



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:23 PM UTC

PDB ID : 1RLI / pdb_00001rli
Title : The Structure of Trp Repressor Binding Protein from *Bacillus subtilis*
Authors : Kim, Y.; Quartey, P.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2003-11-25
Resolution : 1.80 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

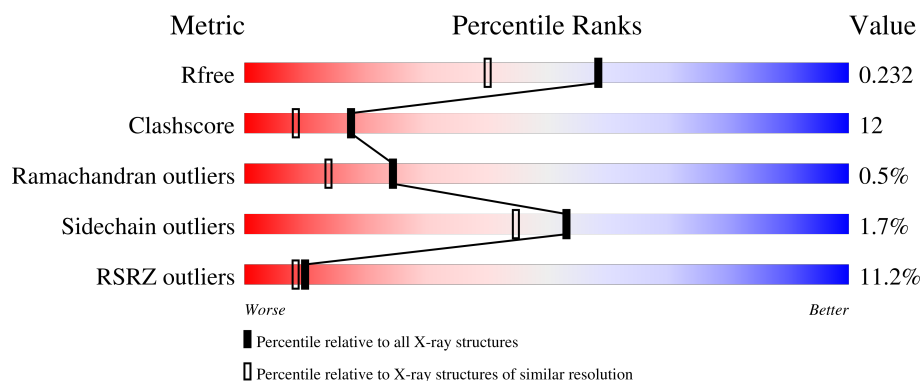
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	184	6% 77% 8% 15%
1	B	184	9% 64% 19% • 15%
1	C	184	16% 73% 15% •• 10%
1	D	184	8% 69% 20% • 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PO4	B	612	-	-	X	-
3	PO4	B	613	-	-	X	-
3	PO4	B	614	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5714 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 Trp Repressor Binding Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1239	797	217	220	5			
1	B	156	Total	C	N	O	S	0	3	0
			1268	813	223	226	6			
1	C	165	Total	C	N	O	S	0	0	0
			1312	846	229	232	5			
1	D	168	Total	C	N	O	S	0	0	0
			1340	864	236	235	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	cloning artifact	UNP P96726
A	-1	ASN	-	cloning artifact	UNP P96726
A	0	ALA	-	cloning artifact	UNP P96726
B	-2	SER	-	cloning artifact	UNP P96726
B	-1	ASN	-	cloning artifact	UNP P96726
B	0	ALA	-	cloning artifact	UNP P96726
C	-2	SER	-	cloning artifact	UNP P96726
C	-1	ASN	-	cloning artifact	UNP P96726
C	0	ALA	-	cloning artifact	UNP P96726
D	-2	SER	-	cloning artifact	UNP P96726
D	-1	ASN	-	cloning artifact	UNP P96726
D	0	ALA	-	cloning artifact	UNP P96726

- Molecule 2 is PLATINUM (II) ION (CCD ID: PT) (formula: Pt).

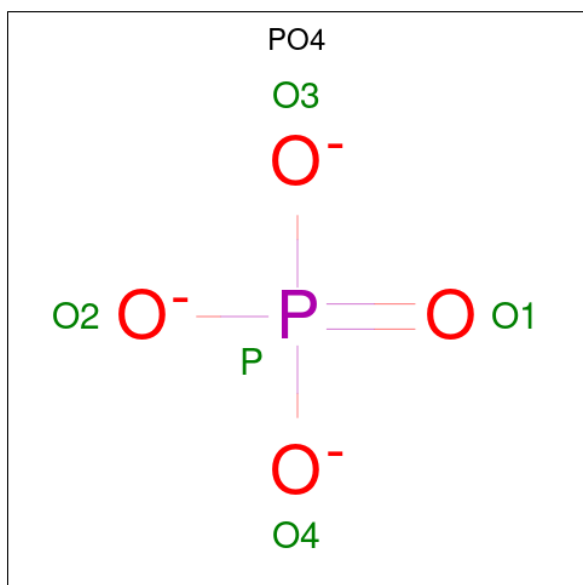
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Pt	0	0
			3	3		
2	B	2	Total	Pt	0	0
			2	2		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	3	Total	Pt	0	0
			3	3		
2	D	3	Total	Pt	0	0
			3	3		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

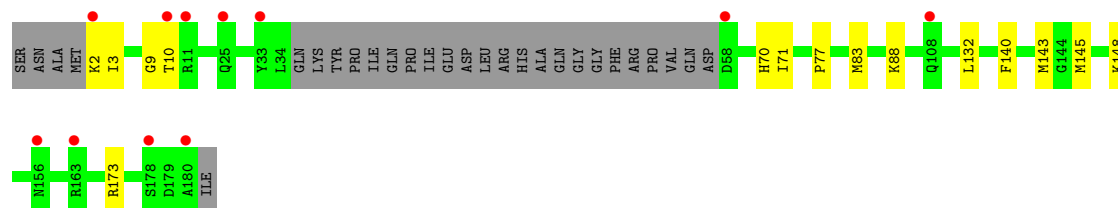
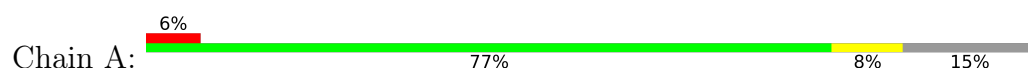
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	129	Total	O	0	0
			129	129		
4	B	137	Total	O	0	0
			137	137		
4	C	109	Total	O	0	0
			109	109		
4	D	154	Total	O	0	0
			154	154		

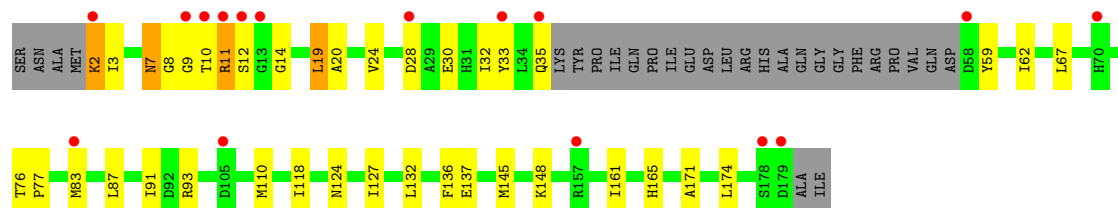
3 Residue-property plots [i](#)

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.

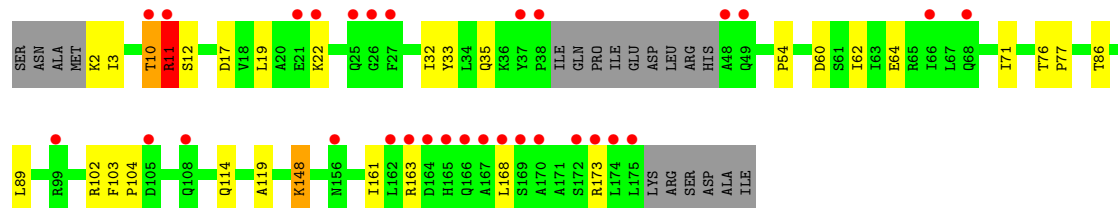
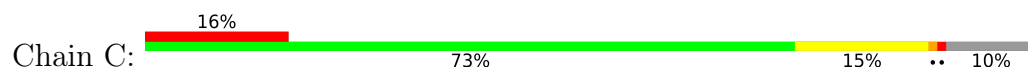
• Molecule 1: Trp Repressor Binding Protein



• Molecule 1: Trp Repressor Binding Protein

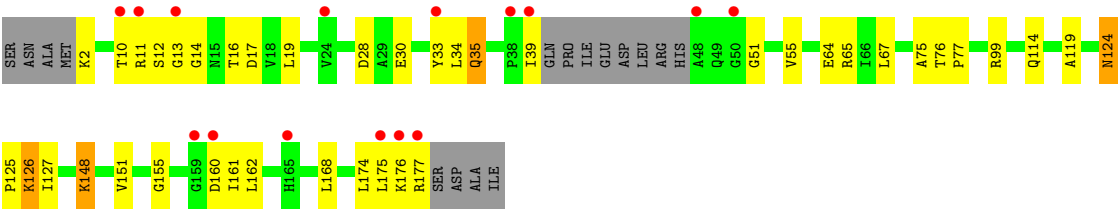


• Molecule 1: Trp Repressor Binding Protein



• Molecule 1: Trp Repressor Binding Protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.99Å 90.67Å 63.45Å 90.00° 97.02° 90.00°	Depositor
Resolution (Å)	36.79 – 1.80 36.79 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.1 (36.79-1.80) 97.2 (36.79-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.230 0.197 , 0.232	Depositor DCC
R_{free} test set	6242 reflections (9.94%)	wwPDB-VP
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.406	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5714	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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: PT, PO4

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	0.48	0/1267	0.87	3/1706 (0.2%)
1	B	0.49	0/1296	0.85	4/1744 (0.2%)
1	C	0.45	0/1344	0.89	5/1812 (0.3%)
1	D	0.52	0/1372	0.92	5/1848 (0.3%)
All	All	0.49	0/5279	0.88	17/7110 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	77	PRO	N-CA-C	-7.45	100.25	111.57
1	C	11	ARG	N-CA-C	7.04	120.07	110.35
1	C	148	LYS	N-CA-C	6.21	120.05	112.23
1	C	103	PHE	CA-C-N	6.16	126.62	119.47
1	C	103	PHE	C-N-CA	6.16	126.62	119.47
1	D	77	PRO	N-CA-C	-5.96	102.52	111.57
1	A	148	LYS	N-CA-C	5.72	119.98	112.89
1	C	77	PRO	N-CA-C	-5.70	102.90	111.57
1	D	28	ASP	N-CA-C	-5.66	101.13	109.62
1	B	77	PRO	N-CA-C	-5.60	103.01	111.41
1	A	9	GLY	N-CA-C	-5.49	104.78	112.13
1	D	124	ASN	CA-C-N	5.37	125.44	119.32
1	D	124	ASN	C-N-CA	5.37	125.44	119.32
1	B	174	LEU	N-CA-C	5.34	117.52	111.11
1	B	24	VAL	N-CA-C	5.33	117.30	112.29
1	D	148	LYS	N-CA-C	5.09	118.64	112.23
1	B	148	LYS	N-CA-C	5.01	119.11	112.89

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1239	0	1237	24	0
1	B	1268	0	1260	43	0
1	C	1312	0	1303	26	0
1	D	1340	0	1340	34	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
3	B	15	0	0	11	0
4	A	129	0	0	14	0
4	B	137	0	0	14	0
4	C	109	0	0	2	0
4	D	154	0	0	7	0
All	All	5714	0	5140	128	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 12.

All (128) 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:C:114:GLN:HE21	1:C:148:LYS:HE3	1.17	1.03
1:A:143:MET:HB2	1:A:145:MET:HE2	1.46	0.94
1:D:12:SER:HB3	4:D:645:HOH:O	1.68	0.94
1:B:83[A]:MET:HE3	1:B:87:LEU:HG	1.52	0.89
1:A:143:MET:CB	1:A:145:MET:HE2	2.05	0.85
1:D:11:ARG:HD3	1:D:33:TYR:CD2	2.12	0.85
1:A:140:PHE:HA	1:A:145:MET:HE3	1.61	0.81
1:D:14:GLY:HA3	4:D:756:HOH:O	1.83	0.79
1:B:9:GLY:HA3	3:B:613:PO4:O4	1.83	0.78
1:B:83[A]:MET:HG3	1:B:132:LEU:HD11	1.68	0.75
3:B:613:PO4:O2	4:B:728:HOH:O	2.05	0.74
1:C:54:PRO:HB3	1:C:86:THR:HG22	1.68	0.74
1:B:83[B]:MET:SD	4:B:702:HOH:O	2.46	0.73
1:B:7[A]:ASN:ND2	4:B:728:HOH:O	2.20	0.73

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83[A]:MET:HE3	1:B:87:LEU:CG	2.18	0.73
1:A:3:ILE:HG22	1:A:71:ILE:HB	1.71	0.73
1:D:11:ARG:HA	1:D:17:ASP:OD1	1.89	0.72
1:B:7[A]:ASN:HD21	3:B:613:PO4:P	2.12	0.72
1:D:99:ARG:HD3	4:D:762:HOH:O	1.90	0.71
1:C:22:LYS:NZ	1:C:168:LEU:HD22	2.06	0.71
1:B:7[A]:ASN:C	1:B:7[A]:ASN:HD22	1.98	0.70
1:C:54:PRO:HG2	1:C:89:LEU:HD22	1.74	0.69
1:A:143:MET:CG	1:A:145:MET:HE2	2.24	0.68
1:B:7[A]:ASN:ND2	1:B:9:GLY:H	1.91	0.68
1:B:83[B]:MET:HG2	4:B:701:HOH:O	1.94	0.67
1:B:7[A]:ASN:ND2	3:B:613:PO4:O3	2.29	0.66
1:A:83:MET:CE	4:A:690:HOH:O	2.44	0.66
1:D:114:GLN:HE21	1:D:148:LYS:HE3	1.61	0.65
1:C:11:ARG:HD3	1:C:33:TYR:CD2	2.32	0.65
1:C:60:ASP:O	1:C:64:GLU:HG3	1.98	0.64
1:C:102:ARG:HG2	1:C:102:ARG:HH11	1.63	0.64
1:C:54:PRO:CG	1:C:89:LEU:HD22	2.29	0.63
1:D:114:GLN:NE2	1:D:148:LYS:HE3	2.13	0.63
1:A:143:MET:HG3	1:A:145:MET:HE2	1.81	0.62
1:B:83[A]:MET:HE1	1:B:91:ILE:HD12	1.82	0.61
1:D:155:GLY:HA2	1:D:160:ASP:OD2	2.00	0.61
1:A:132:LEU:CD1	4:A:690:HOH:O	2.47	0.61
1:D:11:ARG:NH2	1:D:34:LEU:O	2.34	0.60
1:B:124:ASN:HD21	1:B:127:ILE:HD12	1.64	0.60
1:A:140:PHE:HD1	1:A:145:MET:HE3	1.66	0.60
1:D:11:ARG:CZ	1:D:33:TYR:HB3	2.32	0.60
1:A:88:LYS:HD3	4:A:689:HOH:O	2.01	0.59
1:D:2:LYS:HG2	4:D:764:HOH:O	2.02	0.59
1:B:59:TYR:CE2	1:B:93:ARG:HD2	2.38	0.58
1:A:83:MET:HG2	4:A:690:HOH:O	2.03	0.58
1:B:132:LEU:CD1	4:B:701:HOH:O	2.52	0.58
1:D:162:LEU:HA	1:D:168:LEU:HD11	1.86	0.57
1:B:93:ARG:HD3	4:B:736:HOH:O	2.03	0.57
1:D:151:VAL:HG23	1:D:174:LEU:HD13	1.88	0.56
1:B:7[A]:ASN:HD22	1:B:9:GLY:H	1.54	0.55
1:C:11:ARG:HA	1:C:17:ASP:CG	2.31	0.55
1:D:175:LEU:HD23	4:D:648:HOH:O	2.07	0.55
1:A:143:MET:HB2	1:A:145:MET:CE	2.30	0.55
1:B:30:GLU:HG3	4:B:750:HOH:O	2.05	0.55
1:A:173:ARG:HD3	4:A:696:HOH:O	2.06	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83[B]:MET:HE3	4:B:701:HOH:O	2.06	0.55
1:A:83:MET:HE3	4:A:689:HOH:O	2.07	0.54
4:B:664:HOH:O	1:D:176:LYS:HE2	2.07	0.54
1:C:22:LYS:HZ2	1:C:168:LEU:HB3	1.72	0.54
1:B:19:LEU:HG	1:B:161:ILE:HG12	1.91	0.53
1:C:2:LYS:NZ	4:C:696:HOH:O	2.40	0.53
1:D:124:ASN:N	1:D:125:PRO:HD3	2.23	0.53
1:C:102:ARG:C	1:C:104:PRO:HD3	2.34	0.53
1:A:83:MET:CE	4:A:689:HOH:O	2.57	0.52
1:C:11:ARG:HG2	1:C:33:TYR:CE2	2.44	0.52
1:D:16:THR:HG23	1:D:75:ALA:HB1	1.92	0.51
1:A:70:HIS:CE1	4:A:706:HOH:O	2.63	0.51
1:A:83:MET:SD	4:A:689:HOH:O	2.60	0.50
1:B:59:TYR:CZ	1:B:93:ARG:HD2	2.46	0.50
3:B:613:PO4:P	4:B:624:HOH:O	2.69	0.49
1:D:11:ARG:HG3	4:D:661:HOH:O	2.11	0.49
1:A:83:MET:HE2	4:A:690:HOH:O	2.08	0.49
1:D:175:LEU:HD22	1:D:175:LEU:N	2.27	0.49
1:B:83[A]:MET:CE	1:B:87:LEU:HG	2.35	0.48
1:C:114:GLN:NE2	1:C:148:LYS:HE3	2.03	0.48
1:B:7[A]:ASN:ND2	1:B:7[A]:ASN:C	2.70	0.48
1:D:10:THR:O	1:D:17:ASP:OD2	2.31	0.47
1:A:140:PHE:HD1	1:A:145:MET:CE	2.26	0.47
1:A:3:ILE:CG2	1:A:71:ILE:HB	2.39	0.47
1:C:10:THR:HA	4:C:682:HOH:O	2.13	0.47
1:B:2:LYS:HB2	3:B:614:PO4:O1	2.15	0.47
1:C:22:LYS:CE	1:C:168:LEU:HD22	2.45	0.47
1:D:39:ILE:HG12	1:D:51:GLY:O	2.15	0.47
1:B:83[A]:MET:SD	1:B:136:PHE:CZ	3.08	0.46
1:C:32:ILE:HD13	1:C:62:ILE:HG23	1.97	0.46
1:B:19:LEU:HD22	1:B:171:ALA:HB2	1.98	0.46
1:A:83:MET:HE3	4:A:690:HOH:O	2.12	0.46
1:C:19:LEU:HG	1:C:161:ILE:HG12	1.97	0.45
1:D:64:GLU:HG3	4:D:752:HOH:O	2.16	0.45
1:B:14:GLY:HA3	3:B:612:PO4:O4	2.15	0.45
1:B:20:ALA:HA	1:B:118:ILE:HD13	1.98	0.45
1:B:2:LYS:CB	3:B:614:PO4:O1	2.65	0.45
1:B:7[A]:ASN:HD22	1:B:8:GLY:N	2.14	0.45
1:C:76:THR:O	1:C:119:ALA:HA	2.16	0.45
1:A:2:LYS:NZ	4:A:707:HOH:O	2.50	0.45
1:B:32:ILE:HG21	1:B:62:ILE:HG23	1.98	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:76:THR:O	1:D:119:ALA:HA	2.16	0.44
1:C:173:ARG:HH11	1:C:173:ARG:HG2	1.82	0.44
1:B:10:THR:N	3:B:613:PO4:O4	2.50	0.44
1:B:76:THR:HG21	1:B:83[A]:MET:HG2	1.99	0.44
1:C:11:ARG:HA	1:C:17:ASP:OD2	2.18	0.44
1:D:13:GLY:O	1:D:14:GLY:C	2.60	0.43
1:B:137:GLU:HG3	1:D:127:ILE:HG21	2.01	0.43
1:B:165:HIS:CD2	1:D:177:ARG:HD3	2.53	0.43
1:D:11:ARG:HA	1:D:17:ASP:CG	2.43	0.43
1:D:11:ARG:HH11	1:D:11:ARG:HG2	1.84	0.43
1:C:163:ARG:HG2	1:C:163:ARG:O	2.18	0.43
1:B:11:ARG:HB2	1:B:33:TYR:CD2	2.54	0.43
1:B:2:LYS:HD3	1:B:3:ILE:N	2.34	0.43
1:C:3:ILE:HG12	1:C:71:ILE:HB	2.01	0.43
1:D:30:GLU:OE2	1:D:65:ARG:NH2	2.49	0.43
1:B:83[A]:MET:CG	4:B:701:HOH:O	2.66	0.43
1:D:126:LYS:HG3	1:D:127:ILE:HG12	2.01	0.43
1:B:10:THR:HG22	4:B:731:HOH:O	2.19	0.42
3:B:612:PO4:P	4:B:624:HOH:O	2.76	0.42
1:D:67:LEU:HD23	1:D:67:LEU:HA	1.90	0.42
1:B:67:LEU:HD23	1:B:67:LEU:HA	1.89	0.42
1:A:10:THR:HG22	4:A:712:HOH:O	2.20	0.42
1:C:11:ARG:HA	1:C:17:ASP:OD1	2.19	0.41
1:C:102:ARG:HG2	1:C:102:ARG:NH1	2.32	0.41
1:B:110:MET:O	1:B:145:MET:HG2	2.19	0.41
1:B:35:GLN:HG3	4:B:665:HOH:O	2.21	0.41
1:B:7[A]:ASN:CG	3:B:613:PO4:O3	2.64	0.41
1:D:19:LEU:HG	1:D:161:ILE:HG12	2.02	0.41
1:D:35:GLN:HB3	1:D:55:VAL:HB	2.03	0.41
1:A:88:LYS:HE2	4:A:646:HOH:O	2.20	0.40
1:C:54:PRO:HG3	1:C:89:LEU:HD22	2.03	0.40
1:D:11:ARG:O	1:D:12:SER:HB2	2.21	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	152/184 (83%)	148 (97%)	4 (3%)	0	100	100
1	B	155/184 (84%)	151 (97%)	2 (1%)	2 (1%)	9	3
1	C	161/184 (88%)	156 (97%)	4 (2%)	1 (1%)	21	11
1	D	164/184 (89%)	160 (98%)	4 (2%)	0	100	100
All	All	632/736 (86%)	615 (97%)	14 (2%)	3 (0%)	24	14

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	12	SER
1	B	12	SER
1	B	11	ARG

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	130/154 (84%)	130 (100%)	0	100	100
1	B	134/154 (87%)	129 (96%)	5 (4%)	30	18
1	C	137/154 (89%)	134 (98%)	3 (2%)	45	34
1	D	140/154 (91%)	138 (99%)	2 (1%)	59	52
All	All	541/616 (88%)	531 (98%)	10 (2%)	53	43

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2	LYS
1	B	7[A]	ASN
1	B	7[B]	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	19	LEU
1	B	28	ASP
1	C	10	THR
1	C	11	ARG
1	C	35	GLN
1	D	35	GLN
1	D	126	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	31	HIS
1	A	68	GLN
1	A	109	GLN
1	A	134	GLN
1	A	141	HIS
1	B	108	GLN
1	B	109	GLN
1	B	114	GLN
1	B	124	ASN
1	B	134	GLN
1	B	141	HIS
1	B	165	HIS
1	C	25	GLN
1	C	96	GLN
1	C	114	GLN
1	C	141	HIS
1	C	165	HIS
1	D	35	GLN
1	D	49	GLN
1	D	109	GLN
1	D	114	GLN
1	D	134	GLN
1	D	141	HIS

5.3.3 RNA ⓘ

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 14 ligands modelled in this entry, 11 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PO4	B	613	-	4,4,4	1.88	2 (50%)	6,6,6	0.52	0
3	PO4	B	614	2	4,4,4	1.77	1 (25%)	6,6,6	0.46	0
3	PO4	B	612	-	4,4,4	1.81	2 (50%)	6,6,6	0.49	0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	612	PO4	P-O4	-2.60	1.47	1.54
3	B	613	PO4	P-O2	-2.46	1.47	1.54
3	B	614	PO4	P-O4	-2.30	1.47	1.54
3	B	613	PO4	P-O4	-2.09	1.48	1.54
3	B	612	PO4	P-O2	-2.07	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	613	PO4	7	0
3	B	614	PO4	2	0
3	B	612	PO4	2	0

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	156/184 (84%)	0.37	11 (7%) 22 20	14, 27, 46, 69	0
1	B	156/184 (84%)	0.33	16 (10%) 12 10	9, 24, 48, 67	3 (1%)
1	C	165/184 (89%)	0.92	30 (18%) 3 3	14, 33, 56, 66	0
1	D	168/184 (91%)	0.45	15 (8%) 15 13	13, 24, 47, 70	0
All	All	645/736 (87%)	0.52	72 (11%) 10 8	9, 27, 53, 70	3 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	179	ASP	5.6
1	C	10	THR	5.3
1	A	180	ALA	4.7
1	B	10	THR	4.6
1	B	12	SER	3.8
1	D	33	TYR	3.7
1	C	11	ARG	3.5
1	D	160	ASP	3.5
1	C	27	PHE	3.4
1	C	22	LYS	3.4
1	C	169	SER	3.3
1	C	166	GLN	3.3
1	D	13	GLY	3.3
1	D	10	THR	3.3
1	C	99	ARG	3.2
1	C	108	GLN	3.2
1	C	168	LEU	3.1
1	C	175	LEU	3.1
1	D	175	LEU	3.1
1	B	33	TYR	3.0
1	C	162	LEU	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	170	ALA	3.0
1	C	37	TYR	2.9
1	B	2	LYS	2.9
1	A	33	TYR	2.9
1	D	11	ARG	2.8
1	C	49	GLN	2.8
1	B	83[A]	MET	2.8
1	C	21	GLU	2.7
1	C	167	ALA	2.7
1	C	105	ASP	2.7
1	C	172	SER	2.7
1	D	159	GLY	2.6
1	D	177	ARG	2.6
1	C	48	ALA	2.6
1	D	48	ALA	2.6
1	A	2	LYS	2.5
1	C	165	HIS	2.5
1	A	58	ASP	2.5
1	C	25	GLN	2.5
1	C	38	PRO	2.5
1	C	174	LEU	2.4
1	B	9	GLY	2.4
1	D	50	GLY	2.4
1	B	70	HIS	2.4
1	D	165	HIS	2.4
1	D	176	LYS	2.4
1	C	68	GLN	2.4
1	A	178	SER	2.3
1	B	28	ASP	2.3
1	C	26	GLY	2.2
1	D	39	ILE	2.2
1	C	156	ASN	2.2
1	A	10	THR	2.2
1	A	25	GLN	2.2
1	B	178	SER	2.2
1	C	164	ASP	2.2
1	A	156	ASN	2.2
1	A	163	ARG	2.2
1	D	24	VAL	2.2
1	B	157	ARG	2.2
1	B	13	GLY	2.1
1	A	11	ARG	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	163	ARG	2.1
1	B	105	ASP	2.1
1	D	38	PRO	2.1
1	A	108	GLN	2.1
1	B	58	ASP	2.1
1	C	66	ILE	2.0
1	B	35	GLN	2.0
1	B	11	ARG	2.0
1	C	173	ARG	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](#)

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(Å ²)	Q<0.9
3	PO4	B	612	5/5	0.83	0.18	36,40,41,43	1
2	PT	A	606	1/1	0.89	0.23	118,118,118,118	1
3	PO4	B	613	5/5	0.90	0.20	27,39,40,40	3
3	PO4	B	614	5/5	0.90	0.12	35,35,41,42	1
2	PT	B	608	1/1	0.92	0.17	113,113,113,113	1
2	PT	D	607	1/1	0.93	0.14	74,74,74,74	1
2	PT	C	609	1/1	0.94	0.16	72,72,72,72	1
2	PT	C	610	1/1	0.96	0.13	83,83,83,83	1
2	PT	A	605	1/1	0.96	0.11	57,57,57,57	1
2	PT	D	611	1/1	0.96	0.12	60,60,60,60	1
2	PT	C	604	1/1	0.97	0.10	47,47,47,47	1
2	PT	A	601	1/1	0.98	0.08	53,53,53,53	1
2	PT	D	602	1/1	0.99	0.08	52,52,52,52	1
2	PT	B	603	1/1	1.00	0.07	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.