



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:05 AM UTC

PDB ID : 1TGH / pdb_00001tgh
Title : TATA BINDING PROTEIN (TBP)/DNA COMPLEX
Authors : Juo, Z.S.; Dickerson, R.E.
Deposited on : 1996-02-13
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

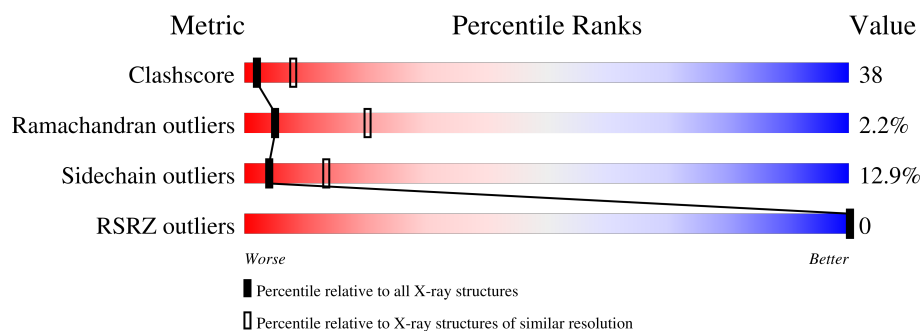
1 Overall quality at a glance

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))
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	B	12	8% 83% 8%
1	C	12	17% 75% 8%
2	A	185	36% 49% 12% ••

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2355 atoms, of which 422 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 (5'-D(*CP*GP*TP*AP*TP*AP*TP*AP*TP*AP*CP*G)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	12	Total	C	H	N	O	P	0	0	0
			267	118	24	44	70	11			
1	C	12	Total	C	H	N	O	P	0	0	0
			267	118	24	44	70	11			

- Molecule 2 is a protein called PROTEIN (TATA BINDING PROTEIN (TBP)).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	A	180	Total	C	H	N	O	S	0	0	0
			1767	927	338	252	243	7			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	151	GLY	ALA	conflict	UNP P20226
A	153	ARG	GLU	conflict	UNP P20226
A	154	GLY	SER	conflict	UNP P20226

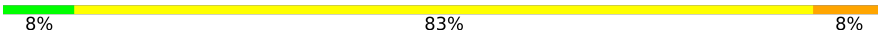
- Molecule 3 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	5	Total	H	O	0	0
			15	10	5		
3	C	1	Total	H	O	0	0
			3	2	1		
3	A	12	Total	H	O	0	0
			36	24	12		

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.

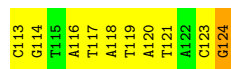
- Molecule 1: DNA (5'-D(*CP*GP*TP*AP*TP*AP*TP*AP*TP*AP*CP*G)-3')

Chain B: 

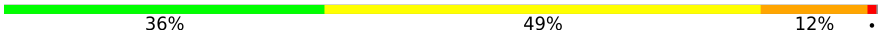


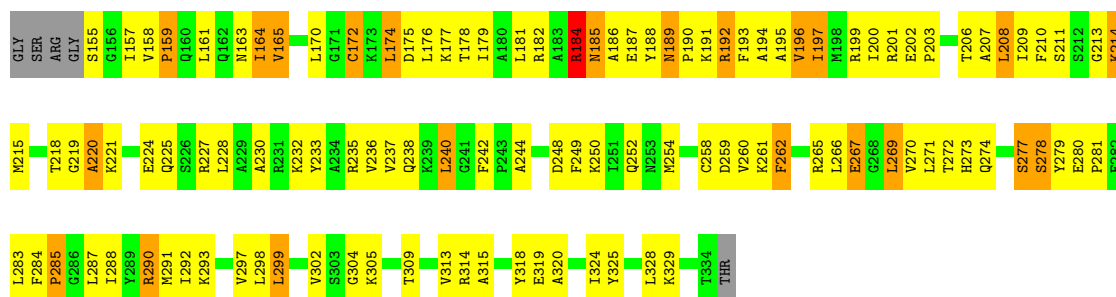
- Molecule 1: DNA (5'-D(*CP*GP*TP*AP*TP*AP*TP*AP*TP*AP*CP*G)-3')

Chain C: 



- Molecule 2: PROTEIN (TATA BINDING PROTEIN (TBP))

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.97Å 67.40Å 86.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.90 8.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.90) 93.6 (8.00-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.15 (at 2.90Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.214 , 0.294 0.213 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	61.7	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 70.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2355	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.73% 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	B	0.58	0/272	1.11	2/418 (0.5%)
1	C	0.53	0/272	1.06	1/418 (0.2%)
2	A	0.85	0/1455	1.29	19/1958 (1.0%)
All	All	0.78	0/1999	1.23	22/2794 (0.8%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	242	PHE	CA-C-N	8.13	130.00	119.84
2	A	242	PHE	C-N-CA	8.13	130.00	119.84
2	A	280	GLU	CA-C-N	8.09	129.95	119.84
2	A	280	GLU	C-N-CA	8.09	129.95	119.84
2	A	165	VAL	N-CA-C	-6.52	98.98	108.11
2	A	182	ARG	N-CA-C	6.35	121.16	113.41
2	A	164	ILE	N-CA-C	-6.19	99.79	108.27
2	A	248	ASP	N-CA-C	6.07	119.76	111.39
2	A	277	SER	N-CA-C	5.78	117.98	110.53
2	A	285	PRO	CA-C-N	-5.76	118.12	122.33
2	A	285	PRO	C-N-CA	-5.76	118.12	122.33
2	A	194	ALA	N-CA-C	5.69	117.82	110.65
1	C	124	DG	C2'-C3'-O3'	-5.57	103.15	111.50
2	A	285	PRO	N-CA-C	5.53	120.43	114.68
1	B	111	DC	C5'-C4'-C3'	-5.52	106.62	114.90
2	A	197	ILE	N-CA-C	-5.38	101.89	109.21
2	A	271	LEU	N-CA-C	5.25	116.86	111.03
2	A	208	LEU	N-CA-C	-5.23	100.20	108.73
2	A	305	LYS	N-CA-C	-5.17	102.63	110.28
2	A	213	GLY	N-CA-C	-5.14	107.56	114.25
2	A	278	SER	N-CA-C	-5.08	100.32	108.55
1	B	108	DA	N9-C1'-C2'	-5.02	105.97	113.50

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	B	243	24	138	32	0
1	C	243	24	138	20	0
2	A	1429	338	1521	105	0
3	A	12	24	0	0	0
3	B	5	10	0	0	0
3	C	1	2	0	0	0
All	All	1933	422	1797	137	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 38.

All (137) 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:B:108:DA:H2''	1:B:109:DT:C5'	1.70	1.20
1:C:123:DC:H2''	1:C:124:DG:C8	1.87	1.10
1:B:108:DA:H2''	1:B:109:DT:H5''	1.20	1.10
1:B:102:DG:H2''	1:B:103:DT:H5'	1.38	1.05
1:B:109:DT:H2''	1:B:110:DA:H5''	1.37	1.01
1:B:102:DG:H2''	1:B:103:DT:C5'	1.91	0.98
1:B:105:DT:H5''	2:A:288:ILE:HD13	1.54	0.90
1:B:109:DT:C6	1:B:109:DT:H5'	2.08	0.89
1:C:117:DT:H5''	2:A:197:ILE:HD13	1.56	0.88
1:B:101:DC:H2''	1:B:102:DG:C8	2.08	0.87
1:B:108:DA:H2''	1:B:109:DT:H5'	1.55	0.87
2:A:277:SER:HB3	2:A:287:LEU:HD11	1.57	0.86
1:B:109:DT:C2'	1:B:110:DA:H5''	2.07	0.85
2:A:176:LEU:HD22	2:A:196:VAL:HG12	1.60	0.84
1:B:108:DA:C2'	1:B:109:DT:H5''	2.07	0.81
1:B:109:DT:H1'	2:A:210:PHE:CE2	2.15	0.81
2:A:192:ARG:HH22	2:A:197:ILE:HD12	1.47	0.78
2:A:196:VAL:HG13	2:A:209:ILE:HB	1.69	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:DT:H5'	1:B:109:DT:H6	1.47	0.74
1:C:123:DC:C2'	1:C:124:DG:C8	2.69	0.73
2:A:313:VAL:HG12	2:A:315:ALA:H	1.56	0.71
1:B:102:DG:H2''	1:B:103:DT:H5''	1.72	0.71
2:A:270:VAL:HG12	2:A:274:GLN:NE2	2.04	0.70
1:B:108:DA:C2'	1:B:109:DT:C5'	2.61	0.70
2:A:266:LEU:O	2:A:270:VAL:HG23	1.92	0.69
2:A:232:LYS:O	2:A:236:VAL:HG23	1.94	0.68
1:B:109:DT:C2'	1:B:110:DA:C5'	2.73	0.67
2:A:290:ARG:HG3	2:A:290:ARG:HH11	1.60	0.66
2:A:155:SER:HB2	2:A:325:TYR:CD2	2.31	0.66
1:C:116:DA:C5'	2:A:192:ARG:HD3	2.25	0.65
1:B:101:DC:C2'	1:B:102:DG:C8	2.79	0.65
1:B:105:DT:C5'	2:A:288:ILE:HD13	2.25	0.65
2:A:266:LEU:HD22	2:A:287:LEU:HB2	1.80	0.63
1:B:105:DT:H5''	2:A:288:ILE:CD1	2.27	0.62
2:A:270:VAL:HG21	2:A:279:TYR:HD2	1.65	0.62
2:A:209:ILE:HD13	2:A:215:MET:HG3	1.82	0.61
1:C:123:DC:C2'	1:C:124:DG:N7	2.63	0.61
2:A:172:CYS:O	2:A:174:LEU:HD13	2.01	0.60
1:B:106:DA:H5''	2:A:297:VAL:HG21	1.84	0.59
2:A:279:TYR:CE1	2:A:281:PRO:HB3	2.37	0.59
1:C:113:DC:H2''	1:C:114:DG:OP2	2.02	0.59
2:A:177:LYS:O	2:A:181:LEU:HG	2.03	0.59
2:A:209:ILE:CD1	2:A:215:MET:SD	2.91	0.58
1:C:123:DC:H5'	1:C:123:DC:H6	1.69	0.57
1:B:102:DG:C2'	1:B:103:DT:C5'	2.77	0.57
2:A:314:ARG:HG2	2:A:318:TYR:CE2	2.38	0.57
2:A:164:ILE:HD13	2:A:254:MET:HG2	1.86	0.57
2:A:192:ARG:NH2	2:A:197:ILE:HD12	2.17	0.57
1:C:123:DC:H2''	1:C:124:DG:N7	2.21	0.56
2:A:260:VAL:HG22	2:A:304:GLY:CA	2.36	0.55
2:A:269:LEU:HD13	2:A:287:LEU:CD2	2.35	0.55
2:A:188:TYR:CZ	2:A:190:PRO:HD3	2.43	0.54
2:A:270:VAL:HG12	2:A:274:GLN:HE22	1.72	0.54
1:C:120:DA:H2''	1:C:121:DT:O5'	2.08	0.53
2:A:262:PHE:CZ	2:A:329:LYS:HA	2.43	0.53
2:A:209:ILE:HD11	2:A:215:MET:SD	2.48	0.53
2:A:291:MET:HE3	2:A:298:LEU:HD11	1.90	0.52
2:A:175:ASP:O	2:A:179:ILE:HG13	2.10	0.52
2:A:191:LYS:NZ	2:A:192:ARG:HB3	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:116:DA:H4'	2:A:192:ARG:CD	2.39	0.52
1:B:105:DT:O2	2:A:299:LEU:HD21	2.08	0.52
2:A:250:LYS:HD3	2:A:252:GLN:HE22	1.75	0.52
1:B:106:DA:H5''	2:A:297:VAL:CG2	2.39	0.52
2:A:209:ILE:HD13	2:A:215:MET:CG	2.40	0.52
2:A:227:ARG:O	2:A:230:ALA:HB3	2.10	0.52
2:A:278:SER:O	2:A:287:LEU:HD12	2.10	0.51
2:A:187:GLU:HB2	2:A:197:ILE:HB	1.92	0.50
2:A:157:ILE:HG22	2:A:158:VAL:N	2.26	0.50
1:C:123:DC:H2''	1:C:124:DG:H8	1.67	0.50
2:A:260:VAL:O	2:A:261:LYS:HB2	2.11	0.49
1:B:102:DG:C2'	1:B:103:DT:H5''	2.41	0.49
1:B:107:DT:C5	1:B:108:DA:N6	2.81	0.48
1:C:116:DA:H4'	2:A:192:ARG:HD3	1.95	0.48
1:C:123:DC:H5'	1:C:123:DC:C6	2.47	0.48
2:A:236:VAL:O	2:A:240:LEU:HD13	2.13	0.48
1:B:108:DA:C2'	1:B:109:DT:H5'	2.36	0.48
1:B:107:DT:H4'	2:A:252:GLN:O	2.14	0.48
2:A:259:ASP:HB3	2:A:261:LYS:NZ	2.29	0.47
2:A:209:ILE:HD13	2:A:215:MET:SD	2.54	0.47
1:B:102:DG:C2'	1:B:103:DT:H5'	2.28	0.47
2:A:324:ILE:O	2:A:328:LEU:HG	2.15	0.47
2:A:233:TYR:O	2:A:237:VAL:HG23	2.14	0.47
2:A:290:ARG:HH11	2:A:290:ARG:CG	2.27	0.47
1:B:108:DA:H4'	2:A:252:GLN:HG3	1.97	0.47
2:A:227:ARG:HD3	2:A:249:PHE:CD2	2.50	0.47
1:B:104:DA:OP1	2:A:283:LEU:HD21	2.16	0.46
2:A:260:VAL:HG22	2:A:304:GLY:HA3	1.97	0.46
1:C:116:DA:O5'	2:A:192:ARG:HD3	2.16	0.46
2:A:238:GLN:HE21	2:A:244:ALA:HB3	1.81	0.46
2:A:269:LEU:HD13	2:A:287:LEU:HD23	1.96	0.46
2:A:220:ALA:O	2:A:221:LYS:HG2	2.16	0.45
2:A:252:GLN:HA	2:A:252:GLN:OE1	2.15	0.45
1:B:108:DA:C4'	2:A:252:GLN:HG3	2.46	0.45
2:A:254:MET:HE2	2:A:254:MET:HB3	1.73	0.45
1:C:119:DT:H6	1:C:119:DT:H2'	1.61	0.45
2:A:210:PHE:HB2	2:A:214:LYS:O	2.17	0.45
2:A:272:THR:OG1	2:A:273:HIS:HD2	1.99	0.45
2:A:206:THR:HG22	2:A:207:ALA:N	2.31	0.44
2:A:284:PHE:CG	2:A:285:PRO:HD2	2.52	0.44
1:C:116:DA:C2	2:A:208:LEU:HD13	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:270:VAL:HG22	2:A:287:LEU:HD13	2.00	0.44
2:A:235:ARG:HG2	2:A:235:ARG:HH11	1.82	0.44
2:A:163:ASN:C	2:A:164:ILE:HG12	2.43	0.43
2:A:200:ILE:HG23	2:A:233:TYR:CE1	2.53	0.43
2:A:224:GLU:O	2:A:225:GLN:C	2.59	0.43
1:C:116:DA:C4'	2:A:192:ARG:HD3	2.49	0.43
2:A:165:VAL:HG13	2:A:218:THR:HG22	2.00	0.43
2:A:159:PRO:HA	2:A:258:CYS:HB3	2.00	0.43
1:C:118:DA:OP1	2:A:199:ARG:NH1	2.51	0.43
2:A:260:VAL:HG22	2:A:304:GLY:O	2.19	0.43
2:A:266:LEU:HB3	2:A:279:TYR:CE2	2.53	0.43
2:A:284:PHE:CD1	2:A:285:PRO:HD2	2.53	0.43
2:A:219:GLY:O	2:A:220:ALA:O	2.36	0.43
2:A:261:LYS:O	2:A:262:PHE:HB3	2.18	0.42
2:A:319:GLU:O	2:A:320:ALA:C	2.61	0.42
2:A:170:LEU:HD12	2:A:174:LEU:HD21	2.00	0.42
1:C:123:DC:H2'	1:C:124:DG:N7	2.34	0.42
2:A:191:LYS:HZ2	2:A:192:ARG:HB3	1.84	0.42
2:A:298:LEU:HD12	2:A:298:LEU:N	2.35	0.42
2:A:184:ARG:C	2:A:186:ALA:H	2.28	0.42
2:A:266:LEU:HD22	2:A:287:LEU:CB	2.48	0.42
1:B:106:DA:N7	1:B:107:DT:C4	2.88	0.42
2:A:195:ALA:HB2	2:A:210:PHE:CE1	2.55	0.42
2:A:185:ASN:OD1	2:A:201:ARG:NH1	2.49	0.41
1:C:116:DA:C2	2:A:193:PHE:CZ	3.08	0.41
2:A:188:TYR:CE2	2:A:190:PRO:HD3	2.55	0.41
2:A:189:ASN:C	2:A:191:LYS:H	2.28	0.41
2:A:297:VAL:O	2:A:309:THR:HG22	2.20	0.41
2:A:189:ASN:OD1	2:A:189:ASN:N	2.54	0.41
2:A:324:ILE:HG13	2:A:328:LEU:HG	2.03	0.41
2:A:265:ARG:HB2	2:A:267:GLU:OE2	2.20	0.41
2:A:177:LYS:HD2	2:A:181:LEU:HD21	2.03	0.41
2:A:221:LYS:HG2	2:A:221:LYS:HZ3	1.65	0.41
2:A:202:GLU:HA	2:A:203:PRO:C	2.46	0.41
2:A:329:LYS:HE2	2:A:329:LYS:HB3	1.82	0.40
2:A:279:TYR:CZ	2:A:281:PRO:HB3	2.57	0.40
2:A:206:THR:CG2	2:A:207:ALA:N	2.84	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	
2	A	178/185 (96%)	150 (84%)	24 (14%)	4 (2%)	5	20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	220	ALA
2	A	172	CYS
2	A	184	ARG
2	A	262	PHE

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	
2	A	155/158 (98%)	135 (87%)	20 (13%)	4	13

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	159	PRO
2	A	161	LEU
2	A	174	LEU
2	A	178	THR
2	A	184	ARG
2	A	185	ASN
2	A	189	ASN

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Mol	Chain	Res	Type
2	A	192	ARG
2	A	196	VAL
2	A	211	SER
2	A	214	LYS
2	A	228	LEU
2	A	240	LEU
2	A	267	GLU
2	A	269	LEU
2	A	290	ARG
2	A	292	ILE
2	A	293	LYS
2	A	299	LEU
2	A	302	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	162	GLN
2	A	273	HIS

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	B	12/12 (100%)	-0.65	0 100 100	15, 28, 35, 41	0
1	C	12/12 (100%)	-0.75	0 100 100	20, 27, 30, 32	0
2	A	180/185 (97%)	-0.52	0 100 100	9, 25, 42, 52	0
All	All	204/209 (97%)	-0.54	0 100 100	9, 25, 42, 52	0

There are no RSRZ outliers to report.

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.