



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 26, 2016 – 10:40 AM EDT

PDB ID : 5I3S
Title : Crystal structure of Staphylococcal IMPase-II
Authors : Dutta, A.; Bhattacharyya, S.; Dutta, D.; Das, A.K.
Deposited on : 2016-02-11
Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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.

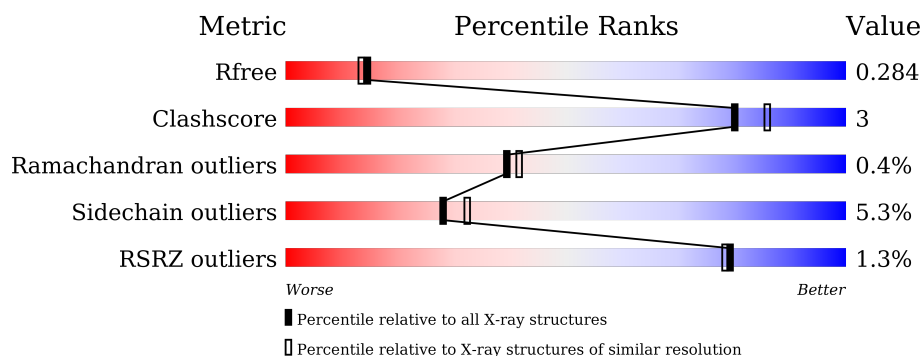
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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-20027939

i

X-RAY DIFFRACTION

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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	275	<div> <div></div> <div>82%</div> <div>11%</div> <div>5%</div> </div>
1	B	275	<div> <div></div> <div>84%</div> <div>8%</div> <div>6%</div> </div>
1	C	275	<div> <div></div> <div>84%</div> <div>9%</div> <div>6%</div> </div>
1	D	275	<div> <div></div> <div>80%</div> <div>8%</div> <div>9%</div> </div>

2 Entry composition [i](#)

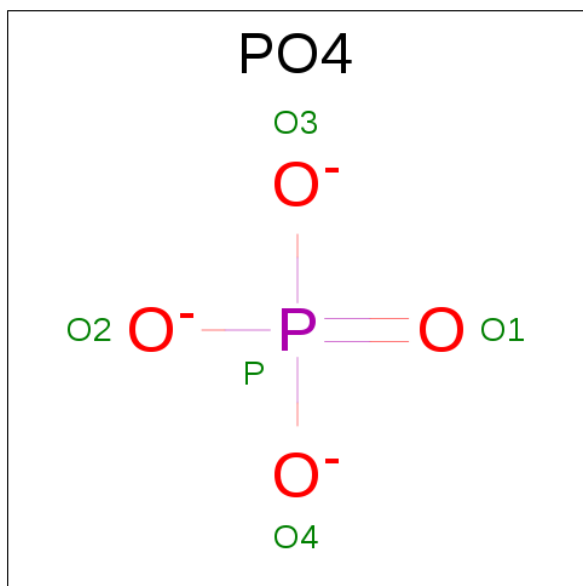
There are 3 unique types of molecules in this entry. The entry contains 8200 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	0	0
			1996	1274	336	383	3			
1	B	258	Total	C	N	O	S	0	0	0
			1942	1241	324	374	3			
1	C	258	Total	C	N	O	S	0	0	0
			1971	1260	332	376	3			
1	D	249	Total	C	N	O	S	0	0	0
			1867	1193	311	360	3			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

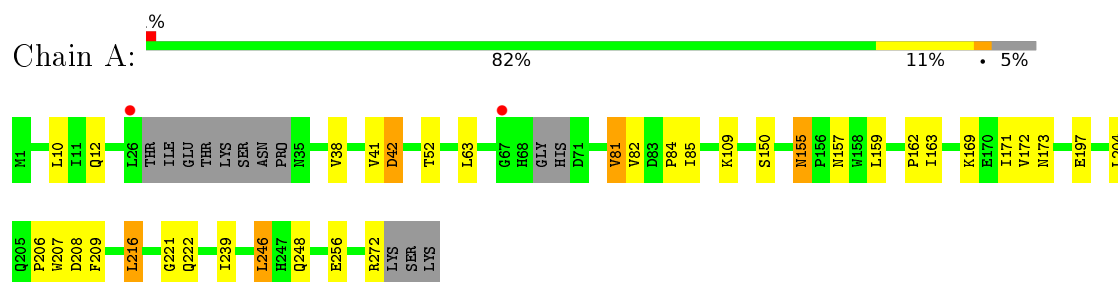
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	107	Total	O	0	0
			107	107		
3	B	95	Total	O	0	0
			95	95		
3	C	98	Total	O	0	0
			98	98		
3	D	104	Total	O	0	0
			104	104		

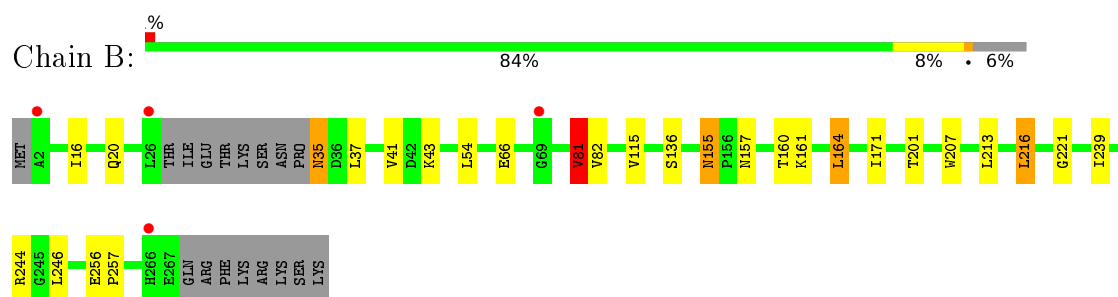
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

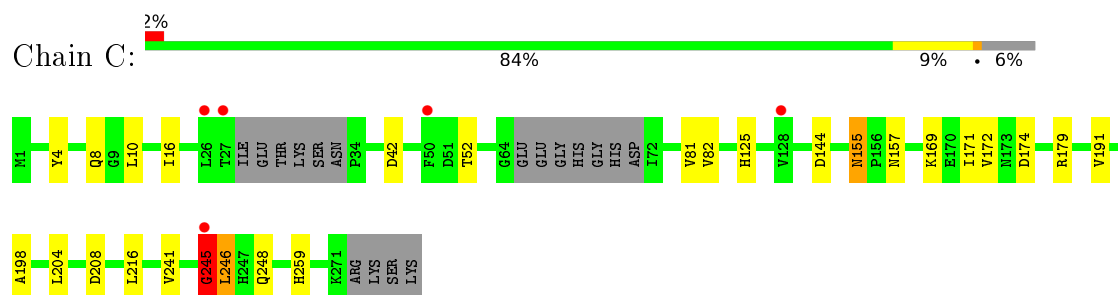
- Molecule 1: Inositol monophosphatase family protein



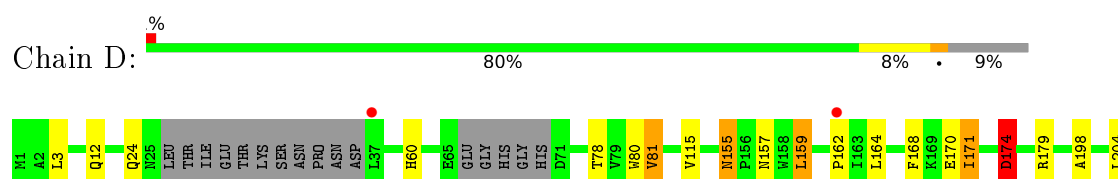
- Molecule 1: Inositol monophosphatase family protein



- Molecule 1: Inositol monophosphatase family protein



- Molecule 1: Inositol monophosphatase family protein



D209	L213	L216	N220	A223	R244	R245	L246	H247	S251	L265	HIS	GLU	GLN	ARG	PHE	LYS	ARG	LYS	SER	LYS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.00 Å 131.39 Å 80.90 Å 90.00° 104.43° 90.00°	Depositor
Resolution (Å)	78.35 – 2.20 19.59 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.1 (78.35-2.20) 97.3 (19.59-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.21 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.229 , 0.283 0.234 , 0.284	Depositor DCC
R_{free} test set	2812 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.4	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	0.90	EDS
Total number of atoms	8200	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: PO4

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	0.78	0/2037	0.87	1/2773 (0.0%)
1	B	0.79	0/1983	0.87	3/2705 (0.1%)
1	C	0.79	1/2012 (0.0%)	0.90	5/2741 (0.2%)
1	D	0.80	1/1905 (0.1%)	0.91	4/2601 (0.2%)
All	All	0.79	2/7937 (0.0%)	0.89	13/10820 (0.1%)

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	C	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	174	ASP	CB-CG	-6.63	1.37	1.51
1	D	174	ASP	CB-CG	-5.83	1.39	1.51

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	174	ASP	CB-CG-OD2	-8.55	110.61	118.30
1	C	144	ASP	CB-CG-OD1	6.98	124.58	118.30
1	D	179	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	D	179	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	B	81	VAL	CB-CA-C	5.50	121.86	111.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	245	GLY	Peptide

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	1996	0	1913	18	0
1	B	1942	0	1855	13	0
1	C	1971	0	1898	14	0
1	D	1867	0	1773	13	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	107	0	0	1	0
3	B	95	0	0	1	0
3	C	98	0	0	0	0
3	D	104	0	0	0	0
All	All	8200	0	7439	52	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 3.

The worst 5 of 52 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:12:GLN:HE22	1:D:12:GLN:HE22	1.32	0.76
1:C:245:GLY:O	1:C:248:GLN:HB2	1.92	0.70
1:B:216:LEU:HD22	1:B:221:GLY:HA3	1.79	0.65
1:A:10:LEU:HD13	1:A:52:THR:HG21	1.83	0.61
1:A:169:LYS:HG3	1:B:160:THR:HG22	1.83	0.60

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	256/275 (93%)	248 (97%)	8 (3%)	0	100	100
1	B	254/275 (92%)	248 (98%)	6 (2%)	0	100	100
1	C	252/275 (92%)	238 (94%)	12 (5%)	2 (1%)	24	22
1	D	243/275 (88%)	237 (98%)	4 (2%)	2 (1%)	24	22
All	All	1005/1100 (91%)	971 (97%)	30 (3%)	4 (0%)	39	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	246	LEU
1	D	24	GLN
1	C	245	GLY
1	D	162	PRO

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	206/232 (89%)	194 (94%)	12 (6%)	25	28
1	B	198/232 (85%)	185 (93%)	13 (7%)	21	22
1	C	205/232 (88%)	199 (97%)	6 (3%)	50	62
1	D	190/232 (82%)	179 (94%)	11 (6%)	25	28
All	All	799/928 (86%)	757 (95%)	42 (5%)	28	32

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	155	ASN
1	B	216	LEU
1	D	216	LEU
1	B	164	LEU
1	B	201	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	173	ASN
1	B	266	HIS
1	D	173	ASN
1	B	222	GLN
1	C	8	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](#)

4 ligands 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	$\# Z > 2$
2	PO4	A	301	-	4,4,4	0.92	0	6,6,6	0.27	0
2	PO4	B	301	-	4,4,4	0.80	0	6,6,6	0.26	0
2	PO4	C	301	-	4,4,4	0.40	0	6,6,6	0.31	0
2	PO4	D	301	-	4,4,4	0.91	0	6,6,6	0.28	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
2	PO4	A	301	-	-	0/0/0/0	0/0/0/0
2	PO4	B	301	-	-	0/0/0/0	0/0/0/0
2	PO4	C	301	-	-	0/0/0/0	0/0/0/0
2	PO4	D	301	-	-	0/0/0/0	0/0/0/0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	262/275 (95%)	-0.16	2 (0%) 87 87	13, 26, 49, 67	0
1	B	258/275 (93%)	-0.09	4 (1%) 74 73	13, 26, 49, 67	0
1	C	258/275 (93%)	-0.08	5 (1%) 70 68	13, 29, 49, 67	0
1	D	249/275 (90%)	-0.18	2 (0%) 87 87	13, 25, 49, 62	0
All	All	1027/1100 (93%)	-0.13	13 (1%) 79 78	13, 26, 49, 67	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	26	LEU	4.4
1	C	50	PHE	2.7
1	A	67	GLY	2.7
1	B	2	ALA	2.6
1	C	27	THR	2.5

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. 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
2	PO4	D	301	5/5	0.98	0.08	-0.95	15,16,17,17	0
2	PO4	C	301	5/5	0.98	0.07	-0.97	15,16,17,18	0
2	PO4	A	301	5/5	0.99	0.07	-1.24	19,19,22,22	0
2	PO4	B	301	5/5	0.99	0.07	-1.48	18,19,20,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.