



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:43 AM UTC

PDB ID : 1LVH / pdb_00001lvh
Title : The Structure of Phosphorylated beta-phosphoglucomutase from *Lactococcus lactis* to 2.3 angstrom resolution
Authors : Lahiri, S.D.; Zhang, G.; Dunaway-Mariano, D.; Allen, K.N.
Deposited on : 2002-05-28
Resolution : 2.30 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

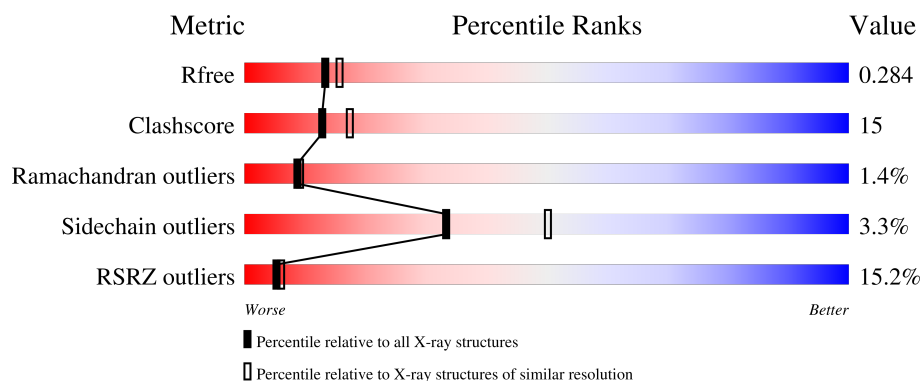
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	221	5% 79% 19% .
1	B	221	25% 65% 30% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3685 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-phosphoglucomutase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	P	S	0	0	0
			1712	1089	285	334	1	3			
1	B	221	Total	C	N	O	P	S	0	0	0
			1712	1089	285	334	1	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	PHD	ASP	modified residue	UNP P71447
B	8	PHD	ASP	modified residue	UNP P71447

- 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		

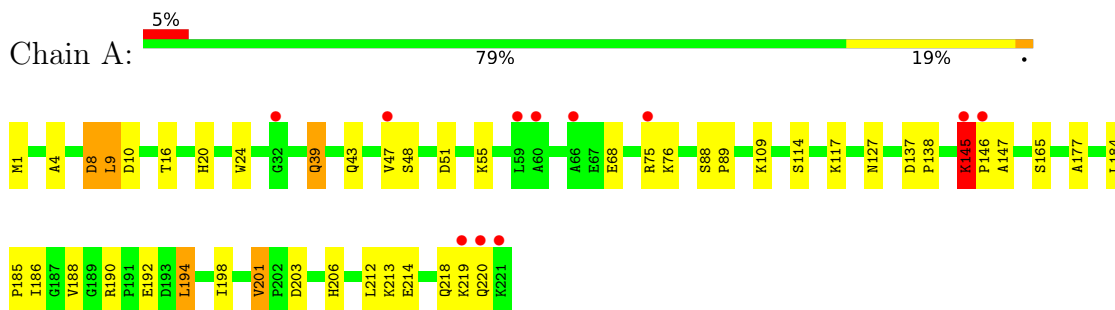
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	167	Total	O	0	0
			167	167		
3	B	92	Total	O	0	0
			92	92		

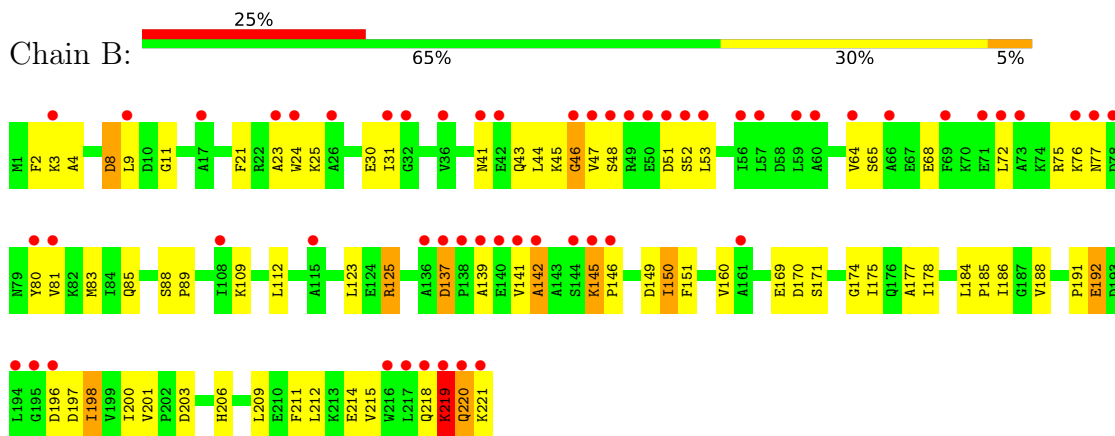
3 Residue-property plots [i](#)

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-phosphoglucosyltransferase



- Molecule 1: beta-phosphoglucosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.67Å 92.78Å 111.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.89 – 2.30 42.89 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.8 (42.89-2.30) 94.8 (42.89-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.82 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.243 , 0.286 0.244 , 0.284	Depositor DCC
R_{free} test set	2390 reflections (9.90%)	wwPDB-VP
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.757	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3685	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.84 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.4461e-03.*

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, PHD

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	0.51	1/1729 (0.1%)	0.98	10/2335 (0.4%)
1	B	0.41	0/1729	0.91	6/2335 (0.3%)
All	All	0.46	1/3458 (0.0%)	0.95	16/4670 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	9	LEU	N-CA	9.93	1.65	1.46

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	LEU	N-CA-CB	10.69	128.68	110.50
1	A	201	VAL	CA-C-N	7.17	127.17	119.28
1	A	201	VAL	C-N-CA	7.17	127.17	119.28
1	A	16	THR	N-CA-C	-6.50	105.92	114.31
1	B	209	LEU	N-CA-C	-5.90	104.85	111.28
1	B	46	GLY	N-CA-C	-5.79	106.51	114.67
1	B	31	ILE	N-CA-C	-5.54	108.45	113.71
1	A	165	SER	N-CA-C	5.40	117.62	109.41
1	A	145	LYS	N-CA-C	5.26	121.44	109.81
1	A	117	LYS	N-CA-C	-5.21	106.88	113.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	142	ALA	N-CA-C	5.16	116.59	110.97
1	A	194	LEU	N-CA-C	5.15	119.64	112.90
1	B	150	ILE	N-CA-C	5.06	115.78	110.62
1	A	137	ASP	CA-C-N	5.04	125.06	119.32
1	A	137	ASP	C-N-CA	5.04	125.06	119.32
1	B	85	GLN	N-CA-C	-5.03	106.25	112.38

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	9	LEU	CA

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	1712	0	1723	33	1
1	B	1712	0	1722	71	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	167	0	0	4	0
3	B	92	0	0	4	0
All	All	3685	0	3445	104	1

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 15.

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:8:PHD:C	1:B:9:LEU:N	1.74	1.46
1:A:184:LEU:HD22	1:A:219:LYS:HD3	1.46	0.97
1:B:169:GLU:HG3	1:B:171:SER:H	1.45	0.81
1:B:191:PRO:HG3	1:B:200:ILE:HD11	1.62	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:GLU:HG3	1:B:218:GLN:HE21	1.48	0.77
1:A:145:LYS:O	1:A:177:ALA:HB2	1.87	0.74
1:A:39:GLN:N	3:A:867:HOH:O	2.21	0.74
1:B:45:LYS:HD3	1:B:46:GLY:N	2.03	0.74
1:B:8:PHD:C	1:B:9:LEU:CA	2.66	0.72
1:B:145:LYS:O	1:B:177:ALA:HB2	1.92	0.69
1:B:203:ASP:OD2	1:B:206:HIS:NE2	2.27	0.68
1:A:1:MET:HE3	1:A:213:LYS:HE2	1.75	0.67
1:B:23:ALA:HB2	1:B:83:MET:HE1	1.75	0.67
1:B:219:LYS:O	1:B:220:GLN:HG2	1.94	0.67
1:B:211:PHE:O	1:B:215:VAL:HG23	1.95	0.66
1:B:8:PHD:O	1:B:9:LEU:N	2.23	0.65
1:B:125:ARG:HA	1:B:125:ARG:HE	1.63	0.63
1:B:191:PRO:HD2	1:B:192:GLU:OE2	1.98	0.63
1:B:169:GLU:HG3	1:B:170:ASP:N	2.13	0.62
1:B:198:ILE:H	1:B:198:ILE:HD13	1.64	0.62
1:A:145:LYS:O	1:A:146:PRO:C	2.42	0.62
1:A:184:LEU:HD12	1:A:185:PRO:HD2	1.82	0.61
1:B:197:ASP:OD1	1:B:198:ILE:HG23	2.02	0.59
1:B:48:SER:O	1:B:52:SER:HB2	2.02	0.59
1:A:190:ARG:HE	1:A:192:GLU:HB2	1.67	0.59
1:B:43:GLN:HB2	3:B:864:HOH:O	2.03	0.59
1:B:77:ASN:O	1:B:81:VAL:HG23	2.03	0.58
1:A:203:ASP:OD2	1:A:206:HIS:HE1	1.86	0.58
1:B:76:LYS:NZ	3:B:814:HOH:O	2.36	0.58
1:B:218:GLN:O	1:B:219:LYS:C	2.48	0.57
1:B:21:PHE:CZ	1:B:25:LYS:HD2	2.40	0.57
1:A:203:ASP:OD2	1:A:206:HIS:CE1	2.59	0.56
1:B:23:ALA:CB	1:B:83:MET:HE1	2.36	0.55
1:A:114:SER:O	1:A:138:PRO:HG3	2.06	0.54
1:B:2:PHE:C	1:B:3:LYS:HD2	2.32	0.54
1:B:142:ALA:HB3	1:B:149:ASP:OD1	2.08	0.54
1:B:80:TYR:HA	1:B:83:MET:HE3	1.89	0.54
1:B:23:ALA:CA	1:B:83:MET:HE1	2.39	0.53
1:B:45:LYS:HD3	1:B:45:LYS:C	2.34	0.53
1:B:8:PHD:O	1:B:9:LEU:CA	2.55	0.52
1:A:24:TRP:CE3	1:A:76:LYS:HG3	2.45	0.52
1:B:171:SER:O	1:B:175:ILE:HG12	2.10	0.52
1:B:41:ASN:HA	1:B:44:LEU:HD12	1.92	0.51
1:B:23:ALA:HB2	1:B:83:MET:CE	2.41	0.51
1:A:20:HIS:HD2	1:A:76:LYS:NZ	2.08	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:LYS:HD3	1:A:76:LYS:C	2.37	0.50
1:A:186:ILE:HD12	1:A:212:LEU:HD22	1.92	0.50
1:B:184:LEU:HD13	1:B:221:LYS:HG2	1.94	0.50
1:A:75:ARG:NE	3:A:851:HOH:O	2.46	0.49
1:B:30:GLU:OE2	1:B:75:ARG:NH2	2.45	0.49
1:B:21:PHE:O	1:B:25:LYS:HG3	2.12	0.49
1:B:112:LEU:HD22	1:B:123:LEU:HD11	1.94	0.49
1:B:24:TRP:CE2	1:B:76:LYS:HE3	2.47	0.49
1:A:8:PHD:OP3	1:A:10:ASP:HB3	2.12	0.49
1:A:190:ARG:NE	1:A:192:GLU:HB2	2.26	0.49
1:B:47:VAL:HG13	1:B:51:ASP:OD2	2.13	0.48
1:B:76:LYS:HD3	1:B:76:LYS:C	2.37	0.48
1:B:45:LYS:C	1:B:47:VAL:H	2.21	0.48
1:B:219:LYS:C	1:B:220:GLN:HG2	2.39	0.48
1:B:64:VAL:HB	1:B:68:GLU:HB2	1.97	0.47
1:A:75:ARG:NH2	3:A:935:HOH:O	2.47	0.47
1:A:145:LYS:O	1:A:147:ALA:O	2.33	0.46
1:B:88:SER:HB2	1:B:89:PRO:CD	2.45	0.46
1:B:21:PHE:HA	1:B:41:ASN:ND2	2.30	0.46
1:B:11:GLY:HA3	1:B:170:ASP:OD2	2.15	0.46
1:B:169:GLU:HG2	1:B:174:GLY:HA3	1.98	0.46
1:B:184:LEU:HB2	1:B:221:LYS:HG2	1.97	0.46
1:B:145:LYS:O	1:B:146:PRO:C	2.58	0.46
1:B:137:ASP:C	1:B:139:ALA:H	2.23	0.46
1:B:8:PHD:O	1:B:9:LEU:HA	2.14	0.45
1:B:65:SER:OG	1:B:68:GLU:HG3	2.17	0.45
1:A:20:HIS:HD2	1:A:76:LYS:HZ3	1.63	0.45
1:B:169:GLU:CD	1:B:174:GLY:HA3	2.42	0.45
1:B:88:SER:HB2	1:B:89:PRO:HD2	1.98	0.44
1:A:75:ARG:HB3	3:A:851:HOH:O	2.17	0.44
1:B:188:VAL:O	1:B:188:VAL:HG13	2.17	0.44
1:A:47:VAL:HG12	1:A:48:SER:N	2.33	0.44
1:B:178:ILE:CG2	1:B:185:PRO:HB3	2.48	0.44
1:B:21:PHE:CD1	1:B:41:ASN:HB2	2.53	0.43
1:A:47:VAL:HG12	1:A:51:ASP:HB2	2.01	0.43
1:A:145:LYS:HB3	1:A:146:PRO:HD3	2.00	0.43
1:A:88:SER:HB2	1:A:89:PRO:CD	2.49	0.43
1:B:11:GLY:HA3	1:B:170:ASP:CG	2.43	0.43
1:B:141:VAL:HG21	1:B:150:ILE:HD13	1.99	0.43
1:B:64:VAL:HG12	3:B:812:HOH:O	2.18	0.43
1:A:214:GLU:HG3	1:A:218:GLN:NE2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ARG:HH11	1:A:75:ARG:HA	1.85	0.42
1:A:4:ALA:HB2	1:A:109:LYS:HB2	2.01	0.42
1:A:188:VAL:HA	1:A:201:VAL:O	2.20	0.42
1:B:21:PHE:CE1	1:B:25:LYS:HD2	2.54	0.42
1:B:188:VAL:HA	1:B:201:VAL:O	2.20	0.42
1:B:3:LYS:HG3	3:B:819:HOH:O	2.20	0.41
1:A:43:GLN:HB3	1:A:55:LYS:HE3	2.03	0.41
1:B:4:ALA:HA	1:B:109:LYS:O	2.20	0.41
1:A:88:SER:HB2	1:A:89:PRO:HD2	2.02	0.41
1:B:23:ALA:HA	1:B:83:MET:HE1	2.03	0.41
1:A:24:TRP:CZ3	1:A:76:LYS:HG3	2.56	0.41
1:B:44:LEU:HD22	1:B:52:SER:OG	2.20	0.41
1:A:145:LYS:HA	1:A:145:LYS:HD3	1.87	0.41
1:B:145:LYS:HE3	1:B:151:PHE:CE1	2.56	0.41
1:B:192:GLU:CD	1:B:192:GLU:H	2.27	0.41
1:B:53:LEU:HD21	1:B:72:LEU:HB2	2.02	0.41
1:B:174:GLY:O	1:B:178:ILE:HG13	2.21	0.40
1:B:186:ILE:HD12	1:B:212:LEU:HD22	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ASP:OD2	1:B:203:ASP:OD2[3_645]	2.11	0.09

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	218/221 (99%)	207 (95%)	8 (4%)	3 (1%)	9	9
1	B	218/221 (99%)	200 (92%)	15 (7%)	3 (1%)	9	9

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	436/442 (99%)	407 (93%)	23 (5%)	6 (1%)	9 9

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	220	GLN
1	A	39	GLN
1	B	219	LYS
1	B	220	GLN
1	A	145	LYS
1	B	145	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	180/180 (100%)	175 (97%)	5 (3%)	38 56
1	B	180/180 (100%)	173 (96%)	7 (4%)	28 43
All	All	360/360 (100%)	348 (97%)	12 (3%)	33 50

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LEU
1	A	68	GLU
1	A	127	ASN
1	A	194	LEU
1	A	198	ILE
1	B	125	ARG
1	B	137	ASP
1	B	160	VAL
1	B	192	GLU
1	B	196	ASP
1	B	198	ILE
1	B	219	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	HIS
1	A	118	ASN
1	A	176	GLN
1	A	206	HIS
1	A	220	GLN
1	B	20	HIS
1	B	41	ASN
1	B	98	GLN
1	B	106	ASN
1	B	218	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	PHD	A	8	2,1	9,11,12	1.44	2 (22%)	9,15,17	2.13	2 (22%)
1	PHD	B	8	2,1	9,11,12	1.57	2 (22%)	9,15,17	1.82	2 (22%)

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	PHD	A	8	2,1	-	3/8/11/13	-
1	PHD	B	8	2,1	-	1/8/11/13	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	8	PHD	P-OD1	-2.78	1.54	1.59
1	A	8	PHD	P-OP2	-2.50	1.45	1.54
1	B	8	PHD	CB-CG	2.47	1.55	1.50
1	A	8	PHD	CB-CG	2.08	1.54	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	PHD	OD1-CG-CB	4.57	121.68	110.95
1	B	8	PHD	OD1-CG-CB	4.06	120.48	110.95
1	A	8	PHD	OD2-CG-CB	-3.66	116.15	124.65
1	B	8	PHD	OD2-CG-CB	-2.88	117.97	124.65

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	8	PHD	O-C-CA-CB
1	A	8	PHD	CG-OD1-P-OP1
1	A	8	PHD	C-CA-CB-CG
1	A	8	PHD	CA-CB-CG-OD2

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	8	PHD	1	0
1	B	8	PHD	5	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	8:PHD	C	9:LEU	N	1.74

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	220/221 (99%)	0.29	11 (5%) 34 35	13, 25, 61, 97	0
1	B	220/221 (99%)	1.37	56 (25%) 1 2	17, 45, 82, 89	0
All	All	440/442 (99%)	0.83	67 (15%) 5 6	13, 36, 79, 97	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	221	LYS	5.4
1	B	146	PRO	4.6
1	B	139	ALA	4.3
1	A	219	LYS	4.3
1	B	52	SER	4.0
1	A	221	LYS	3.9
1	A	220	GLN	3.9
1	B	145	LYS	3.8
1	B	72	LEU	3.8
1	B	46	GLY	3.8
1	B	141	VAL	3.8
1	B	73	ALA	3.7
1	B	142	ALA	3.7
1	A	145	LYS	3.5
1	B	216	TRP	3.4
1	A	47	VAL	3.3
1	A	32	GLY	3.1
1	B	195	GLY	3.0
1	B	24	TRP	3.0
1	B	26	ALA	3.0
1	B	217	LEU	3.0
1	B	76	LYS	3.0
1	B	51	ASP	3.0
1	B	50	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	47	VAL	2.9
1	B	80	TYR	2.9
1	B	220	GLN	2.9
1	B	23	ALA	2.8
1	B	42	GLU	2.7
1	B	219	LYS	2.7
1	A	60	ALA	2.6
1	B	81	VAL	2.6
1	B	137	ASP	2.6
1	B	59	LEU	2.6
1	B	218	GLN	2.6
1	B	31	ILE	2.6
1	B	48	SER	2.6
1	B	136	ALA	2.5
1	B	196	ASP	2.5
1	A	66	ALA	2.5
1	B	161	ALA	2.5
1	B	49	ARG	2.4
1	B	194	LEU	2.4
1	B	9	LEU	2.4
1	B	66	ALA	2.4
1	B	69	PHE	2.4
1	B	32	GLY	2.3
1	B	144	SER	2.3
1	A	75	ARG	2.3
1	B	56	ILE	2.3
1	B	64	VAL	2.3
1	B	138	PRO	2.3
1	B	53	LEU	2.2
1	B	71	GLU	2.2
1	B	3	LYS	2.2
1	B	36	VAL	2.2
1	B	60	ALA	2.2
1	B	77	ASN	2.2
1	B	115	ALA	2.1
1	B	57	LEU	2.1
1	B	140	GLU	2.0
1	A	59	LEU	2.0
1	B	78	ASP	2.0
1	B	17	ALA	2.0
1	B	41	ASN	2.0
1	A	146	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	108	ILE	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	PHD	A	8	12/13	0.86	0.15	14,18,22,25	0
1	PHD	B	8	12/13	0.91	0.12	20,27,31,33	0

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(Å ²)	Q<0.9
2	MG	A	801	1/1	0.95	0.26	1,1,1,1	0
2	MG	B	800	1/1	0.99	0.18	1,1,1,1	0

6.5 Other polymers [i](#)

There are no such residues in this entry.