



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:50 AM UTC

PDB ID : 1D3U / pdb_00001d3u
Title : TATA-BINDING P R O T E I N / T R A N S C R I P T I O N F A C T O R
(II)B/BRE+TATA-BOX COMPLEX FROM PYROCOCCLUS WOESEI
Authors : Littlefield, O.; Korkhin, Y.; Sigler, P.B.
Deposited on : 1999-10-01
Resolution : 2.40 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

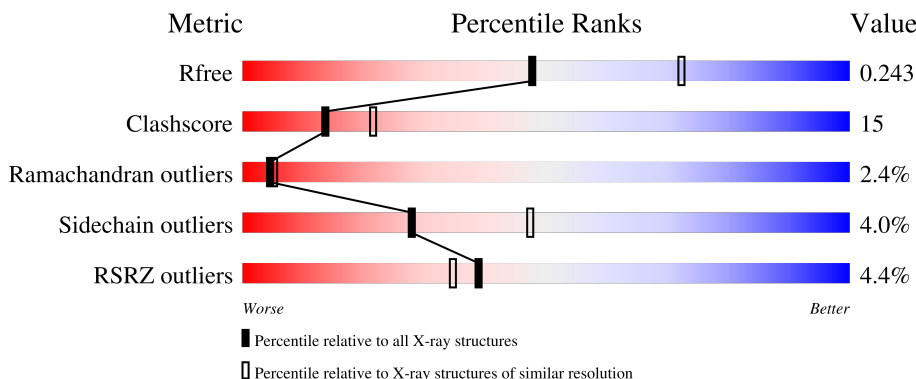
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	C	24	
2	D	23	
3	A	181	
4	B	201	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4147 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a DNA chain called DNA 24-MER: BRE+TATA-BOX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	24	Total	C	N	O	P	0	0	0
			493	239	91	140	23			

- Molecule 2 is a DNA chain called DNA 23-MER: BRE+TATA-BOX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	23	Total	C	N	O	P	0	0	0
			464	227	76	139	22			

- Molecule 3 is a protein called TATA-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	181	Total	C	N	O	S	0	0	0
			1415	910	235	263	7			

- Molecule 4 is a protein called TRANSCRIPTION INITIATION FACTOR IIB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	201	Total	C	N	O	S	0	0	0
			1595	1006	296	289	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1100	MET	ARG	engineered mutation	UNP P61999

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	16	Total	O	0	0
			16	16		

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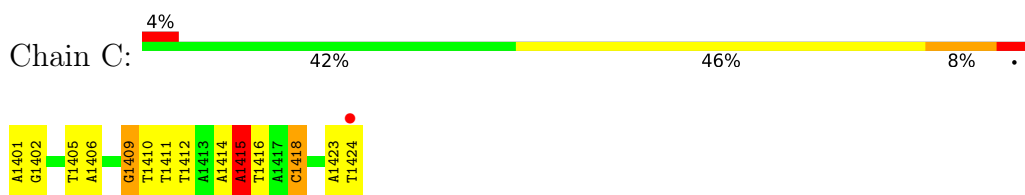
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	19	Total 19	O 19	0	0
5	A	72	Total 72	O 72	0	0
5	B	73	Total 73	O 73	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

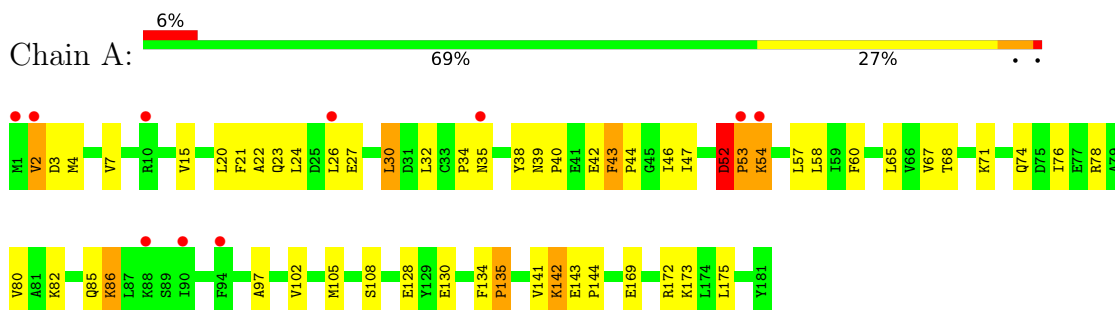
- Molecule 1: DNA 24-MER: BRE+TATA-BOX



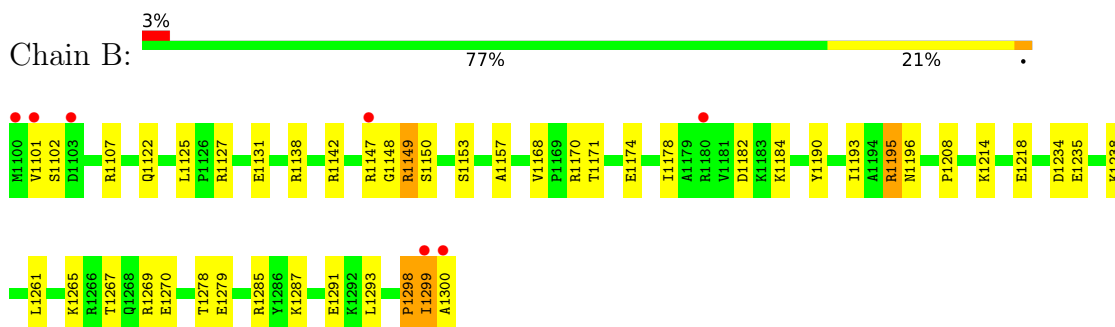
- Molecule 2: DNA 23-MER: BRE+TATA-BOX



- Molecule 3: TATA-BINDING PROTEIN



- Molecule 4: TRANSCRIPTION INITIATION FACTOR IIB



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	174.20Å 174.20Å 174.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 20.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.7 (20.00-2.40) 98.6 (20.00-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 2.42Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.206 , 0.245 0.205 , 0.243	Depositor DCC
R_{free} test set	3325 reflections (9.83%)	wwPDB-VP
Wilson B-factor (Å ²)	50.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.021 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4147	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.61% 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	C	0.39	0/554	1.09	1/854 (0.1%)
2	D	0.42	0/518	1.04	1/797 (0.1%)
3	A	0.73	0/1438	1.14	11/1939 (0.6%)
4	B	0.84	0/1612	1.07	5/2165 (0.2%)
All	All	0.71	0/4122	1.09	18/5755 (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	C	0	4
2	D	0	3
All	All	0	7

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	52	ASP	CA-C-N	-10.61	108.76	119.90
3	A	52	ASP	C-N-CA	-10.61	108.76	119.90
3	A	43	PHE	CA-C-N	7.96	127.62	119.82
3	A	43	PHE	C-N-CA	7.96	127.62	119.82
4	B	1168	VAL	N-CA-C	-7.11	100.51	107.55
3	A	130	GLU	CA-C-N	6.87	126.55	119.82
3	A	130	GLU	C-N-CA	6.87	126.55	119.82
3	A	53	PRO	N-CA-C	-6.82	100.50	110.80
3	A	108	SER	N-CA-C	-6.11	100.67	110.14
3	A	135	PRO	N-CA-C	5.91	121.25	113.86
2	D	1432	DT	N1-C1'-C2'	-5.74	104.90	113.50
3	A	169	GLU	N-CA-C	-5.55	105.13	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1101	VAL	N-CA-C	-5.46	97.99	109.34
4	B	1196	ASN	N-CA-C	5.45	119.19	112.54
4	B	1125	LEU	CA-C-N	5.27	125.21	119.78
4	B	1125	LEU	C-N-CA	5.27	125.21	119.78
1	C	1415	DA	N9-C1'-C2'	-5.24	105.64	113.50
3	A	97	ALA	N-CA-C	-5.15	103.54	109.93

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1409	DG	Sidechain
1	C	1414	DA	Sidechain
1	C	1415	DA	Sidechain
1	C	1418	DC	Sidechain
2	D	1431	DG	Sidechain
2	D	1432	DT	Sidechain
2	D	1445	DC	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	493	0	275	14	0
2	D	464	0	266	25	1
3	A	1415	0	1473	41	1
4	B	1595	0	1695	37	0
5	A	72	0	0	0	0
5	B	73	0	0	6	0
5	C	16	0	0	0	0
5	D	19	0	0	1	0
All	All	4147	0	3709	109	2

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 15.

All (109) 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)
4:B:1149:ARG:HH11	4:B:1149:ARG:HB3	0.99	1.11
1:C:1415:DA:H2''	1:C:1416:DT:H5'	1.31	1.11
3:A:15:VAL:HG22	3:A:68:THR:HG22	1.45	0.98
4:B:1149:ARG:HB3	4:B:1149:ARG:NH1	1.82	0.93
2:D:1444:DA:H2''	2:D:1445:DC:H5''	1.48	0.91
4:B:1149:ARG:HH11	4:B:1149:ARG:CB	1.83	0.91
1:C:1415:DA:H2''	1:C:1416:DT:C5'	2.03	0.88
4:B:1195:ARG:HH11	4:B:1195:ARG:HG3	1.40	0.87
3:A:54:LYS:HD2	3:A:54:LYS:O	1.74	0.86
4:B:1299:ILE:HG22	4:B:1300:ALA:H	1.38	0.86
3:A:52:ASP:HB3	3:A:53:PRO:HD3	1.59	0.83
2:D:1429:DA:H2''	2:D:1430:DA:H8	1.48	0.78
2:D:1434:DT:H2''	2:D:1435:DT:H5'	1.67	0.74
2:D:1445:DC:H2'	2:D:1446:DT:H72	1.69	0.73
4:B:1214:LYS:HE2	4:B:1218:GLU:OE1	1.90	0.71
1:C:1416:DT:H1'	3:A:60:PHE:CE2	2.26	0.71
3:A:38:TYR:CZ	3:A:40:PRO:HG3	2.26	0.70
1:C:1423:DA:H2''	1:C:1424:DT:H5'	1.72	0.70
2:D:1444:DA:C2'	2:D:1445:DC:H5''	2.22	0.69
2:D:1426:DT:H2''	2:D:1427:DA:C8	2.27	0.69
2:D:1429:DA:H2''	2:D:1430:DA:C8	2.29	0.68
1:C:1418:DC:H5'	3:A:44:PRO:HB3	1.78	0.65
4:B:1195:ARG:HG3	4:B:1195:ARG:NH1	2.11	0.65
3:A:2:VAL:O	3:A:4:MET:HG3	1.97	0.64
3:A:32:LEU:HD22	3:A:86:LYS:HB3	1.79	0.63
3:A:23:GLN:O	3:A:24:LEU:HD23	1.99	0.63
4:B:1142:ARG:HG2	4:B:1142:ARG:HH11	1.64	0.63
4:B:1298:PRO:O	4:B:1299:ILE:HD13	2.00	0.62
4:B:1149:ARG:N	4:B:1149:ARG:HD2	2.13	0.62
3:A:76:ILE:O	3:A:80:VAL:HG23	1.99	0.61
3:A:52:ASP:HB3	3:A:53:PRO:CD	2.29	0.61
2:D:1445:DC:H2'	2:D:1446:DT:C7	2.30	0.61
3:A:82:LYS:N	3:A:82:LYS:HD3	2.15	0.61
3:A:134:PHE:CD1	3:A:135:PRO:HD2	2.36	0.60
3:A:47:ILE:HD13	3:A:58:LEU:HD22	1.84	0.60
2:D:1427:DA:H1'	2:D:1428:DT:H5''	1.83	0.60
3:A:54:LYS:HE2	3:A:71:LYS:HG2	1.83	0.60
2:D:1440:DC:H2'	2:D:1441:DT:H72	1.84	0.59
4:B:1182:ASP:OD2	4:B:1184:LYS:HB3	2.02	0.59
4:B:1299:ILE:HG22	4:B:1300:ALA:N	2.14	0.59
2:D:1447:DC:H2''	2:D:1448:DT:H5''	1.84	0.58
3:A:26:LEU:HB3	3:A:46:ILE:HD11	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1405:DT:H2''	1:C:1406:DA:C8	2.38	0.58
4:B:1150:SER:OG	4:B:1153:SER:HB2	2.04	0.57
4:B:1171:THR:OG1	4:B:1174:GLU:HG3	2.05	0.56
4:B:1287:LYS:O	4:B:1291:GLU:HG2	2.05	0.56
3:A:7:VAL:HG21	3:A:175:LEU:HD11	1.88	0.55
1:C:1416:DT:H1'	3:A:60:PHE:CZ	2.41	0.54
4:B:1142:ARG:HG2	4:B:1142:ARG:NH1	2.23	0.54
4:B:1234:ASP:O	4:B:1238:LYS:HG3	2.07	0.54
2:D:1428:DT:H6	2:D:1428:DT:H5'	1.72	0.53
2:D:1446:DT:H71	5:D:250:HOH:O	2.09	0.52
3:A:57:LEU:HD23	3:A:67:VAL:HG13	1.90	0.52
3:A:15:VAL:CG2	3:A:68:THR:HG22	2.31	0.51
4:B:1149:ARG:HD2	4:B:1149:ARG:H	1.76	0.51
2:D:1430:DA:H4'	4:B:1148:GLY:O	2.11	0.51
4:B:1147:ARG:HG3	5:B:179:HOH:O	2.11	0.51
4:B:1102:SER:O	4:B:1107:ARG:NH1	2.44	0.51
3:A:141:VAL:HG13	3:A:173:LYS:HE2	1.93	0.50
1:C:1411:DT:H2''	1:C:1412:DT:H5'	1.93	0.50
4:B:1122:GLN:HG2	4:B:1178:ILE:HG21	1.94	0.50
2:D:1430:DA:H5'	4:B:1148:GLY:HA3	1.94	0.50
3:A:143:GLU:HA	3:A:144:PRO:C	2.36	0.49
1:C:1415:DA:C2'	1:C:1416:DT:H5'	2.22	0.49
3:A:57:LEU:CD2	3:A:67:VAL:HG13	2.42	0.48
3:A:20:LEU:O	3:A:21:PHE:HB2	2.12	0.48
4:B:1170:ARG:NE	5:B:222:HOH:O	2.10	0.48
2:D:1431:DG:H2'	2:D:1432:DT:H72	1.96	0.48
3:A:52:ASP:CB	3:A:53:PRO:HD3	2.37	0.47
4:B:1157:ALA:HB1	4:B:1190:TYR:HA	1.96	0.47
2:D:1430:DA:H2''	2:D:1431:DG:H8	1.79	0.47
3:A:52:ASP:OD2	3:A:78:ARG:NH2	2.40	0.47
3:A:15:VAL:HG13	3:A:68:THR:CG2	2.44	0.47
3:A:39:ASN:HB2	3:A:42:GLU:HG2	1.97	0.47
1:C:1401:DA:H2''	1:C:1402:DG:O5'	2.15	0.47
1:C:1418:DC:H5'	3:A:44:PRO:CB	2.44	0.47
3:A:30:LEU:HD21	3:A:38:TYR:HB2	1.97	0.46
3:A:142:LYS:H	3:A:142:LYS:HD2	1.80	0.46
4:B:1170:ARG:NH2	5:B:222:HOH:O	2.45	0.46
1:C:1405:DT:H2''	1:C:1406:DA:N7	2.31	0.46
2:D:1432:DT:H2''	2:D:1433:DA:O5'	2.15	0.46
2:D:1447:DC:C2'	2:D:1448:DT:H5''	2.45	0.45
2:D:1428:DT:H5'	2:D:1428:DT:C6	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1127:ARG:HG3	4:B:1131:GLU:OE2	2.15	0.45
3:A:52:ASP:O	3:A:53:PRO:C	2.56	0.45
3:A:52:ASP:CB	3:A:53:PRO:CD	2.95	0.45
2:D:1448:DT:H6	2:D:1448:DT:H5'	1.82	0.45
4:B:1279:GLU:HG3	5:B:206:HOH:O	2.15	0.45
3:A:57:LEU:O	3:A:58:LEU:HD23	2.17	0.44
2:D:1440:DC:H2'	2:D:1441:DT:C7	2.46	0.44
2:D:1429:DA:C2'	2:D:1430:DA:C8	3.00	0.44
4:B:1138:ARG:HD2	5:B:127:HOH:O	2.17	0.43
4:B:1149:ARG:NH1	5:B:256:HOH:O	2.51	0.43
3:A:82:LYS:HD3	3:A:82:LYS:H	1.84	0.43
3:A:134:PHE:CG	3:A:135:PRO:HD2	2.54	0.43
1:C:1409:DG:C8	1:C:1410:DT:H72	2.54	0.43
3:A:128:GLU:OE1	4:B:1195:ARG:HD3	2.19	0.42
3:A:43:PHE:CD1	3:A:44:PRO:HD2	2.53	0.42
1:C:1410:DT:H2''	1:C:1411:DT:H5'	2.02	0.41
4:B:1195:ARG:NH1	4:B:1195:ARG:CG	2.76	0.41
4:B:1285:ARG:HA	4:B:1285:ARG:HD3	1.86	0.41
2:D:1426:DT:C2'	2:D:1427:DA:C8	3.01	0.41
2:D:1442:DT:H2'	4:B:1278:THR:HG21	2.02	0.41
3:A:20:LEU:HA	3:A:20:LEU:HD23	1.84	0.41
3:A:102:VAL:HG11	3:A:105:MET:HE3	2.02	0.41
3:A:54:LYS:H	3:A:54:LYS:HG3	1.67	0.40
4:B:1267:THR:OG1	4:B:1270:GLU:HG3	2.21	0.40
4:B:1235:GLU:HG3	4:B:1293:LEU:HD22	2.04	0.40
4:B:1269:ARG:HG2	4:B:1269:ARG:HH11	1.86	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:1426:DT:O4'	2:D:1426:DT:O4'[16_546]	2.09	0.11
3:A:172:ARG:NH1	3:A:172:ARG:NH1[15_556]	2.14	0.06

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	A	179/181 (99%)	163 (91%)	10 (6%)	6 (3%)	3	2
4	B	199/201 (99%)	193 (97%)	3 (2%)	3 (2%)	8	12
All	All	378/382 (99%)	356 (94%)	13 (3%)	9 (2%)	4	5

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	52	ASP
4	B	1265	LYS
4	B	1298	PRO
3	A	30	LEU
3	A	34	PRO
3	A	22	ALA
3	A	85	GLN
3	A	86	LYS
4	B	1299	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	A	159/159 (100%)	151 (95%)	8 (5%)	22	38
4	B	167/167 (100%)	162 (97%)	5 (3%)	36	58
All	All	326/326 (100%)	313 (96%)	13 (4%)	28	47

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	2	VAL
3	A	3	ASP
3	A	27	GLU

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Mol	Chain	Res	Type
3	A	35	ASN
3	A	54	LYS
3	A	65	LEU
3	A	74	GLN
3	A	142	LYS
4	B	1149	ARG
4	B	1193	ILE
4	B	1195	ARG
4	B	1208	PRO
4	B	1261	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
4	B	1122	GLN
4	B	1128	HIS
4	B	1213	ASN

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 ⓘ

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	C	24/24 (100%)	0.07	1 (4%) 40 36	34, 59, 87, 100	0
2	D	23/23 (100%)	0.15	1 (4%) 40 36	34, 59, 93, 100	0
3	A	181/181 (100%)	0.12	10 (5%) 30 27	27, 53, 97, 100	0
4	B	201/201 (100%)	-0.14	7 (3%) 47 43	25, 44, 77, 100	0
All	All	429/429 (100%)	-0.00	19 (4%) 39 34	25, 49, 95, 100	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	B	1100	MET	12.6
4	B	1299	ILE	7.7
4	B	1300	ALA	4.9
3	A	1	MET	3.5
3	A	2	VAL	3.3
1	C	1424	DT	3.1
3	A	10	ARG	2.9
4	B	1101	VAL	2.9
4	B	1103	ASP	2.8
2	D	1426	DT	2.8
3	A	54	LYS	2.5
4	B	1147	ARG	2.3
3	A	26	LEU	2.2
3	A	90	ILE	2.2
3	A	88	LYS	2.2
3	A	53	PRO	2.2
4	B	1180	ARG	2.1
3	A	94	PHE	2.0
3	A	35	ASN	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.