



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 11:57 am BST

PDB ID : 6PWJ
Title : Vibrio cholerae LapD S helix-GGDEF-EAL (apo)
Authors : Giglio, K.M.; Cooley, R.B.; Sondermann, H.
Deposited on : 2019-07-23
Resolution : 2.70 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

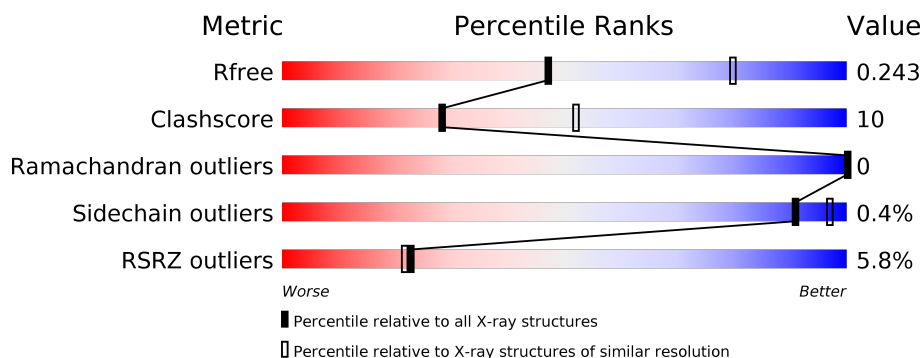
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 on 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	418	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>23%</div> <div></div> </div> </div>
1	B	418	<div> <div>7%</div> <div> <div></div> <div>73%</div> <div>24%</div> <div></div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6523 atoms, of which 14 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 GGDEF and EAL domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	408	Total	C	N	O	S	0	0	0
			3259	2077	554	621	7			
1	B	406	Total	C	N	O	S	0	0	0
			3227	2054	549	617	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	637	ALA	-	expression tag	UNP A0A0H6T0A6
A	638	ALA	-	expression tag	UNP A0A0H6T0A6
B	637	ALA	-	expression tag	UNP A0A0H6T0A6
B	638	ALA	-	expression tag	UNP A0A0H6T0A6

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	0	0
			24	6	14	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	5	Total	O	0	0
			5	5		

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.83Å 129.83Å 114.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	12.69 – 2.70 19.53 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.1 (12.69-2.70) 90.4 (19.53-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.88 (at 2.71Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.199 , 0.243 0.199 , 0.243	Depositor DCC
R_{free} test set	1974 reflections (6.79%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 58.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6523	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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: MG, PGE

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	0.24	0/3325	0.39	0/4496
1	B	0.24	0/3293	0.39	0/4457
All	All	0.24	0/6618	0.39	0/8953

There are no bond length outliers.

There are no bond angle outliers.

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	3259	0	3183	62	0
1	B	3227	0	3118	66	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	B	10	14	14	0	0
4	A	6	0	0	0	0
4	B	5	0	0	0	0
All	All	6509	14	6315	126	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 (126) 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:226:GLN:HG3	1:B:599:THR:HG22	1.38	1.03
1:A:636:LYS:NZ	1:B:585:ASP:OD2	2.00	0.94
1:A:632:PHE:CE2	1:A:636:LYS:HD3	2.16	0.80
1:B:270:LEU:HD11	1:B:355:LEU:HD11	1.64	0.80
1:A:524:GLU:HG2	1:A:554:ASP:HB2	1.66	0.78
1:B:524:GLU:HG2	1:B:554:ASP:HB2	1.69	0.74
1:B:410:LEU:HD21	1:B:443:PHE:HZ	1.53	0.74
1:B:266:GLY:HA3	1:B:329:LEU:HD11	1.69	0.74
1:B:313:SER:HB3	1:B:316:GLU:HB2	1.71	0.73
1:A:463:ASN:OD1	1:A:500:GLN:NE2	2.21	0.73
1:A:381:LYS:HD2	1:A:393:ILE:HD11	1.70	0.73
1:B:277:GLU:OE2	1:B:354:ASN:ND2	2.23	0.71
1:A:608:ILE:HG12	1:A:628:VAL:HB	1.71	0.71
1:B:418:ASP:OD1	1:B:475:ARG:NH1	2.23	0.70
1:A:498:ILE:HD11	1:A:525:ILE:HD12	1.73	0.70
1:B:413:GLU:OE1	1:B:445:LYS:NZ	2.24	0.70
1:A:313:SER:HB3	1:A:316:GLU:HB2	1.75	0.68
1:B:567:ASN:O	1:B:570:ARG:NH1	2.24	0.68
1:A:226:GLN:NE2	1:A:229:GLU:OE1	2.27	0.68
1:A:632:PHE:O	1:A:636:LYS:HG3	1.93	0.68
1:B:584:LEU:HD12	1:B:620:PHE:HE2	1.58	0.67
1:B:595:SER:O	1:B:599:THR:HG23	1.96	0.66
1:B:608:ILE:HG12	1:B:628:VAL:HB	1.77	0.65
1:B:513:LYS:HG3	1:B:514:HIS:CD2	2.32	0.65
1:B:401:ILE:O	1:B:401:ILE:HD12	1.97	0.64
1:A:462:LEU:O	1:A:463:ASN:HB3	1.97	0.64
1:B:223:PHE:CE2	1:B:567:ASN:HB2	2.32	0.64
1:B:304:ILE:CB	1:B:307:ILE:HB	2.28	0.63
1:A:230:ALA:HB1	1:A:596:ILE:HD12	1.81	0.63
1:B:402:LEU:HB2	1:B:407:TRP:CE2	2.34	0.62
1:A:567:ASN:OD1	1:A:603:LEU:HD11	2.00	0.62
1:A:495:GLN:HG3	1:A:529:CYS:SG	2.40	0.62
1:A:632:PHE:CZ	1:A:636:LYS:HD3	2.35	0.62
1:B:271:GLN:HG2	1:B:316:GLU:HG2	1.81	0.61
1:B:328:GLU:O	1:B:331:ILE:HG12	2.01	0.60
1:A:437:ARG:NH2	1:A:487:ASP:OD2	2.36	0.58
1:A:373:SER:O	1:A:377:ASN:ND2	2.36	0.58
1:A:504:ILE:O	1:A:507:ILE:HG22	2.04	0.57
1:A:558:ARG:HD3	1:A:577:ASP:OD1	2.05	0.57
1:B:407:TRP:CE3	1:B:455:PHE:HB3	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:585:ASP:OD2	1:B:636:LYS:NZ	2.38	0.56
1:B:544:VAL:HG13	1:B:549:ALA:HB3	1.85	0.56
1:B:498:ILE:HD11	1:B:525:ILE:HD12	1.87	0.56
1:B:302:ILE:HG22	1:B:303:THR:H	1.70	0.55
1:B:504:ILE:O	1:B:507:ILE:HG22	2.06	0.55
1:A:246:GLY:HA3	1:A:250:TYR:CD1	2.43	0.54
1:A:418:ASP:OD1	1:A:475:ARG:NH1	2.40	0.54
1:B:235:GLU:HG3	1:B:239:LEU:HD12	1.90	0.52
1:A:245:LEU:HD13	1:A:309:ILE:HG22	1.91	0.52
1:A:415:ILE:HG23	1:A:471:TYR:CD1	2.45	0.52
1:A:563:LEU:HD11	1:A:596:ILE:HD11	1.92	0.51
1:B:256:SER:OG	1:B:368:THR:HG21	2.11	0.51
1:A:404:LYS:HE3	1:A:461:GLN:HE22	1.76	0.51
1:A:244:HIS:HB2	1:A:292:ARG:NH1	2.25	0.51
1:B:410:LEU:CD2	1:B:443:PHE:HZ	2.22	0.51
1:B:460:GLU:OE2	1:B:496:GLY:N	2.37	0.50
1:A:325:ASP:OD2	1:A:327:THR:HG22	2.12	0.50
1:B:270:LEU:CD1	1:B:355:LEU:HD11	2.40	0.49
1:A:255:LEU:HD11	1:A:371:LEU:CD2	2.42	0.49
1:A:505:ARG:O	1:A:509:GLN:HG2	2.12	0.49
1:B:490:ALA:HA	1:B:522:HIS:O	2.12	0.49
1:B:437:ARG:NH1	1:B:487:ASP:OD2	2.45	0.49
1:A:563:LEU:HD11	1:A:596:ILE:CD1	2.42	0.49
1:B:495:GLN:HG3	1:B:529:CYS:SG	2.53	0.48
1:A:333:ALA:O	1:A:337:ILE:HG12	2.13	0.48
1:A:403:GLY:H	1:A:406:GLN:NE2	2.11	0.48
1:B:269:ILE:HG13	1:B:375:LEU:HD13	1.94	0.48
1:A:274:PHE:HD1	1:A:354:ASN:O	1.96	0.48
1:B:473:ILE:O	1:B:477:ILE:HG13	2.12	0.48
1:A:251:TYR:HB2	1:A:310:ALA:CB	2.44	0.48
1:B:246:GLY:HA3	1:B:250:TYR:CD2	2.49	0.48
1:B:452:ALA:HA	1:B:455:PHE:CZ	2.49	0.48
1:B:506:TRP:CE2	1:B:510:THR:HG21	2.49	0.48
1:A:327:THR:O	1:A:331:ILE:HG12	2.14	0.47
1:B:267:VAL:HG21	1:B:371:LEU:HD21	1.96	0.47
1:A:354:ASN:OD1	1:A:355:LEU:N	2.47	0.47
1:B:302:ILE:HG22	1:B:303:THR:N	2.29	0.47
1:A:340:VAL:HG11	1:A:355:LEU:HD23	1.97	0.47
1:A:395:SER:HA	1:A:396:ASP:HA	1.52	0.46
1:A:527:GLU:HG3	1:A:565:TYR:OH	2.15	0.46
1:A:308:SER:O	1:A:319:ILE:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:396:ASP:O	1:B:399:LYS:N	2.48	0.45
1:B:327:THR:O	1:B:331:ILE:HG23	2.15	0.45
1:A:276:LYS:HE2	1:A:315:TYR:CZ	2.51	0.45
1:A:228:LYS:O	1:A:232:GLN:HG3	2.16	0.45
1:A:293:GLU:OE1	1:A:343:ILE:HG23	2.16	0.45
1:B:464:ALA:O	1:B:467:ILE:HG12	2.16	0.45
1:A:580:PHE:O	1:A:593:LEU:HD13	2.17	0.44
1:A:403:GLY:H	1:A:406:GLN:HE21	1.66	0.44
1:A:321:MET:HB3	1:A:324:MET:HG3	2.00	0.44
1:A:552:GLY:HA3	1:A:573:TYR:CZ	2.53	0.44
1:A:234:ARG:CZ	1:A:560:PHE:HA	2.48	0.44
1:B:527:GLU:HG3	1:B:565:TYR:OH	2.16	0.44
1:A:410:LEU:HD21	1:A:450:TYR:CE2	2.53	0.44
1:B:530:PHE:CE1	1:B:553:VAL:HG22	2.53	0.43
1:B:533:GLU:N	1:B:534:PRO:HD3	2.33	0.43
1:B:471:TYR:OH	1:B:475:ARG:NH1	2.52	0.43
1:A:452:ALA:HA	1:A:455:PHE:CZ	2.54	0.43
1:B:410:LEU:HD21	1:B:443:PHE:CZ	2.43	0.43
1:A:360:VAL:HG23	1:A:375:LEU:HD23	2.01	0.43
1:A:456:LEU:HA	1:A:456:LEU:HD22	1.90	0.43
1:B:452:ALA:HA	1:B:455:PHE:CE2	2.54	0.42
1:A:509:GLN:O	1:A:513:LYS:HG3	2.19	0.42
1:B:307:ILE:HG13	1:B:320:ILE:O	2.19	0.42
1:B:323:ASN:HB3	1:B:457:PHE:CE1	2.55	0.42
1:A:343:ILE:O	1:A:344:ASN:C	2.58	0.42
1:A:234:ARG:HG2	1:A:560:PHE:CE2	2.55	0.42
1:A:382:ALA:HB2	1:A:391:GLY:H	1.85	0.42
1:B:251:TYR:HB2	1:B:310:ALA:CB	2.49	0.42
1:B:431:TRP:HE1	1:B:433:LYS:CG	2.33	0.42
1:B:266:GLY:HA3	1:B:329:LEU:CD1	2.43	0.41
1:A:438:GLU:HA	1:A:490:ALA:HB3	2.01	0.41
1:A:613:GLU:HA	1:A:632:PHE:HB3	2.01	0.41
1:B:400:ILE:O	1:B:400:ILE:HG13	2.20	0.41
1:B:407:TRP:O	1:B:411:VAL:HG23	2.19	0.41
1:A:452:ALA:HA	1:A:455:PHE:CE2	2.55	0.41
1:B:226:GLN:HG2	1:B:602:ASN:ND2	2.36	0.41
1:A:408:LYS:O	1:A:412:GLU:HG3	2.21	0.41
1:B:506:TRP:CZ2	1:B:510:THR:HG21	2.56	0.41
1:B:525:ILE:HD11	1:B:540:PHE:HE2	1.85	0.40
1:B:330:LYS:HG2	1:B:390:TYR:OH	2.21	0.40
1:B:478:GLN:O	1:B:482:LYS:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:525:ILE:HD11	1:A:540:PHE:CE2	2.56	0.40
1:B:529:CYS:HB3	1:B:537:THR:OG1	2.21	0.40
1:B:552:GLY:HA3	1:B:573:TYR:CZ	2.56	0.40
1:B:553:VAL:HB	1:B:574:VAL:HG12	2.02	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	402/418 (96%)	393 (98%)	9 (2%)	0	100	100
1	B	400/418 (96%)	392 (98%)	8 (2%)	0	100	100
All	All	802/836 (96%)	785 (98%)	17 (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	349/360 (97%)	348 (100%)	1 (0%)	92	98
1	B	343/360 (95%)	341 (99%)	2 (1%)	86	95
All	All	692/720 (96%)	689 (100%)	3 (0%)	91	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	456	LEU
1	B	410	LEU
1	B	509	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	377	ASN
1	A	406	GLN
1	A	461	GLN
1	B	231	GLN
1	B	354	ASN
1	B	389	ASN
1	B	617	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 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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
3	PGE	B	702	-	9,9,9	0.30	0	8,8,8	0.34	0

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
3	PGE	B	702	-	-	4/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	702	PGE	O3-C5-C6-O4
3	B	702	PGE	O2-C3-C4-O3
3	B	702	PGE	C4-C3-O2-C2
3	B	702	PGE	O1-C1-C2-O2

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	408/418 (97%)	0.11	19 (4%) 31 30	26, 51, 88, 130	0
1	B	406/418 (97%)	0.22	28 (6%) 16 15	28, 56, 99, 157	0
All	All	814/836 (97%)	0.17	47 (5%) 23 22	26, 53, 95, 157	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	446	ASP	5.9
1	B	396	ASP	5.5
1	B	306	ASP	5.0
1	B	509	GLN	4.8
1	B	384	SER	4.3
1	A	446	ASP	4.3
1	A	484	GLU	4.0
1	A	387	GLU	3.7
1	A	400	ILE	3.7
1	A	416	HIS	3.6
1	A	384	SER	3.5
1	B	281	GLU	3.4
1	B	380	ALA	3.3
1	A	431	TRP	3.3
1	B	221	LYS	3.2
1	B	405	GLN	3.1
1	B	433	LYS	2.9
1	B	419	TRP	2.9
1	A	417	ASN	2.9
1	B	268	ALA	2.9
1	A	418	ASP	2.8
1	B	418	ASP	2.8
1	A	388	LEU	2.8
1	B	344	ASN	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	341	ASP	2.7
1	A	405	GLN	2.7
1	A	300	ASN	2.7
1	A	303	THR	2.6
1	A	587	GLU	2.6
1	A	596	ILE	2.5
1	A	396	ASP	2.5
1	B	376	ASP	2.5
1	A	540	PHE	2.4
1	A	228	LYS	2.3
1	A	353	ALA	2.3
1	B	403	GLY	2.2
1	B	521	LEU	2.2
1	B	546	ASN	2.2
1	B	417	ASN	2.2
1	B	402	LEU	2.1
1	B	270	LEU	2.1
1	B	431	TRP	2.1
1	B	596	ILE	2.1
1	B	587	GLU	2.1
1	B	353	ALA	2.1
1	B	228	LYS	2.0
1	B	297	ARG	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 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. 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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MG	A	1000	1/1	0.96	0.06	53,53,53,53	0
3	PGE	B	702	10/10	0.96	0.09	41,49,58,59	24
2	MG	B	701	1/1	0.97	0.09	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.