



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:26 PM UTC

PDB ID : 1BCM / pdb_00001bcm
Title : BACTERIOPHAGE MU TRANSPOSASE CORE DOMAIN WITH 2
MONOMERS PER ASYMMETRIC UNIT
Authors : Rice, P.A.; Mizuuchi, K.
Deposited on : 1995-05-26
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

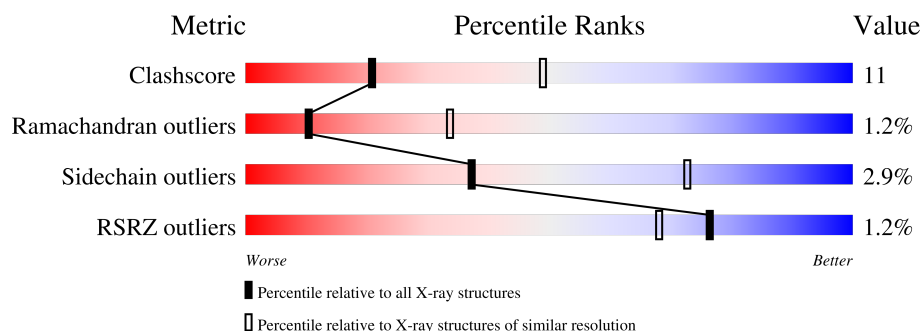
1 Overall quality at a glance

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))
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	327	68% 23% • 7%
1	B	327	67% 25% • 7%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5815 atoms, of which 1059 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.

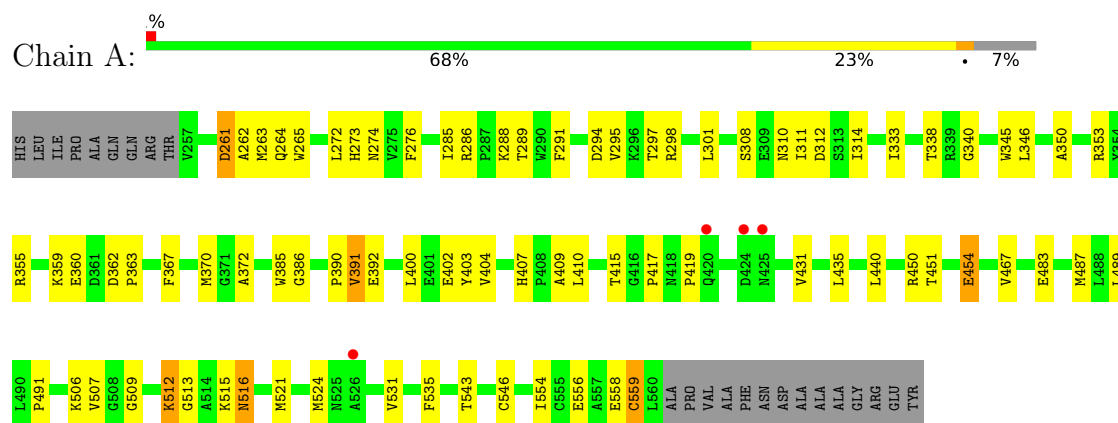
- Molecule 1 is a protein called BACTERIOPHAGE MU TRANSPOSASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	304	Total	C	H	N	O	S	0	0	0
			2902	1520	525	412	431	14			
1	B	303	Total	C	H	N	O	S	0	0	0
			2913	1520	534	416	429	14			

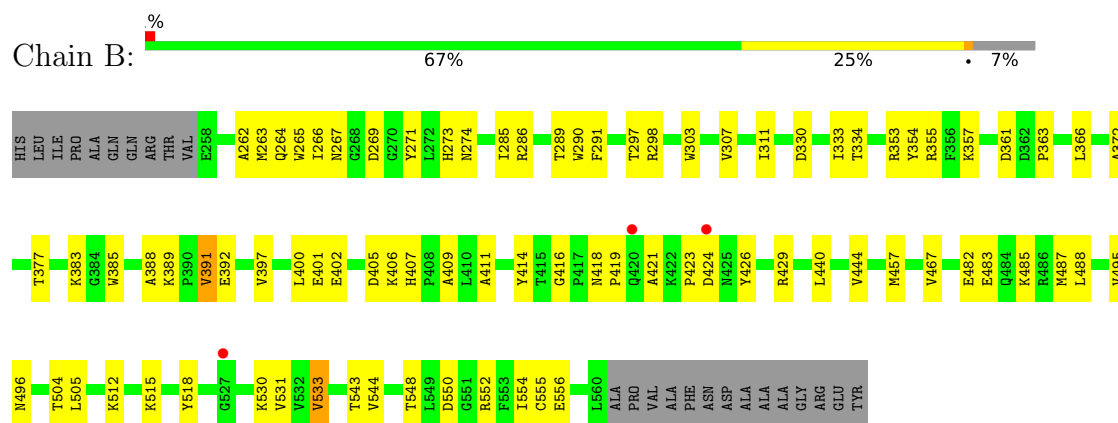
3 Residue-property plots

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: BACTERIOPHAGE MU TRANSPOSASE



• Molecule 1: BACTERIOPHAGE MU TRANSPOSASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	172.20Å 52.20Å 101.60Å 90.00° 115.50° 90.00°	Depositor
Resolution (Å)	8.00 – 2.80 8.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.0 (8.00-2.80) 89.5 (8.00-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.69Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.214 , 0.276 0.203 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 70.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5815	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.47	0/2436	0.90	2/3300 (0.1%)
1	B	0.49	1/2438 (0.0%)	0.92	2/3300 (0.1%)
All	All	0.48	1/4874 (0.0%)	0.91	4/6600 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	266	ILE	CA-CB	5.42	1.61	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	340	GLY	N-CA-C	-6.92	105.27	115.72
1	B	411	ALA	N-CA-C	5.74	119.10	111.75
1	B	533	VAL	N-CA-C	5.51	115.80	107.80
1	A	295	VAL	N-CA-C	5.07	115.79	110.62

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	2377	525	2304	53	0
1	B	2379	534	2317	52	0
All	All	4756	1059	4621	104	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 11.

All (104) 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:286:ARG:NH2	1:B:421:ALA:HB3	1.95	0.81
1:B:363:PRO:HA	1:B:366:LEU:HG	1.75	0.69
1:B:504:THR:HG21	1:B:515:LYS:HE2	1.74	0.69
1:B:273:HIS:HE1	1:B:289:THR:OG1	1.77	0.68
1:A:531:VAL:HG21	1:A:546:CYS:HB3	1.78	0.66
1:A:350:ALA:O	1:A:353:ARG:HG3	1.96	0.66
1:B:285:ILE:HG12	1:B:307:VAL:HG13	1.77	0.65
1:A:391:VAL:HG23	1:A:392:GLU:H	1.62	0.65
1:B:286:ARG:HH21	1:B:421:ALA:HB3	1.60	0.65
1:A:273:HIS:HE1	1:A:289:THR:OG1	1.80	0.63
1:B:333:ILE:HD12	1:B:372:ALA:HB1	1.80	0.62
1:B:426:TYR:HA	1:B:429:ARG:HB2	1.79	0.62
1:B:407:HIS:HD2	1:B:409:ALA:H	1.47	0.62
1:A:431:VAL:CG2	1:A:435:LEU:HD23	2.30	0.61
1:B:290:TRP:O	1:B:303:TRP:HA	2.01	0.61
1:A:288:LYS:HD3	1:A:308:SER:HA	1.82	0.60
1:B:391:VAL:HG23	1:B:392:GLU:H	1.66	0.60
1:A:311:ILE:HG21	1:A:363:PRO:HG3	1.84	0.59
1:B:267:ASN:HD21	1:B:389:LYS:HE3	1.68	0.59
1:A:355:ARG:HE	1:A:506:LYS:HB3	1.68	0.58
1:A:262:ALA:O	1:A:263:MET:HB2	2.04	0.57
1:B:543:THR:CG2	1:B:556:GLU:HG3	2.34	0.57
1:A:297:THR:HB	1:A:467:VAL:HG23	1.86	0.56
1:A:531:VAL:CG2	1:A:546:CYS:HB3	2.35	0.56
1:B:548:THR:HG22	1:B:550:ASP:H	1.71	0.56
1:A:512:LYS:HD3	1:A:512:LYS:H	1.71	0.56
1:B:273:HIS:HD2	1:B:405:ASP:OD1	1.89	0.55
1:B:271:TYR:HB2	1:B:397:VAL:HG11	1.90	0.53
1:B:543:THR:HG22	1:B:544:VAL:N	2.24	0.53
1:A:274:ASN:ND2	1:A:286:ARG:HH21	2.07	0.53
1:A:346:LEU:HD13	1:B:552:ARG:HH21	1.74	0.52
1:A:417:PRO:O	1:A:419:PRO:HD3	2.10	0.52
1:A:276:PHE:CD2	1:A:415:THR:HG21	2.46	0.51
1:B:262:ALA:O	1:B:263:MET:HB2	2.10	0.51
1:B:414:TYR:CZ	1:B:416:GLY:HA2	2.46	0.51
1:A:311:ILE:CG2	1:A:363:PRO:HG3	2.41	0.50
1:A:311:ILE:HG21	1:A:363:PRO:CG	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:TRP:CZ3	1:A:385:TRP:HB3	2.47	0.50
1:B:264:GLN:OE1	1:B:385:TRP:CZ2	2.65	0.50
1:A:333:ILE:HG12	1:A:372:ALA:HB1	1.94	0.49
1:A:404:VAL:O	1:A:410:LEU:HD12	2.12	0.49
1:A:298:ARG:NH1	1:A:390:PRO:HD3	2.26	0.49
1:A:274:ASN:HA	1:A:286:ARG:HH21	1.78	0.49
1:A:507:VAL:O	1:A:513:GLY:HA2	2.13	0.49
1:B:505:LEU:HD12	1:B:518:TYR:CE2	2.47	0.49
1:A:291:PHE:CZ	1:A:400:LEU:HD13	2.47	0.49
1:B:496:ASN:HD22	1:B:496:ASN:N	2.11	0.48
1:A:407:HIS:HD2	1:A:409:ALA:H	1.60	0.48
1:A:489:LEU:HB3	1:A:535:PHE:O	2.13	0.48
1:B:401:GLU:HG3	1:B:405:ASP:HB2	1.96	0.48
1:A:294:ASP:HB2	1:A:301:LEU:HD11	1.95	0.47
1:B:311:ILE:HG12	1:B:363:PRO:HG3	1.96	0.47
1:B:483:GLU:O	1:B:487:MET:HG2	2.15	0.47
1:B:334:THR:HG23	1:B:377:THR:O	2.15	0.46
1:A:516:ASN:OD1	1:A:516:ASN:N	2.49	0.46
1:B:297:THR:HB	1:B:467:VAL:HG23	1.97	0.46
1:A:261:ASP:HB2	1:A:264:GLN:OE1	2.16	0.46
1:A:359:LYS:HD3	1:A:491:PRO:HB2	1.98	0.46
1:B:482:GLU:HA	1:B:485:LYS:HE3	1.98	0.46
1:B:303:TRP:CZ3	1:B:444:VAL:HG11	2.51	0.46
1:A:273:HIS:CE1	1:A:289:THR:OG1	2.65	0.45
1:B:269:ASP:HB3	1:B:397:VAL:HG21	1.97	0.45
1:B:297:THR:O	1:B:298:ARG:HB2	2.16	0.45
1:B:273:HIS:CE1	1:B:289:THR:OG1	2.63	0.45
1:A:515:LYS:HE3	1:A:515:LYS:HB3	1.85	0.45
1:A:483:GLU:O	1:A:487:MET:HG2	2.17	0.45
1:B:286:ARG:HH22	1:B:421:ALA:HB3	1.80	0.44
1:B:495:VAL:O	1:B:530:LYS:HA	2.17	0.44
1:A:355:ARG:NE	1:A:506:LYS:HE3	2.32	0.44
1:B:383:LYS:HD2	1:B:385:TRP:CH2	2.53	0.44
1:B:423:PRO:O	1:B:426:TYR:N	2.50	0.44
1:A:310:ASN:HD21	1:A:312:ASP:HB2	1.83	0.44
1:B:265:TRP:CZ2	1:B:388:ALA:HB2	2.53	0.44
1:B:402:GLU:HA	1:B:406:LYS:HE3	1.99	0.44
1:A:272:LEU:HD13	1:A:288:LYS:HE3	1.99	0.44
1:B:457:MET:HA	1:B:457:MET:HE2	2.00	0.43
1:A:391:VAL:HG23	1:A:392:GLU:N	2.31	0.43
1:B:504:THR:CG2	1:B:515:LYS:HE2	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:533:VAL:HG13	1:B:544:VAL:HB	2.00	0.43
1:A:367:PHE:N	1:A:367:PHE:CD1	2.85	0.43
1:A:370:MET:HE2	1:A:370:MET:HB3	1.84	0.43
1:A:531:VAL:HG23	1:A:554:ILE:HD12	2.00	0.43
1:B:298:ARG:HA	1:B:298:ARG:HD3	1.76	0.43
1:A:274:ASN:ND2	1:A:286:ARG:NH2	2.67	0.42
1:B:274:ASN:O	1:B:414:TYR:HD1	2.02	0.42
1:B:554:ILE:O	1:B:555:CYS:HB3	2.19	0.42
1:B:353:ARG:HG3	1:B:354:TYR:CD2	2.55	0.42
1:A:450:ARG:HG2	1:A:451:THR:O	2.20	0.42
1:A:353:ARG:O	1:A:509:GLY:HA3	2.20	0.42
1:B:264:GLN:HG3	1:B:330:ASP:O	2.19	0.41
1:A:310:ASN:ND2	1:A:312:ASP:HB2	2.34	0.41
1:A:512:LYS:H	1:A:512:LYS:CD	2.34	0.41
1:A:543:THR:HG21	1:A:556:GLU:OE2	2.20	0.41
1:B:291:PHE:CZ	1:B:400:LEU:HD13	2.55	0.41
1:A:276:PHE:CE2	1:A:415:THR:HG21	2.55	0.41
1:A:314:ILE:HD11	1:A:345:TRP:HB3	2.03	0.41
1:A:558:GLU:HG2	1:A:559:CYS:N	2.36	0.41
1:B:505:LEU:HD12	1:B:518:TYR:HE2	1.85	0.41
1:B:418:ASN:HA	1:B:419:PRO:HD2	1.90	0.41
1:B:488:LEU:HD23	1:B:488:LEU:HA	1.93	0.41
1:A:285:ILE:HD12	1:A:285:ILE:C	2.47	0.40
1:A:402:GLU:HG3	1:A:403:TYR:CD2	2.57	0.40
1:B:505:LEU:HD23	1:B:505:LEU:HA	1.94	0.40
1:A:362:ASP:HA	1:A:363:PRO:HD2	1.88	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	302/327 (92%)	275 (91%)	22 (7%)	5 (2%)	7	25
1	B	301/327 (92%)	274 (91%)	25 (8%)	2 (1%)	18	47
All	All	603/654 (92%)	549 (91%)	47 (8%)	7 (1%)	10	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	424	ASP
1	A	454	GLU
1	A	521	MET
1	A	524	MET
1	A	386	GLY
1	B	355	ARG
1	A	338	THR

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	241/267 (90%)	233 (97%)	8 (3%)	33	69
1	B	242/267 (91%)	236 (98%)	6 (2%)	42	76
All	All	483/534 (90%)	469 (97%)	14 (3%)	37	73

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	261	ASP
1	A	360	GLU
1	A	391	VAL
1	A	440	LEU
1	A	454	GLU
1	A	512	LYS
1	A	516	ASN
1	A	559	CYS
1	B	357	LYS

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Mol	Chain	Res	Type
1	B	361	ASP
1	B	391	VAL
1	B	440	LEU
1	B	512	LYS
1	B	531	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	273	HIS
1	A	274	ASN
1	A	407	HIS
1	A	496	ASN
1	A	520	ASN
1	A	525	ASN
1	A	541	HIS
1	B	264	GLN
1	B	267	ASN
1	B	273	HIS
1	B	387	GLN
1	B	407	HIS
1	B	425	ASN
1	B	496	ASN
1	B	516	ASN
1	B	541	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	304/327 (92%)	-0.36	4 (1%) 75 66	5, 22, 38, 42	0
1	B	303/327 (92%)	-0.42	3 (0%) 79 72	4, 19, 36, 44	0
All	All	607/654 (92%)	-0.39	7 (1%) 76 68	4, 20, 37, 44	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	420	GLN	3.5
1	A	424	ASP	3.3
1	B	420	GLN	3.0
1	B	424	ASP	3.0
1	A	526	ALA	3.0
1	A	425	ASN	2.5
1	B	527	GLY	2.3

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.