



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2017 – 10:27 PM EDT

PDB ID : 3DZC
Title : 2.35 Angstrom resolution structure of WecB (VC0917), a UDP-N-acetylglucosamine 2-epimerase from *Vibrio cholerae*.
Authors : Minasov, G.; Shuvalova, L.; Dubrovskaya, I.; Winsor, J.; Papazisi, L.; Kwon, K.; Hasseman, J.; Peterson, S.N.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : unknown
Resolution : 2.35 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

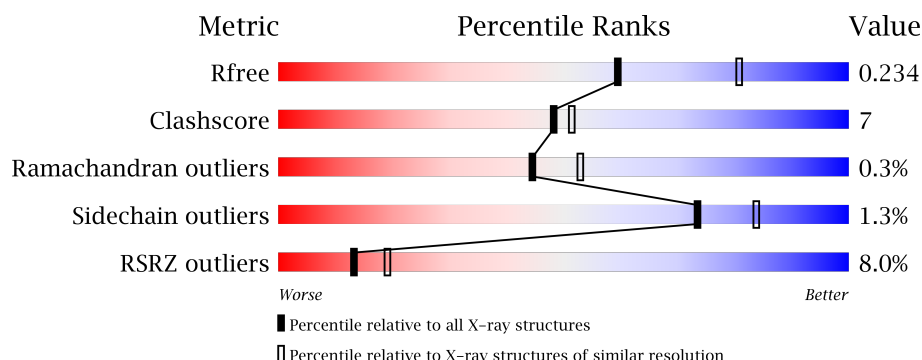
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	100719	1522 (2.38-2.34)
Clashscore	112137	1626 (2.38-2.34)
Ramachandran outliers	110173	1605 (2.38-2.34)
Sidechain outliers	110143	1606 (2.38-2.34)
RSRZ outliers	101464	1528 (2.38-2.34)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	396	<div> <div>9%</div> <div> <div></div> <div>80%</div> <div>14%</div> <div>6%</div> </div> </div>
1	B	396	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>6%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6616 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 UDP-N-acetylglucosamine 2-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	373	Total	C	N	O	S	0	25	0
			3092	1944	539	592	17			
1	B	373	Total	C	N	O	S	0	16	0
			3020	1903	524	576	17			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	EXPRESSION TAG	UNP Q9KTI5
A	-22	HIS	-	EXPRESSION TAG	UNP Q9KTI5
A	-21	HIS	-	EXPRESSION TAG	UNP Q9KTI5
A	-20	HIS	-	EXPRESSION TAG	UNP Q9KTI5
A	-19	HIS	-	EXPRESSION TAG	UNP Q9KTI5
A	-18	HIS	-	EXPRESSION TAG	UNP Q9KTI5
A	-17	HIS	-	EXPRESSION TAG	UNP Q9KTI5
A	-16	SER	-	EXPRESSION TAG	UNP Q9KTI5
A	-15	SER	-	EXPRESSION TAG	UNP Q9KTI5
A	-14	GLY	-	EXPRESSION TAG	UNP Q9KTI5
A	-13	VAL	-	EXPRESSION TAG	UNP Q9KTI5
A	-12	ASP	-	EXPRESSION TAG	UNP Q9KTI5
A	-11	LEU	-	EXPRESSION TAG	UNP Q9KTI5
A	-10	GLY	-	EXPRESSION TAG	UNP Q9KTI5
A	-9	THR	-	EXPRESSION TAG	UNP Q9KTI5
A	-8	GLU	-	EXPRESSION TAG	UNP Q9KTI5
A	-7	ASN	-	EXPRESSION TAG	UNP Q9KTI5
A	-6	LEU	-	EXPRESSION TAG	UNP Q9KTI5
A	-5	TYR	-	EXPRESSION TAG	UNP Q9KTI5
A	-4	PHE	-	EXPRESSION TAG	UNP Q9KTI5
A	-3	GLN	-	EXPRESSION TAG	UNP Q9KTI5
A	-2	SER	-	EXPRESSION TAG	UNP Q9KTI5
A	-1	ASN	-	EXPRESSION TAG	UNP Q9KTI5
A	0	ALA	-	EXPRESSION TAG	UNP Q9KTI5
B	-23	MET	-	EXPRESSION TAG	UNP Q9KTI5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	HIS	-	EXPRESSION TAG	UNP Q9KTI5
B	-21	HIS	-	EXPRESSION TAG	UNP Q9KTI5
B	-20	HIS	-	EXPRESSION TAG	UNP Q9KTI5
B	-19	HIS	-	EXPRESSION TAG	UNP Q9KTI5
B	-18	HIS	-	EXPRESSION TAG	UNP Q9KTI5
B	-17	HIS	-	EXPRESSION TAG	UNP Q9KTI5
B	-16	SER	-	EXPRESSION TAG	UNP Q9KTI5
B	-15	SER	-	EXPRESSION TAG	UNP Q9KTI5
B	-14	GLY	-	EXPRESSION TAG	UNP Q9KTI5
B	-13	VAL	-	EXPRESSION TAG	UNP Q9KTI5
B	-12	ASP	-	EXPRESSION TAG	UNP Q9KTI5
B	-11	LEU	-	EXPRESSION TAG	UNP Q9KTI5
B	-10	GLY	-	EXPRESSION TAG	UNP Q9KTI5
B	-9	THR	-	EXPRESSION TAG	UNP Q9KTI5
B	-8	GLU	-	EXPRESSION TAG	UNP Q9KTI5
B	-7	ASN	-	EXPRESSION TAG	UNP Q9KTI5
B	-6	LEU	-	EXPRESSION TAG	UNP Q9KTI5
B	-5	TYR	-	EXPRESSION TAG	UNP Q9KTI5
B	-4	PHE	-	EXPRESSION TAG	UNP Q9KTI5
B	-3	GLN	-	EXPRESSION TAG	UNP Q9KTI5
B	-2	SER	-	EXPRESSION TAG	UNP Q9KTI5
B	-1	ASN	-	EXPRESSION TAG	UNP Q9KTI5
B	0	ALA	-	EXPRESSION TAG	UNP Q9KTI5

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	1	Total Ca 1 1	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	247	Total 250	O 250	0	8
4	B	241	Total 250	O 250	0	16

- Molecule 1: UDP-N-acetylglucosamine 2-epimerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.39Å 88.13Å 132.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.35 29.90 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-2.35) 99.9 (29.90-2.35)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.40 (at 2.36Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.172 , 0.234 0.173 , 0.234	Depositor DCC
R_{free} test set	2013 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6616	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CL

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	0.62	0/3156	0.76	1/4282 (0.0%)
1	B	0.61	0/3074	0.76	1/4175 (0.0%)
All	All	0.62	0/6230	0.76	2/8457 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	95	GLY	N-CA-C	6.06	128.26	113.10
1	A	283	ARG	NE-CZ-NH1	-5.37	117.61	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3092	0	3092	54	0
1	B	3020	0	3032	32	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	250	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	250	0	0	4	0
All	All	6616	0	6124	85	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 7.

All (85) 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:66:GLN:HE22	1:B:74:LYS:HD2	1.36	0.88
1:A:258:LYS:HA	1:A:261:LYS:HE3	1.63	0.79
1:A:165[A]:GLU:HG3	4:A:595:HOH:O	1.86	0.74
1:B:69:ASN:HD22	1:B:69:ASN:H	1.38	0.72
1:A:369:ILE:O	1:A:372[A]:LYS:HG2	1.90	0.71
1:A:182[A]:ARG:HD3	1:A:301:LEU:HD22	1.75	0.68
1:B:45:ASP:O	1:B:49:GLU:HG3	1.94	0.68
1:A:191:LEU:HD11	1:A:275:LEU:HD21	1.75	0.67
1:A:228:GLN:HG2	4:A:411:HOH:O	1.93	0.67
1:A:41:ARG:HH21	1:A:43[B]:MET:HG2	1.58	0.66
1:B:201:MET:HG3	1:B:267:VAL:HG11	1.77	0.66
1:A:41:ARG:HG2	1:A:42:GLU:H	1.58	0.66
1:B:39:GLN:HA	1:B:62[A]:MET:HG3	1.77	0.65
1:A:69:ASN:HD22	1:A:69:ASN:H	1.45	0.64
1:B:202:LEU:HD21	1:B:243:LEU:HD22	1.80	0.64
1:A:41:ARG:HH21	1:A:43[A]:MET:HG3	1.63	0.62
1:A:251:ASN:ND2	4:A:617:HOH:O	2.32	0.61
1:A:41:ARG:HG2	1:A:42:GLU:N	2.15	0.61
1:B:329:ASN:HD21	1:B:332:GLN:CD	2.05	0.60
1:A:41:ARG:CG	1:A:42:GLU:H	2.16	0.57
1:B:329:ASN:HD22	1:B:329:ASN:H	1.54	0.56
1:A:11:ARG:NH2	1:A:43[A]:MET:HE3	2.21	0.55
1:B:244:TYR:HB3	1:B:268:LEU:HD23	1.89	0.55
1:B:136:ARG:NH2	4:B:531[B]:HOH:O	2.28	0.54
1:A:62:MET:HB3	4:A:486:HOH:O	2.07	0.54
1:B:66:GLN:NE2	1:B:74:LYS:HD2	2.16	0.53
1:B:313[B]:GLU:HG2	1:B:314:ARG:HD2	1.89	0.53
1:A:62:MET:HA	1:A:66[A]:GLN:HE22	1.74	0.53
1:B:219:PHE:CZ	1:B:224:GLU:HB3	2.44	0.53
1:A:80[A]:GLN:NE2	1:B:67:THR:OG1	2.44	0.51
1:B:328:THR:O	1:B:328:THR:HG22	2.09	0.51
1:A:3[B]:LYS:HG3	1:A:88:PRO:HA	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:372[A]:LYS:HG3	4:A:423:HOH:O	2.11	0.50
1:B:11:ARG:HB3	1:B:12:PRO:HD3	1.93	0.50
1:A:313[B]:GLU:HG2	1:A:314:ARG:HD3	1.94	0.50
1:A:244:TYR:HB3	1:A:268:LEU:HD22	1.93	0.50
1:A:312:THR:HB	1:A:325:LEU:HD11	1.94	0.50
1:B:75:ILE:HD13	1:B:99:THR:HA	1.94	0.49
1:A:22[B]:GLN:CA	1:A:23:GLN:N	2.75	0.49
1:B:313[B]:GLU:HG2	1:B:314:ARG:HH11	1.76	0.49
1:A:66[B]:GLN:HG3	4:B:465:HOH:O	2.12	0.48
1:A:295:GLN:HB3	1:A:323:VAL:HG21	1.93	0.48
1:A:372[A]:LYS:CG	4:A:423:HOH:O	2.61	0.48
1:B:372:LYS:HE3	1:B:372:LYS:HB2	1.62	0.48
1:A:63:GLU:H	1:A:66[A]:GLN:NE2	2.12	0.47
1:B:219:PHE:HB2	1:B:252:VAL:HG23	1.96	0.47
1:A:41:ARG:CZ	1:A:41:ARG:H	2.28	0.47
1:B:20:LEU:HD21	1:B:92[B]:LEU:HD21	1.97	0.47
1:A:66[B]:GLN:HE22	1:A:74:LYS:HD2	1.80	0.46
1:B:244:TYR:CE1	1:B:246:VAL:HB	2.50	0.46
1:A:314:ARG:HG3	1:A:357:TYR:OH	2.15	0.46
1:B:154:ARG:HG3	1:B:167:ILE:HG21	1.96	0.46
1:A:121[B]:LEU:HD12	4:A:403:HOH:O	2.16	0.46
1:A:212:THR:O	1:A:289:THR:HA	2.16	0.46
1:B:244:TYR:HB3	1:B:268:LEU:CD2	2.46	0.46
1:A:69:ASN:H	1:A:69:ASN:ND2	2.12	0.45
1:A:216:ARG:NH2	1:A:290:ASP:OD1	2.49	0.45
1:B:38:GLY:N	4:B:691:HOH:O	2.48	0.45
1:A:254[A]:GLU:HB2	1:A:255:PRO:HD3	1.97	0.45
1:A:43[A]:MET:HE1	4:A:441:HOH:O	2.17	0.45
1:A:92:LEU:HD21	1:A:370:LEU:HD11	1.98	0.45
1:A:66[A]:GLN:NE2	4:A:535:HOH:O	2.48	0.45
1:B:22:GLN:HG2	4:B:540:HOH:O	2.16	0.45
1:B:313[B]:GLU:CG	1:B:314:ARG:HH11	2.30	0.45
1:B:69:ASN:ND2	1:B:69:ASN:H	2.10	0.45
1:B:14:ALA:HB2	1:B:44:LEU:HD22	2.00	0.44
1:A:316[B]:GLU:CD	1:A:316[B]:GLU:H	2.22	0.43
1:A:41:ARG:CG	1:A:42:GLU:N	2.78	0.43
1:A:307:VAL:HG23	1:A:323:VAL:HG13	2.01	0.43
1:A:38:GLY:HA2	4:A:609:HOH:O	2.19	0.42
1:A:329:ASN:OD1	1:A:331:GLN:HB3	2.19	0.42
1:B:237:HIS:NE2	1:B:338:SER:HB3	2.34	0.42
1:A:254[B]:GLU:HB3	1:A:255:PRO:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163[B]:ASN:ND2	1:A:165[B]:GLU:OE2	2.52	0.42
1:A:43[B]:MET:HE2	1:A:43[B]:MET:HA	2.01	0.42
1:A:244:TYR:CE1	1:A:246:VAL:HB	2.55	0.42
1:B:4[A]:VAL:HG22	1:B:90:VAL:CG1	2.50	0.41
1:A:11:ARG:NH2	1:A:43[A]:MET:CE	2.83	0.41
1:A:211:VAL:HA	1:A:288:LEU:O	2.21	0.41
1:A:307:VAL:HB	1:A:325:LEU:HD23	2.03	0.41
1:A:17:MET:SD	1:A:94:HIS:HB2	2.60	0.41
1:A:38:GLY:O	1:A:62:MET:HB2	2.21	0.40
1:B:69:ASN:ND2	1:B:69:ASN:N	2.69	0.40
1:A:309:ARG:HD3	1:A:313[B]:GLU:OE2	2.21	0.40
1:A:226:ILE:O	1:A:230:LEU:HG	2.22	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	395/396 (100%)	385 (98%)	10 (2%)	0	100	100
1	B	387/396 (98%)	379 (98%)	6 (2%)	2 (0%)	32	37
All	All	782/792 (99%)	764 (98%)	16 (2%)	2 (0%)	44	53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	41	ARG
1	B	261	LYS

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	342/338 (101%)	339 (99%)	3 (1%)	82	90
1	B	333/338 (98%)	328 (98%)	5 (2%)	70	81
All	All	675/676 (100%)	667 (99%)	8 (1%)	73	86

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	ARG
1	A	69	ASN
1	A	244	TYR
1	B	69	ASN
1	B	176	ASP
1	B	201	MET
1	B	244	TYR
1	B	329	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	46	GLN
1	A	69	ASN
1	A	198	GLN
1	A	214	HIS
1	A	228	GLN
1	A	236	GLN
1	A	251	ASN
1	A	332	GLN
1	B	22	GLN
1	B	39	GLN
1	B	40	HIS
1	B	66	GLN
1	B	69	ASN

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Mol	Chain	Res	Type
1	B	81	GLN
1	B	110	GLN
1	B	161	ASN
1	B	329	ASN
1	B	364	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	373/396 (94%)	0.33	34 (9%) 10 15	20, 40, 72, 83	0
1	B	373/396 (94%)	0.24	26 (6%) 17 24	20, 38, 71, 85	0
All	All	746/792 (94%)	0.29	60 (8%) 13 19	20, 39, 72, 85	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	311	THR	6.0
1	A	189	MET	5.1
1	A	214	HIS	4.7
1	B	41	ARG	4.6
1	B	261	LYS	4.3
1	A	345	GLN	4.3
1	A	191	LEU	4.2
1	A	310	GLU	4.1
1	A	190	ASP	4.0
1	A	261	LYS	4.0
1	B	38	GLY	4.0
1	B	40	HIS	3.7
1	B	190	ASP	3.7
1	B	42	GLU	3.6
1	B	64	PRO	3.6
1	B	62[A]	MET	3.6
1	B	39	GLN	3.6
1	A	215	ARG	3.4
1	A	312	THR	3.4
1	B	189	MET	3.2
1	A	313[A]	GLU	3.1
1	B	0	ALA	3.1
1	A	372[A]	LYS	3.0
1	A	42	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	354[A]	HIS	2.9
1	A	258	LYS	2.9
1	B	187	THR	2.9
1	A	262	GLY	2.9
1	A	220	GLY	2.8
1	A	242	ILE	2.8
1	A	331	GLN	2.7
1	B	65	GLY	2.6
1	B	193	ALA	2.5
1	B	248	LEU	2.5
1	A	309	ARG	2.4
1	B	191	LEU	2.4
1	B	186	HIS	2.4
1	B	311	THR	2.4
1	A	93	VAL	2.4
1	B	253	ARG	2.4
1	A	117	VAL	2.4
1	A	46	GLN	2.3
1	B	63	GLU	2.3
1	A	188	ASP	2.3
1	A	41	ARG	2.3
1	A	209	ILE	2.2
1	B	194	THR	2.2
1	B	252	VAL	2.2
1	A	0	ALA	2.2
1	A	210	LEU	2.2
1	A	49	GLU	2.2
1	B	93	VAL	2.2
1	B	354	HIS	2.1
1	A	344	PRO	2.1
1	B	1	MET	2.1
1	A	352	GLN	2.1
1	A	101	PHE	2.1
1	B	372	LYS	2.0
1	A	286	ILE	2.0
1	A	102	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	CL	A	374	1/1	1.00	0.13	-0.53	25,25,25,25	0
2	CA	A	373	1/1	1.00	0.09	-1.06	29,29,29,29	0
3	CL	B	374	1/1	1.00	0.11	-1.47	27,27,27,27	0
2	CA	B	373	1/1	0.73	0.17	-	35,35,35,35	1

6.5 Other polymers [i](#)

There are no such residues in this entry.