



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:48 AM UTC

PDB ID : 1IME / pdb_00001ime
Title : STRUCTURAL STUDIES OF METAL BINDING BY INOSITOL MONOPHOSPHATASE: EVIDENCE FOR TWO-METAL ION CATALYSIS
Authors : Bone, R.
Deposited on : 1994-02-08
Resolution : 2.25 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

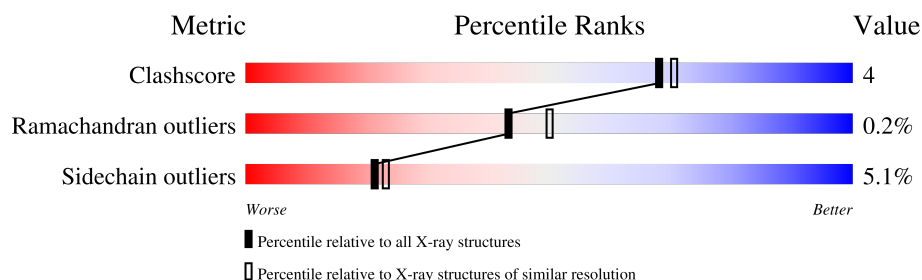
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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))
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	277	
1	B	277	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4331 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 INOSITOL MONOPHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	55	0	1
			2073	1307	352	396	18			
1	B	273	Total	C	N	O	S	62	0	1
			2073	1307	352	396	18			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		
3	B	91	Total	O	0	0
			91	91		

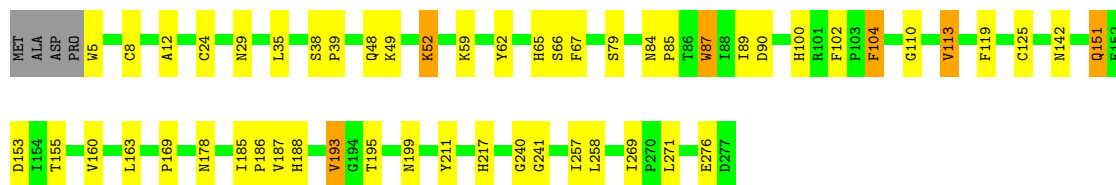
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

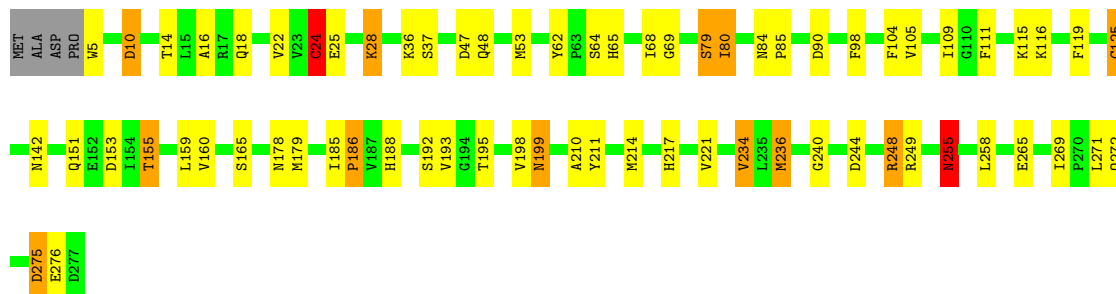
• Molecule 1: INOSITOL MONOPHOSPHATASE

Chain A: 



• Molecule 1: INOSITOL MONOPHOSPHATASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	86.00Å 86.00Å 154.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.25	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.25)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.184 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4331	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

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

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	1.01	5/2106 (0.2%)	1.68	36/2849 (1.3%)
1	B	1.02	5/2106 (0.2%)	1.75	45/2849 (1.6%)
All	All	1.02	10/4212 (0.2%)	1.72	81/5698 (1.4%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	188	HIS	CD2-NE2	-7.23	1.29	1.37
1	A	276	GLU	C-N	-7.08	1.23	1.33
1	B	188	HIS	CD2-NE2	-6.88	1.30	1.37
1	B	65	HIS	CD2-NE2	-6.75	1.30	1.37
1	B	217	HIS	CD2-NE2	-6.54	1.30	1.37
1	A	100	HIS	CD2-NE2	-6.54	1.30	1.37
1	B	276	GLU	C-N	-6.47	1.24	1.33
1	A	65	HIS	CD2-NE2	-6.07	1.31	1.37
1	B	221	VAL	CA-CB	5.61	1.61	1.54
1	A	217	HIS	CD2-NE2	-5.25	1.32	1.37

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	HIS	CB-CG-CD2	-8.79	119.77	131.20
1	B	119	PHE	CA-CB-CG	8.64	122.44	113.80
1	A	104	PHE	CA-CB-CG	-7.84	105.96	113.80
1	A	142	ASN	OD1-CG-ND2	-7.70	114.90	122.60
1	B	217	HIS	CB-CG-ND1	7.45	133.87	122.70
1	B	188	HIS	CB-CG-CD2	-7.29	121.73	131.20
1	B	199	ASN	OD1-CG-ND2	-7.27	115.33	122.60
1	A	269	ILE	CA-C-O	-7.19	114.52	119.19
1	B	105	VAL	CG1-CB-CG2	-7.14	95.10	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	236	MET	CG-SD-CE	-7.07	85.34	100.90
1	B	244	ASP	CA-CB-CG	6.84	119.44	112.60
1	A	119	PHE	CA-CB-CG	6.75	120.55	113.80
1	B	62	TYR	CA-C-N	6.73	127.28	119.47
1	B	62	TYR	C-N-CA	6.73	127.28	119.47
1	A	38	SER	N-CA-C	-6.68	102.12	108.07
1	B	178	ASN	OD1-CG-ND2	-6.58	116.03	122.60
1	A	178	ASN	OD1-CG-ND2	-6.51	116.09	122.60
1	A	102	PHE	N-CA-C	-6.51	101.58	109.72
1	B	48	GLN	OE1-CD-NE2	-6.50	116.09	122.60
1	A	185	ILE	O-C-N	-6.43	116.31	120.42
1	B	188	HIS	CB-CG-ND1	6.41	132.32	122.70
1	A	48	GLN	OE1-CD-NE2	-6.35	116.25	122.60
1	B	234	VAL	N-CA-CB	-6.35	100.05	111.92
1	A	217	HIS	CB-CG-CD2	-6.32	122.98	131.20
1	A	151	GLN	OE1-CD-NE2	-6.28	116.32	122.60
1	B	65	HIS	N-CA-C	6.28	119.03	110.55
1	B	68	ILE	N-CA-C	-6.27	99.34	108.11
1	A	29	ASN	OD1-CG-ND2	-6.24	116.36	122.60
1	B	255	ASN	OD1-CG-ND2	-6.23	116.37	122.60
1	B	155	THR	CA-C-O	6.09	126.91	119.11
1	A	100	HIS	CB-CG-CD2	-5.95	123.47	131.20
1	B	248	ARG	CA-CB-CG	-5.95	102.21	114.10
1	A	8	CYS	N-CA-C	-5.93	104.72	111.07
1	B	125	CYS	N-CA-C	5.84	117.31	111.07
1	A	89	ILE	N-CA-C	5.82	116.86	108.48
1	A	113	VAL	N-CA-C	-5.79	99.84	108.46
1	B	79	SER	O-C-N	-5.75	115.52	122.65
1	A	188	HIS	CB-CG-CD2	-5.75	123.73	131.20
1	B	151	GLN	N-CA-C	5.65	118.52	110.10
1	B	198	VAL	N-CA-C	-5.65	105.01	110.72
1	B	248	ARG	CB-CG-CD	5.65	124.30	111.30
1	A	84	ASN	OD1-CG-ND2	-5.64	116.96	122.60
1	B	24	CYS	CA-CB-SG	-5.64	101.43	114.40
1	A	240	GLY	N-CA-C	-5.64	107.88	115.21
1	B	5	TRP	CG-CD2-CE3	5.63	139.53	133.90
1	B	69	GLY	O-C-N	-5.58	118.39	123.36
1	A	29	ASN	CB-CG-ND2	5.57	124.76	116.40
1	A	62	TYR	CA-C-N	5.55	126.78	119.84
1	A	62	TYR	C-N-CA	5.55	126.78	119.84
1	A	110	GLY	O-C-N	-5.54	116.92	123.18
1	A	90	ASP	CA-C-N	5.48	124.83	118.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	ASP	C-N-CA	5.48	124.83	118.85
1	B	142	ASN	OD1-CG-ND2	-5.48	117.12	122.60
1	B	10	ASP	CA-CB-CG	5.48	118.08	112.60
1	B	153	ASP	CA-CB-CG	5.43	118.03	112.60
1	A	187	VAL	N-CA-C	-5.41	101.28	108.96
1	A	217	HIS	CB-CG-ND1	5.39	130.79	122.70
1	A	257	ILE	N-CA-C	-5.33	105.34	110.72
1	A	241	GLY	CA-C-N	5.31	125.25	119.78
1	A	241	GLY	C-N-CA	5.31	125.25	119.78
1	B	37	SER	N-CA-C	5.29	119.58	113.18
1	B	90	ASP	CA-C-N	5.25	124.57	118.85
1	B	90	ASP	C-N-CA	5.25	124.57	118.85
1	B	47	ASP	CA-CB-CG	5.23	117.83	112.60
1	B	80	ILE	N-CA-C	-5.22	100.86	108.17
1	A	199	ASN	OD1-CG-ND2	-5.20	117.40	122.60
1	A	5	TRP	CE2-CD2-CG	-5.16	101.01	107.20
1	B	98	PHE	CA-C-N	5.14	127.50	120.46
1	B	98	PHE	C-N-CA	5.14	127.50	120.46
1	A	87	TRP	CG-CD2-CE3	5.14	139.04	133.90
1	B	275	ASP	CA-CB-CG	5.13	117.73	112.60
1	A	193	VAL	N-CA-CB	-5.12	100.33	111.81
1	A	5	TRP	CB-CG-CD1	-5.10	119.25	126.90
1	B	269	ILE	CA-C-N	5.10	125.34	119.83
1	B	269	ILE	C-N-CA	5.10	125.34	119.83
1	B	271	LEU	O-C-N	-5.09	117.44	123.25
1	B	5	TRP	CB-CG-CD1	-5.08	119.28	126.90
1	B	192	SER	CB-CA-C	-5.07	103.74	110.79
1	B	276	GLU	CA-CB-CG	5.04	124.18	114.10
1	A	269	ILE	N-CA-CB	-5.04	104.81	111.21
1	B	36	LYS	O-C-N	-5.03	115.89	122.59

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	2073	0	2090	12	0
1	B	2073	0	2090	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	92	0	0	0	0
3	B	91	0	0	0	0
All	All	4331	0	4180	29	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 4.

All (29) 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:151:GLN:HE21	1:A:153:ASP:H	1.28	0.79
1:A:49:LYS:HA	1:A:52:LYS:HD2	1.63	0.79
1:B:236:MET:HE3	1:B:240:GLY:HA2	1.68	0.75
1:B:22:VAL:HG21	1:B:53:MET:HE1	1.76	0.68
1:B:214:MET:HG2	1:B:249:ARG:HG2	1.79	0.65
1:B:80:ILE:HD12	1:B:275:ASP:HB2	1.79	0.64
1:B:185:ILE:HD11	1:B:265:GLU:HG3	1.82	0.62
1:B:25:GLU:O	1:B:28:LYS:HG2	2.03	0.58
1:B:160:VAL:HG23	1:B:211:TYR:HB2	1.86	0.57
1:A:155:THR:HA	1:A:186:PRO:O	2.08	0.54
1:B:24:CYS:SG	1:B:125:CYS:HB3	2.46	0.54
1:A:24:CYS:SG	1:A:125:CYS:HB3	2.49	0.53
1:B:159:LEU:HD22	1:B:210:ALA:HB3	1.92	0.52
1:B:85:PRO:HB3	1:B:111:PHE:CZ	2.45	0.52
1:A:151:GLN:NE2	1:A:153:ASP:HB3	2.28	0.49
1:A:160:VAL:HG23	1:A:211:TYR:HB2	1.94	0.49
1:B:193:VAL:HG22	1:B:199:ASN:ND2	2.27	0.48
1:B:248:ARG:HD3	1:B:272:GLN:O	2.14	0.47
1:A:59:LYS:HE2	1:A:67:PHE:CE1	2.50	0.46
1:B:16:ALA:HB2	1:B:109:ILE:HD12	1.97	0.46
1:A:35:LEU:HD13	1:A:39:PRO:HA	1.96	0.46
1:A:151:GLN:HE21	1:A:153:ASP:N	2.05	0.45
1:B:155:THR:HA	1:B:186:PRO:O	2.19	0.42
1:B:84:ASN:OD1	1:B:85:PRO:HD2	2.19	0.42
1:A:12:ALA:HB2	1:A:87:TRP:HZ3	1.84	0.42
1:A:104:PHE:CZ	1:B:104:PHE:HZ	2.37	0.42
1:B:255:ASN:ND2	1:B:258:LEU:H	2.18	0.41
1:A:163:LEU:HD11	1:B:179:MET:HE3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:ARG:HD3	1:B:248:ARG:HH11	1.70	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	271/277 (98%)	263 (97%)	7 (3%)	1 (0%)	30	31
1	B	271/277 (98%)	260 (96%)	11 (4%)	0	100	100
All	All	542/554 (98%)	523 (96%)	18 (3%)	1 (0%)	43	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	85	PRO

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	227/231 (98%)	218 (96%)	9 (4%)	28	34
1	B	227/231 (98%)	213 (94%)	14 (6%)	16	16
All	All	454/462 (98%)	431 (95%)	23 (5%)	21	23

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LYS
1	A	66	SER
1	A	79	SER
1	A	113	VAL
1	A	169	PRO
1	A	193	VAL
1	A	195	THR
1	A	258	LEU
1	A	271	LEU
1	B	10	ASP
1	B	14	THR
1	B	18	GLN
1	B	24	CYS
1	B	28	LYS
1	B	64	SER
1	B	79	SER
1	B	115	LYS
1	B	116	LYS
1	B	165	SER
1	B	186	PRO
1	B	195	THR
1	B	234	VAL
1	B	255	ASN

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

Mol	Chain	Res	Type
1	A	114	ASN
1	A	142	ASN
1	A	147	GLN
1	A	151	GLN
1	B	151	GLN
1	B	199	ASN
1	B	255	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.