



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:43 PM UTC

PDB ID : 4S2P / pdb_00004s2p
Title : Crystal structure of unbound OXA-48
Authors : King, D.T.; Strynadka, N.C.J.
Deposited on : 2015-01-21
Resolution : 1.70 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

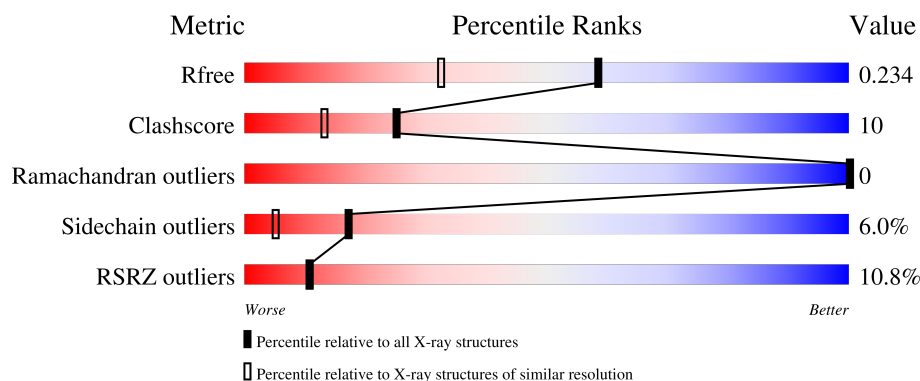
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	265	10% 70% 19% • 9%
1	B	265	9% 70% 19% • 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4311 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 Beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	1	0
			1983	1261	350	365	7			
1	B	241	Total	C	N	O	S	0	2	0
			1986	1263	350	366	7			

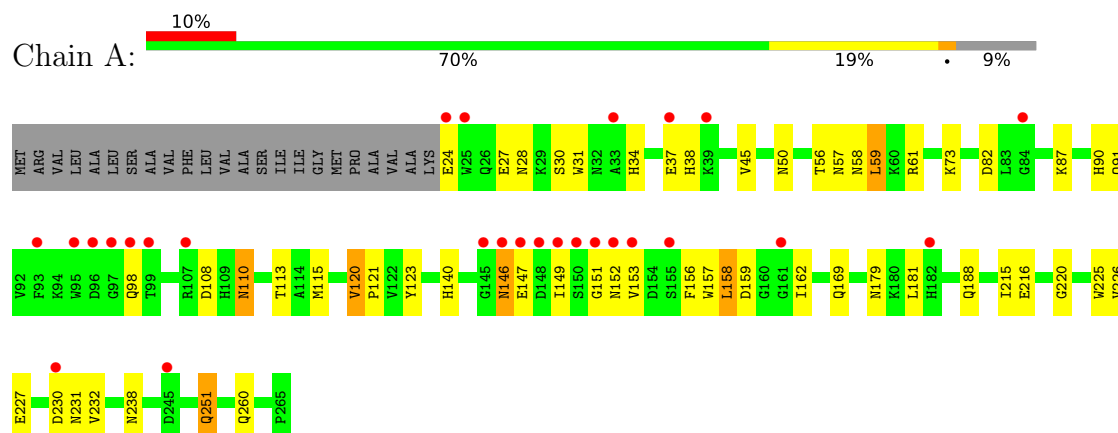
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	191	Total	O	0	0
			191	191		
2	B	151	Total	O	0	0
			151	151		

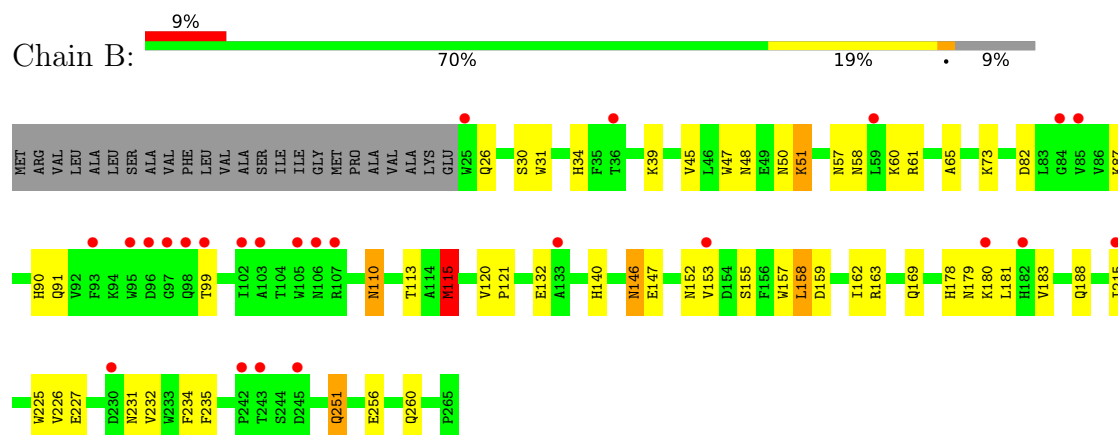
3 Residue-property plots [i](#)

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: Beta-lactamase



• Molecule 1: Beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	43.41 Å 102.87 Å 124.73 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.60 – 1.70 41.60 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (41.60-1.70) 99.6 (41.60-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 1.70 Å)	Xtriage
Refinement program	REFMAC 5.8.0071	Depositor
R, R_{free}	0.192 , 0.226 0.205 , 0.234	Depositor DCC
R_{free} test set	3151 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4311	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.60% 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: KCX

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	0.90	5/2019 (0.2%)	0.85	2/2730 (0.1%)
1	B	0.87	4/2022 (0.2%)	0.89	3/2734 (0.1%)
All	All	0.89	9/4041 (0.2%)	0.87	5/5464 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	226	VAL	C-O	-6.41	1.17	1.24
1	B	227	GLU	C-O	-6.33	1.16	1.24
1	A	226	VAL	C-O	-6.17	1.17	1.24
1	A	45	VAL	C-O	-5.92	1.17	1.24
1	B	115	MET	C-O	-5.88	1.17	1.24
1	A	225	TRP	C-O	-5.66	1.17	1.23
1	B	225	TRP	C-O	-5.62	1.17	1.23
1	A	227	GLU	C-O	-5.43	1.17	1.24
1	A	108	ASP	C-O	-5.12	1.17	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	256	GLU	CA-C-N	-7.33	111.32	120.56
1	B	256	GLU	C-N-CA	-7.33	111.32	120.56
1	A	153	VAL	N-CA-C	6.25	116.92	110.36
1	B	115	MET	CG-SD-CE	-5.73	88.30	100.90
1	A	120	VAL	CB-CA-C	-5.25	108.71	113.70

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	1983	0	1925	40	0
1	B	1986	0	1934	36	0
2	A	191	0	0	8	0
2	B	151	0	0	3	0
All	All	4311	0	3859	76	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 10.

All (76) 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:149:ILE:HG22	1:A:156:PHE:CE2	2.15	0.81
1:B:153:VAL:HG23	2:B:433:HOH:O	1.84	0.76
1:B:87:LYS:H	1:B:91:GLN:NE2	1.85	0.73
1:A:251:GLN:H	1:A:251:GLN:HE21	1.34	0.73
1:B:39:LYS:HG2	2:B:441:HOH:O	1.89	0.71
1:A:87:LYS:H	1:A:91:GLN:NE2	1.92	0.68
1:A:146:ASN:HD21	1:A:162:ILE:HA	1.59	0.68
1:B:251:GLN:HE21	1:B:251:GLN:H	1.42	0.67
1:B:50:ASN:H	1:B:231:ASN:HD21	1.41	0.66
1:A:50:ASN:H	1:A:231:ASN:HD21	1.42	0.66
1:B:39:LYS:NZ	2:B:441:HOH:O	2.28	0.66
1:B:152:ASN:HB2	1:B:155[B]:SER:HB3	1.76	0.66
1:B:159:ASP:HA	1:B:215:ILE:HD12	1.77	0.65
1:A:37:GLU:OE2	1:A:38:HIS:NE2	2.31	0.64
1:A:151:GLY:HA3	2:A:304:HOH:O	1.99	0.63
1:A:149:ILE:O	1:A:149:ILE:HG13	1.99	0.62
1:B:140:HIS:HD2	1:B:147:GLU:OE1	1.82	0.62
1:B:146:ASN:HD21	1:B:162:ILE:HA	1.66	0.60
1:B:110:ASN:HD22	1:B:110:ASN:C	2.09	0.60
1:A:37:GLU:HG3	2:A:426:HOH:O	2.01	0.59
1:A:157:TRP:CD1	1:A:158:LEU:HD13	2.37	0.59
1:B:157:TRP:CD1	1:B:158:LEU:HD13	2.38	0.59
1:A:157:TRP:HA	1:A:162:ILE:CG2	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:343:HOH:O	1:B:90:HIS:HE1	1.85	0.59
1:A:110:ASN:HD22	1:A:110:ASN:C	2.10	0.59
1:B:58:ASN:C	1:B:58:ASN:HD22	2.10	0.58
1:A:58:ASN:HD22	1:A:58:ASN:C	2.12	0.57
1:B:152:ASN:HB2	1:B:155[A]:SER:OG	2.05	0.57
1:A:90:HIS:NE2	2:A:313:HOH:O	2.33	0.56
1:B:31:TRP:HB2	1:B:57:ASN:HB3	1.88	0.56
1:B:48:ASN:ND2	1:B:51:LYS:HG3	2.21	0.56
1:B:152:ASN:HB2	1:B:155[B]:SER:CB	2.38	0.54
1:A:146:ASN:HD22	1:A:146:ASN:C	2.14	0.54
1:A:140:HIS:HD2	1:A:147:GLU:OE2	1.92	0.53
1:B:58:ASN:ND2	1:B:61:ARG:H	2.07	0.53
1:A:28:ASN:HD22	1:A:28:ASN:C	2.18	0.52
2:A:415:HOH:O	1:B:90:HIS:HD2	1.93	0.52
1:B:120:VAL:N	1:B:121:PRO:HD2	2.25	0.51
1:A:27:GLU:HA	1:A:56:THR:O	2.10	0.51
1:A:149:ILE:HG22	1:A:156:PHE:CD2	2.45	0.51
1:A:149:ILE:CG2	1:A:156:PHE:CE2	2.92	0.49
1:A:146:ASN:C	1:A:146:ASN:ND2	2.69	0.49
1:A:251:GLN:H	1:A:251:GLN:NE2	2.05	0.49
1:A:98:GLN:NE2	2:A:449:HOH:O	2.24	0.49
1:A:115:MET:HG2	1:A:123:TYR:OH	2.13	0.48
1:A:179:ASN:HA	1:A:188:GLN:NE2	2.28	0.48
1:A:56:THR:CG2	1:A:59:LEU:HD22	2.44	0.47
1:A:120:VAL:N	1:A:121:PRO:HD2	2.29	0.47
1:A:159:ASP:HA	1:A:215:ILE:HD12	1.95	0.47
1:B:115:MET:HE3	1:B:115:MET:HB2	1.64	0.47
1:B:110:ASN:ND2	1:B:113:THR:H	2.13	0.46
1:B:34:HIS:NE2	1:B:260:GLN:NE2	2.63	0.46
1:A:58:ASN:ND2	1:A:61:ARG:H	2.14	0.46
1:B:87:LYS:H	1:B:91:GLN:HE22	1.59	0.46
1:B:181:LEU:H	1:B:188:GLN:HE22	1.64	0.46
1:A:31:TRP:HB2	1:A:57:ASN:HB3	1.97	0.45
1:A:181:LEU:H	1:A:188:GLN:HE22	1.65	0.45
1:B:146:ASN:ND2	1:B:163:ARG:H	2.15	0.44
1:B:120:VAL:N	1:B:121:PRO:CD	2.81	0.44
1:A:34:HIS:NE2	1:A:260:GLN:NE2	2.66	0.43
2:A:343:HOH:O	1:B:90:HIS:CE1	2.66	0.43
1:B:45:VAL:O	1:B:235:PHE:HA	2.18	0.43
1:A:87:LYS:HD2	1:A:87:LYS:HA	1.66	0.43
1:A:231:ASN:ND2	1:A:232:VAL:H	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:ASN:C	1:B:146:ASN:HD22	2.28	0.42
1:A:28:ASN:HD22	1:A:30:SER:H	1.68	0.42
1:B:179:ASN:HA	1:B:188:GLN:NE2	2.35	0.41
1:A:115:MET:HE3	1:A:115:MET:HB3	1.77	0.41
1:B:178:HIS:HB3	1:B:180:LYS:HE3	2.01	0.41
1:A:220:GLY:O	1:A:238:ASN:HA	2.20	0.41
1:A:24:GLU:HA	2:A:437:HOH:O	2.20	0.41
1:B:65:ALA:HB1	1:B:163:ARG:HB3	2.03	0.41
1:B:157:TRP:HA	1:B:162:ILE:CG2	2.51	0.41
1:B:47:TRP:HB3	1:B:234:PHE:HB2	2.03	0.40
1:A:110:ASN:ND2	1:A:113:THR:H	2.19	0.40
1:A:120:VAL:N	1:A:121:PRO:CD	2.84	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	240/265 (91%)	233 (97%)	7 (3%)	0	100	100
1	B	240/265 (91%)	235 (98%)	5 (2%)	0	100	100
All	All	480/530 (91%)	468 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	209/227 (92%)	199 (95%)	10 (5%)	23	8
1	B	211/227 (93%)	196 (93%)	15 (7%)	13	3
All	All	420/454 (92%)	395 (94%)	25 (6%)	17	5

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	LEU
1	A	82	ASP
1	A	110	ASN
1	A	146	ASN
1	A	152	ASN
1	A	158	LEU
1	A	169	GLN
1	A	216	GLU
1	A	230	ASP
1	A	251	GLN
1	B	26	GLN
1	B	30	SER
1	B	51	LYS
1	B	60	LYS
1	B	82	ASP
1	B	99	THR
1	B	110	ASN
1	B	115	MET
1	B	132	GLU
1	B	146	ASN
1	B	158	LEU
1	B	169	GLN
1	B	183	VAL
1	B	232	VAL
1	B	251	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	58	ASN
1	A	63	ASN

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Mol	Chain	Res	Type
1	A	64	GLN
1	A	76	ASN
1	A	91	GLN
1	A	110	ASN
1	A	140	HIS
1	A	146	ASN
1	A	152	ASN
1	A	188	GLN
1	A	193	GLN
1	A	231	ASN
1	A	251	GLN
1	A	260	GLN
1	B	28	ASN
1	B	52	GLN
1	B	58	ASN
1	B	63	ASN
1	B	76	ASN
1	B	90	HIS
1	B	91	GLN
1	B	98	GLN
1	B	110	ASN
1	B	124	GLN
1	B	140	HIS
1	B	146	ASN
1	B	188	GLN
1	B	193	GLN
1	B	231	ASN
1	B	251	GLN
1	B	260	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
1	KCX	B	73	1	10,11,12	0.75	0	6,12,14	1.28	1 (16%)
1	KCX	A	73	1	10,11,12	0.77	0	6,12,14	1.49	2 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	KCX	B	73	1	-	0/9/10/12	-
1	KCX	A	73	1	-	0/9/10/12	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	KCX	OQ1-CX-NZ	-2.68	120.85	124.92
1	A	73	KCX	CE-NZ-CX	2.14	125.61	121.98
1	B	73	KCX	CE-NZ-CX	2.09	125.53	121.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	A	241/265 (90%)	0.55	27 (11%)	10 10	9, 25, 46, 77	1 (0%)
1	B	240/265 (90%)	0.56	25 (10%)	11 12	9, 24, 47, 61	2 (0%)
All	All	481/530 (90%)	0.56	52 (10%)	11 11	9, 24, 47, 77	3 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	25	TRP	6.8
1	A	149	ILE	6.2
1	A	99	THR	5.5
1	B	99	THR	5.0
1	B	95	TRP	4.6
1	A	95	TRP	4.2
1	A	24	GLU	4.2
1	B	97	GLY	4.0
1	A	98	GLN	3.8
1	A	152	ASN	3.7
1	B	245	ASP	3.5
1	A	97	GLY	3.4
1	A	153	VAL	3.3
1	B	153	VAL	3.2
1	A	151	GLY	3.2
1	A	148	ASP	3.1
1	A	161	GLY	3.0
1	A	37	GLU	2.9
1	B	230	ASP	2.8
1	A	147	GLU	2.8
1	B	107	ARG	2.7
1	B	98	GLN	2.7
1	A	150	SER	2.6
1	B	133	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	243	THR	2.5
1	B	102	ILE	2.5
1	B	103	ALA	2.5
1	A	39	LYS	2.5
1	A	96	ASP	2.5
1	B	84	GLY	2.5
1	B	106	ASN	2.4
1	B	36	THR	2.4
1	B	96	ASP	2.4
1	A	107	ARG	2.4
1	B	105	TRP	2.3
1	A	33	ALA	2.3
1	B	215	ILE	2.2
1	A	146	ASN	2.2
1	A	93	PHE	2.2
1	B	182	HIS	2.2
1	B	59	LEU	2.2
1	A	230	ASP	2.1
1	A	84	GLY	2.1
1	A	155	SER	2.1
1	A	25	TRP	2.1
1	B	180	LYS	2.1
1	A	145	GLY	2.0
1	B	85	VAL	2.0
1	B	242	PRO	2.0
1	B	93	PHE	2.0
1	A	245	ASP	2.0
1	A	182	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	KCX	A	73	12/13	0.94	0.09	15,20,30,38	0
1	KCX	B	73	12/13	0.96	0.07	15,20,25,29	0

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.