



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:29 AM UTC

PDB ID : 4RGV / pdb_00004rgv
Title : Crystal structure of the Methanocaldococcus jannaschii G1PDH
Authors : Carbone, V.; Ronimus, R.S.; Schofield, L.R.; Sutherland-Smith, A.J.
Deposited on : 2014-09-30
Resolution : 2.45 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

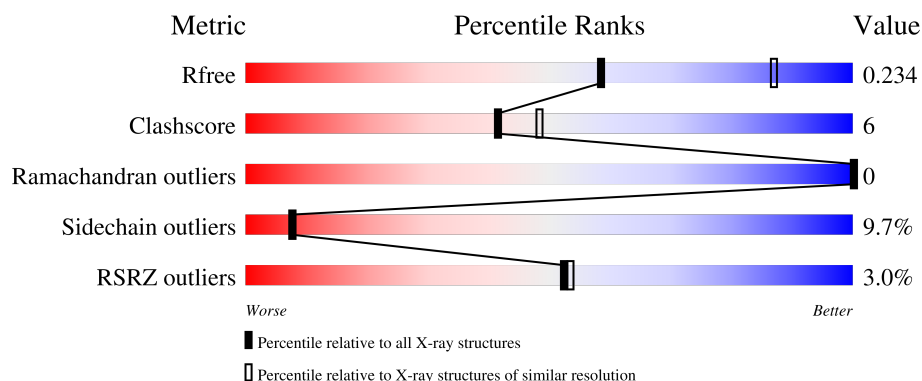
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	368	3% 74% 14% • 11%
1	B	368	2% 68% 19% • 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5188 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 Glycerol-1-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2560	1637	428	486	9			
1	B	333	Total	C	N	O	S	0	0	0
			2604	1665	436	493	10			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-32	MET	-	expression tag	UNP Q58122
A	-31	HIS	-	expression tag	UNP Q58122
A	-30	HIS	-	expression tag	UNP Q58122
A	-29	HIS	-	expression tag	UNP Q58122
A	-28	HIS	-	expression tag	UNP Q58122
A	-27	HIS	-	expression tag	UNP Q58122
A	-26	HIS	-	expression tag	UNP Q58122
A	-25	GLY	-	expression tag	UNP Q58122
A	-24	LYS	-	expression tag	UNP Q58122
A	-23	PRO	-	expression tag	UNP Q58122
A	-22	ILE	-	expression tag	UNP Q58122
A	-21	PRO	-	expression tag	UNP Q58122
A	-20	ASN	-	expression tag	UNP Q58122
A	-19	PRO	-	expression tag	UNP Q58122
A	-18	LEU	-	expression tag	UNP Q58122
A	-17	LEU	-	expression tag	UNP Q58122
A	-16	GLY	-	expression tag	UNP Q58122
A	-15	LEU	-	expression tag	UNP Q58122
A	-14	ASP	-	expression tag	UNP Q58122
A	-13	SER	-	expression tag	UNP Q58122
A	-12	THR	-	expression tag	UNP Q58122
A	-11	GLU	-	expression tag	UNP Q58122
A	-10	ASN	-	expression tag	UNP Q58122
A	-9	LEU	-	expression tag	UNP Q58122
A	-8	TYR	-	expression tag	UNP Q58122

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	PHE	-	expression tag	UNP Q58122
A	-6	GLN	-	expression tag	UNP Q58122
A	-5	GLY	-	expression tag	UNP Q58122
A	-4	ILE	-	expression tag	UNP Q58122
A	-3	ASP	-	expression tag	UNP Q58122
A	-2	PRO	-	expression tag	UNP Q58122
A	-1	PHE	-	expression tag	UNP Q58122
A	0	THR	-	expression tag	UNP Q58122
B	-32	MET	-	expression tag	UNP Q58122
B	-31	HIS	-	expression tag	UNP Q58122
B	-30	HIS	-	expression tag	UNP Q58122
B	-29	HIS	-	expression tag	UNP Q58122
B	-28	HIS	-	expression tag	UNP Q58122
B	-27	HIS	-	expression tag	UNP Q58122
B	-26	HIS	-	expression tag	UNP Q58122
B	-25	GLY	-	expression tag	UNP Q58122
B	-24	LYS	-	expression tag	UNP Q58122
B	-23	PRO	-	expression tag	UNP Q58122
B	-22	ILE	-	expression tag	UNP Q58122
B	-21	PRO	-	expression tag	UNP Q58122
B	-20	ASN	-	expression tag	UNP Q58122
B	-19	PRO	-	expression tag	UNP Q58122
B	-18	LEU	-	expression tag	UNP Q58122
B	-17	LEU	-	expression tag	UNP Q58122
B	-16	GLY	-	expression tag	UNP Q58122
B	-15	LEU	-	expression tag	UNP Q58122
B	-14	ASP	-	expression tag	UNP Q58122
B	-13	SER	-	expression tag	UNP Q58122
B	-12	THR	-	expression tag	UNP Q58122
B	-11	GLU	-	expression tag	UNP Q58122
B	-10	ASN	-	expression tag	UNP Q58122
B	-9	LEU	-	expression tag	UNP Q58122
B	-8	TYR	-	expression tag	UNP Q58122
B	-7	PHE	-	expression tag	UNP Q58122
B	-6	GLN	-	expression tag	UNP Q58122
B	-5	GLY	-	expression tag	UNP Q58122
B	-4	ILE	-	expression tag	UNP Q58122
B	-3	ASP	-	expression tag	UNP Q58122
B	-2	PRO	-	expression tag	UNP Q58122
B	-1	PHE	-	expression tag	UNP Q58122
B	0	THR	-	expression tag	UNP Q58122

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

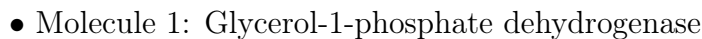
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total	O	0	0
			11	11		
4	B	9	Total	O	0	0
			9	9		

- Molecule 1: Glycerol-1-phosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.77 Å 59.55 Å 119.83 Å 90.00° 90.39° 90.00°	Depositor
Resolution (Å)	46.07 – 2.45 46.07 – 2.46	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.07-2.45) 97.8 (46.07-2.46)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.45 Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.193 , 0.231 0.197 , 0.234	Depositor DCC
R_{free} test set	1293 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	47.4	Xtriage
Anisotropy	0.762	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5188	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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: ZN, MG

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	1.08	1/2600 (0.0%)	1.01	6/3501 (0.2%)
1	B	1.12	0/2645	1.02	5/3562 (0.1%)
All	All	1.10	1/5245 (0.0%)	1.01	11/7063 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	ARG	C-O	-6.06	1.18	1.24

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	ILE	CB-CA-C	-9.04	100.31	111.88
1	A	107	ILE	CB-CA-C	-8.15	101.44	111.88
1	B	219	ARG	CA-C-N	-6.08	113.55	119.87
1	B	219	ARG	C-N-CA	-6.08	113.55	119.87
1	A	98	VAL	CA-C-N	-5.75	113.78	119.99
1	A	98	VAL	C-N-CA	-5.75	113.78	119.99
1	A	319	GLY	CA-C-O	5.37	126.39	121.66
1	A	123	ALA	CA-C-N	-5.22	114.40	119.78
1	A	123	ALA	C-N-CA	-5.22	114.40	119.78
1	B	285	ALA	CA-C-N	-5.12	114.51	119.78
1	B	285	ALA	C-N-CA	-5.12	114.51	119.78

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	2560	0	2647	24	0
1	B	2604	0	2701	44	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	11	0	0	0	0
4	B	9	0	0	0	0
All	All	5188	0	5348	67	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 6.

All (67) 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:257:MET:HG2	1:B:332:THR:HG21	1.47	0.96
1:B:328:LEU:O	1:B:332:THR:HB	1.78	0.82
1:B:332:THR:CG2	1:B:334:VAL:HG23	2.21	0.71
1:A:227:LEU:HD21	1:A:312:ARG:HE	1.56	0.71
1:B:2:ILE:C	1:B:2:ILE:HD12	2.19	0.67
1:B:265:GLU:OE2	1:B:317:ARG:NH1	2.26	0.67
1:B:30:ASN:ND2	1:B:68:ALA:O	2.29	0.66
1:A:288:THR:OG1	1:A:291:GLU:HG3	1.99	0.63
1:A:266:ASN:OD1	1:A:268:LYS:HB3	1.98	0.62
1:A:10:ILE:HD11	1:A:22:ILE:HD11	1.82	0.61
1:B:332:THR:HG23	1:B:334:VAL:HG23	1.83	0.61
1:B:321:SER:OG	1:B:324:GLU:HG3	2.01	0.60
1:A:265:GLU:OE2	1:A:317:ARG:NH1	2.26	0.58
1:B:313:TRP:CZ2	1:B:317:ARG:HG2	2.38	0.58
1:B:10:ILE:HD11	1:B:22:ILE:HD11	1.86	0.57
1:A:313:TRP:CZ2	1:A:317:ARG:HG2	2.41	0.55
1:B:165:GLU:OE2	1:B:317:ARG:NH2	2.34	0.54
1:A:159:TRP:HA	1:A:219:ARG:HH11	1.74	0.53
1:B:300:ILE:O	1:B:304:THR:HG23	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:TRP:CD1	1:A:219:ARG:HD2	2.43	0.53
1:A:140:ARG:HD3	1:A:246:LEU:HD11	1.91	0.53
1:A:326:ARG:O	1:A:330:GLU:HG2	2.10	0.52
1:B:257:MET:CG	1:B:332:THR:HG21	2.32	0.52
1:B:111:ILE:CD1	1:B:120:MET:HE2	2.39	0.52
1:B:233:ASP:HA	1:B:236:LYS:HD2	1.91	0.52
1:B:321:SER:CB	1:B:324:GLU:HG3	2.40	0.52
1:A:312:ARG:O	1:A:312:ARG:CG	2.58	0.51
1:B:265:GLU:CD	1:B:317:ARG:HH12	2.18	0.50
1:B:288:THR:HG22	1:B:333:GLY:O	2.11	0.50
1:B:323:GLU:O	1:B:327:LYS:HB2	2.12	0.49
1:B:116:GLN:HA	1:B:117:PRO:C	2.37	0.48
1:B:85:LYS:NZ	1:B:109:SER:O	2.48	0.47
1:B:147:GLY:HA3	1:B:248:GLY:O	2.15	0.47
1:B:288:THR:OG1	1:B:291:GLU:HG3	2.15	0.46
1:A:120:MET:HE3	1:A:120:MET:HB3	1.86	0.46
1:A:288:THR:HG22	1:A:333:GLY:O	2.16	0.46
1:B:48:ASP:OD2	1:B:69:TYR:OH	2.30	0.45
1:B:99:PRO:HG3	1:B:107:ILE:HG23	1.97	0.45
1:A:5:THR:O	1:B:7:ARG:HD3	2.17	0.45
1:A:313:TRP:HZ2	1:A:317:ARG:HG2	1.81	0.45
1:A:266:ASN:C	1:A:268:LYS:H	2.25	0.45
1:A:2:ILE:O	1:A:2:ILE:HG23	2.17	0.44
1:B:116:GLN:HG3	1:B:117:PRO:HA	1.98	0.44
1:B:317:ARG:HB3	1:B:318:ASP:H	1.64	0.44
1:B:332:THR:HG22	1:B:334:VAL:HG23	1.98	0.43
1:B:111:ILE:HD11	1:B:120:MET:HE2	2.00	0.43
1:A:219:ARG:N	1:A:220:PRO:HD2	2.34	0.43
1:A:31:PRO:HB3	1:A:71:CYS:SG	2.58	0.43
1:B:56:LEU:HA	1:B:59:LEU:HD21	2.01	0.42
1:B:121:VAL:HG23	1:B:122:ASP:N	2.34	0.42
1:B:2:ILE:HD12	1:B:3:ILE:N	2.34	0.42
1:B:120:MET:HE3	1:B:120:MET:HB2	1.72	0.42
1:B:210:ILE:HG22	1:B:214:ILE:HD12	2.02	0.42
1:B:151:SER:HB2	1:B:225:GLU:HG3	2.02	0.42
1:A:266:ASN:C	1:A:268:LYS:N	2.77	0.42
1:B:114:ILE:CG2	1:B:115:ARG:N	2.84	0.41
1:B:107:ILE:HG21	1:B:107:ILE:HD13	1.87	0.41
1:B:193:ASP:O	1:B:194:LEU:HB2	2.20	0.41
1:A:296:GLU:OE2	1:A:326:ARG:NH2	2.41	0.41
1:B:326:ARG:O	1:B:330:GLU:HG2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:LYS:O	1:A:163:TYR:HB3	2.21	0.40
1:A:206:VAL:O	1:A:210:ILE:HD13	2.22	0.40
1:A:232:LEU:HA	1:A:232:LEU:HD23	1.83	0.40
1:B:259:SER:OG	1:B:272:LEU:HD23	2.20	0.40
1:B:283:VAL:O	1:B:284:ASP:CB	2.70	0.40
1:B:5:THR:HB	1:B:6:PRO:HD2	2.04	0.40
1:B:115:ARG:O	1:B:117:PRO:O	2.39	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	322/368 (88%)	311 (97%)	11 (3%)	0	100	100
1	B	329/368 (89%)	318 (97%)	11 (3%)	0	100	100
All	All	651/736 (88%)	629 (97%)	22 (3%)	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	281/318 (88%)	256 (91%)	25 (9%)	9	10

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	286/318 (90%)	256 (90%)	30 (10%)	6	6
All	All	567/636 (89%)	512 (90%)	55 (10%)	8	8

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	37	LYS
1	A	39	THR
1	A	45	PHE
1	A	53	ASP
1	A	54	GLU
1	A	58	ASN
1	A	63	LEU
1	A	65	LYS
1	A	81	ILE
1	A	82	ASP
1	A	90	LYS
1	A	96	ILE
1	A	113	SER
1	A	120	MET
1	A	139	ARG
1	A	164	LYS
1	A	199	ASN
1	A	240	ASN
1	A	268	LYS
1	A	270	SER
1	A	273	HIS
1	A	284	ASP
1	A	290	LYS
1	A	312	ARG
1	B	1	MET
1	B	2	ILE
1	B	7	ARG
1	B	16	ILE
1	B	27	ASN
1	B	29	LYS
1	B	30	ASN
1	B	63	LEU
1	B	65	LYS
1	B	81	ILE
1	B	82	ASP

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Mol	Chain	Res	Type
1	B	90	LYS
1	B	115	ARG
1	B	118	SER
1	B	120	MET
1	B	160	LYS
1	B	166	LYS
1	B	189	VAL
1	B	240	ASN
1	B	245	SER
1	B	268	LYS
1	B	270	SER
1	B	275	LYS
1	B	277	LYS
1	B	278	MET
1	B	284	ASP
1	B	321	SER
1	B	324	GLU
1	B	326	ARG
1	B	332	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	A	116	GLN
1	A	250	GLN
1	A	307	HIS
1	B	273	HIS
1	B	307	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 [i](#)

There are no oligosaccharides 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	328/368 (89%)	0.07	12 (3%) 45 45	36, 56, 83, 119	46 (14%)
1	B	333/368 (90%)	0.08	8 (2%) 59 62	37, 56, 84, 103	45 (13%)
All	All	661/736 (89%)	0.07	20 (3%) 52 53	36, 56, 84, 119	91 (13%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	63	LEU	4.6
1	B	45	PHE	4.2
1	B	63	LEU	3.7
1	A	312	ARG	3.3
1	B	40	LYS	2.8
1	A	37	LYS	2.7
1	A	267	LYS	2.6
1	A	2	ILE	2.6
1	A	217	SER	2.5
1	A	55	ILE	2.4
1	B	2	ILE	2.4
1	B	1	MET	2.4
1	B	311	ASN	2.4
1	B	116	GLN	2.3
1	A	292	LEU	2.2
1	A	117	PRO	2.2
1	A	116	GLN	2.2
1	A	310	ARG	2.1
1	B	41	LYS	2.1
1	A	64	LYS	2.1

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](#)

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($\text{\AA}^2$)	Q<0.9
3	ZN	A	402	1/1	0.92	0.10	59,59,59,59	1
2	MG	A	401	1/1	0.93	0.06	25,25,25,25	0
3	ZN	B	402	1/1	0.93	0.10	58,58,58,58	1
2	MG	B	401	1/1	0.95	0.10	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.