A:144:LEU:HA	1.80	0.43
1:B:196:VAL:HG12	1:B:197:THR:N	2.34	0.43
1:A:65:GLY:HA2	1:A:70:LEU:CD1	2.42	0.43
1:A:239:LYS:NZ	1:A:243:GLU:OE1	2.45	0.43
1:B:128:LEU:HD23	1:B:129:THR:H	1.84	0.43
1:A:97:SER:C	1:A:99:GLU:H	2.26	0.42
1:B:56:GLU:OE2	1:A:29:ARG:NH1	2.41	0.42
1:B:150:VAL:HG12	1:B:151:THR:H	1.75	0.42
1:A:99:GLU:O	1:A:100:PRO:C	2.62	0.42
1:B:283:VAL:O	1:A:257:VAL:HB	2.19	0.42
1:A:169:ASP:CG	1:A:170:LEU:N	2.76	0.42
1:A:255:LEU:N	1:A:255:LEU:CD2	2.83	0.42
1:B:98:GLN:HA	1:A:16:MET:HE2	2.00	0.42
1:A:274:ARG:HD2	1:A:274:ARG:HA	1.76	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:18:DT:H2'	3:D:18:DT:O5'	2.20	0.42
1:B:43:ALA:O	1:B:47:LYS:HG3	2.19	0.42
1:A:7:LEU:O	1:A:11:LEU:HB2	2.18	0.42
1:A:9:LYS:HB3	1:A:9:LYS:HE2	1.84	0.42
1:A:150:VAL:HG13	1:A:154:GLU:OE2	2.19	0.42
1:B:80:GLU:OE1	1:B:80:GLU:HA	2.20	0.42
1:B:103:LEU:HD23	1:B:103:LEU:HA	1.77	0.42
1:A:133:LYS:CA	1:A:136:ILE:HB	2.50	0.42
1:B:117:ILE:HG22	1:B:120:ARG:HD3	2.03	0.41
1:A:91:LEU:HD13	1:A:91:LEU:HA	1.78	0.41
1:A:149:LEU:HD22	1:A:209:ALA:HB2	2.02	0.41
1:A:161:PHE:CD1	1:A:161:PHE:C	2.98	0.41
1:B:5:ILE:HD13	1:A:245:ARG:NH2	2.32	0.41
1:B:70:LEU:HD12	1:B:70:LEU:N	2.34	0.41
1:B:131:GLU:O	1:B:134:ARG:HB3	2.19	0.41
1:B:266:MET:O	1:B:267:GLN:C	2.64	0.41
1:A:167:GLN:HA	1:A:170:LEU:HD12	2.02	0.41
1:B:83:ARG:O	1:B:87:MET:HG2	2.20	0.41
1:B:239:LYS:HA	1:A:227:MET:CE	2.40	0.41
1:A:87:MET:HA	1:A:87:MET:CE	2.41	0.41
1:B:67:LEU:HD23	1:B:67:LEU:HA	1.84	0.41
1:B:269:LEU:HB2	1:A:276:TRP:CZ2	2.56	0.41
1:B:288:MET:C	1:B:289:ILE:HG12	2.45	0.41
1:A:44:THR:HG23	1:A:92:LEU:HD13	2.02	0.41
1:A:100:PRO:HG2	1:A:103:LEU:HD12	2.03	0.41
1:A:150:VAL:HG11	1:A:154:GLU:C	2.45	0.41
1:B:259:PRO:HD3	1:A:286:ILE:CD1	2.50	0.41
5:B:408:HOH:O	1:A:98:GLN:HB2	2.19	0.41
1:A:11:LEU:HB3	1:A:13:PRO:HD3	2.03	0.41
1:A:13:PRO:HB2	1:A:17:ASP:HB2	2.03	0.41
1:A:136:ILE:HD13	1:A:136:ILE:HA	1.87	0.41
1:A:51:TYR:CG	1:A:82:VAL:HG11	2.56	0.41
1:B:107:PRO:HG3	1:B:195:THR:HG21	2.02	0.41
1:B:273:ILE:HD13	1:B:273:ILE:HA	1.79	0.41
1:B:288:MET:HE3	1:B:288:MET:HB3	1.71	0.41
1:A:117:ILE:HG23	1:A:118:PRO:HD2	2.02	0.41
1:A:150:VAL:HG11	1:A:155:PHE:N	2.35	0.41
1:A:276:TRP:O	1:A:279:LYS:HB3	2.21	0.41
1:B:29:ARG:NH2	1:A:56:GLU:OE2	2.50	0.40
1:B:91:LEU:HD23	1:B:91:LEU:HA	1.89	0.40
1:A:107:PRO:HG2	1:A:214:PHE:HA	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:ARG:C	1:A:136:ILE:N	2.78	0.40
1:B:244:LYS:HG2	1:A:223:ARG:NH2	2.36	0.40
1:A:90:LYS:HD3	1:A:90:LYS:HA	1.91	0.40
1:B:237:MET:HB3	1:A:293:VAL:HG21	2.03	0.40
1:A:60:ASN:CG	1:A:63:MET:HB2	2.45	0.40
1:B:11:LEU:HD22	1:B:11:LEU:HA	1.68	0.40
1:B:62:ARG:HE	1:B:62:ARG:HB2	1.71	0.40
1:B:184:GLU:OE2	1:B:184:GLU:HA	2.22	0.40
1:A:103:LEU:HD23	1:A:103:LEU:HA	1.85	0.40
2:C:3:DG:H2'	2:C:4:DA:C8	2.56	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	282/307 (92%)	251 (89%)	28 (10%)	3 (1%)	11	36
1	B	280/307 (91%)	259 (92%)	20 (7%)	1 (0%)	30	60
All	All	562/614 (92%)	510 (91%)	48 (8%)	4 (1%)	18	47

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	THR
1	A	90	LYS
1	A	258	PRO
1	B	89	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	255/269 (95%)	214 (84%)	41 (16%)	2	8
1	B	253/269 (94%)	222 (88%)	31 (12%)	4	16
All	All	508/538 (94%)	436 (86%)	72 (14%)	3	12

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	4	ASP
1	B	6	ASP
1	B	10	ARG
1	B	11	LEU
1	B	23	LEU
1	B	70	LEU
1	B	79	VAL
1	B	86	LEU
1	B	98	GLN
1	B	117	ILE
1	B	128	LEU
1	B	132	GLU
1	B	139	LYS
1	B	152	SER
1	B	162	LEU
1	B	173	GLU
1	B	177	GLU
1	B	198	ARG
1	B	201	SER
1	B	203	TRP
1	B	222	GLU
1	B	226	ILE
1	B	227	MET
1	B	245	ARG
1	B	252	LEU
1	B	257	VAL
1	B	260	GLU

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Mol	Chain	Res	Type
1	B	270	ASP
1	B	273	ILE
1	B	275	THR
1	B	297	LYS
1	A	4	ASP
1	A	8	ILE
1	A	10	ARG
1	A	11	LEU
1	A	23	LEU
1	A	29	ARG
1	A	70	LEU
1	A	86	LEU
1	A	87	MET
1	A	88	GLU
1	A	91	LEU
1	A	101	ARG
1	A	104	ILE
1	A	117	ILE
1	A	120	ARG
1	A	129	THR
1	A	132	GLU
1	A	133	LYS
1	A	134	ARG
1	A	136	ILE
1	A	148	GLN
1	A	156	LEU
1	A	162	LEU
1	A	170	LEU
1	A	205	MET
1	A	207	PHE
1	A	211	THR
1	A	212	ARG
1	A	214	PHE
1	A	221	GLN
1	A	222	GLU
1	A	227	MET
1	A	245	ARG
1	A	253	GLU
1	A	255	LEU
1	A	260	GLU
1	A	266	MET
1	A	267	GLN

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Mol	Chain	Res	Type
1	A	271	GLU
1	A	289	ILE
1	A	296	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	33	HIS
1	B	143	ASN
1	A	98	GLN
1	A	148	GLN
1	A	267	GLN
1	A	282	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	288/307 (93%)	0.81	28 (9%) 13 10	53, 85, 149, 177	0
1	B	286/307 (93%)	0.62	22 (7%) 19 14	52, 84, 132, 149	0
2	C	21/21 (100%)	0.92	3 (14%) 6 5	20, 70, 79, 97	0
3	D	21/21 (100%)	0.52	2 (9%) 14 10	20, 69, 87, 90	0
All	All	616/656 (93%)	0.72	55 (8%) 15 11	20, 83, 137, 177	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	GLU	4.6
1	A	126	THR	4.3
2	C	18	DT	4.3
1	A	257	VAL	4.3
1	B	99	GLU	4.3
1	A	136	ILE	4.2
1	B	287	PRO	4.0
3	D	1	DA	3.9
1	A	119	GLY	3.7
2	C	9	DT	3.7
1	B	225	TYR	3.6
3	D	18	DT	3.5
1	A	207	PHE	3.4
1	A	129	THR	3.2
1	B	144	LEU	3.1
2	C	21	DT	3.0
1	A	262	TRP	3.0
1	A	96	GLY	2.9
1	B	11	LEU	2.9
1	A	138	HIS	2.9
1	A	199	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	12	GLY	2.9
1	A	214	PHE	2.8
1	A	283	VAL	2.7
1	B	203	TRP	2.7
1	A	204	GLY	2.7
1	A	258	PRO	2.6
1	A	133	LYS	2.5
1	A	178	LEU	2.5
1	B	283	VAL	2.5
1	A	273	ILE	2.4
1	A	128	LEU	2.4
1	B	206	PRO	2.4
1	A	274	ARG	2.3
1	B	172	PRO	2.3
1	B	89	GLY	2.2
1	B	98	GLN	2.2
1	B	126	THR	2.2
1	B	290	LEU	2.2
1	A	140	LEU	2.2
1	B	247	ASN	2.2
1	B	96	GLY	2.1
1	A	130	SER	2.1
1	A	221	GLN	2.1
1	B	275	THR	2.1
1	B	13	PRO	2.1
1	B	122	ARG	2.1
1	A	229	GLU	2.1
1	A	150	VAL	2.0
1	B	282	GLN	2.0
1	B	293	VAL	2.0
1	B	248	ALA	2.0
1	A	132	GLU	2.0
1	A	264	GLU	2.0
1	A	176	MET	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CA	C	101	1/1	0.81	0.34	112,112,112,112	0
4	CA	B	301	1/1	0.83	0.12	107,107,107,107	0
4	CA	D	101	1/1	0.91	0.28	96,96,96,96	0
4	CA	C	102	1/1	0.92	0.34	97,97,97,97	0

6.5 Other polymers [i](#)

There are no such residues in this entry.