



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 8, 2016 – 07:32 PM EST

PDB ID : 5LZQ  
Title : Crystal structure of Thermotoga maritima sodium pumping membrane integral pyrophosphatase in complex with imidodiphosphate and magnesium, and with bound sodium ion  
Authors : Wilkinson, C.; Kellosalo, J.; Kajander, T.; Goldman, A.  
Deposited on : 2016-10-01  
Resolution : 3.50 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](mailto: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.

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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-20028442  
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-20028442

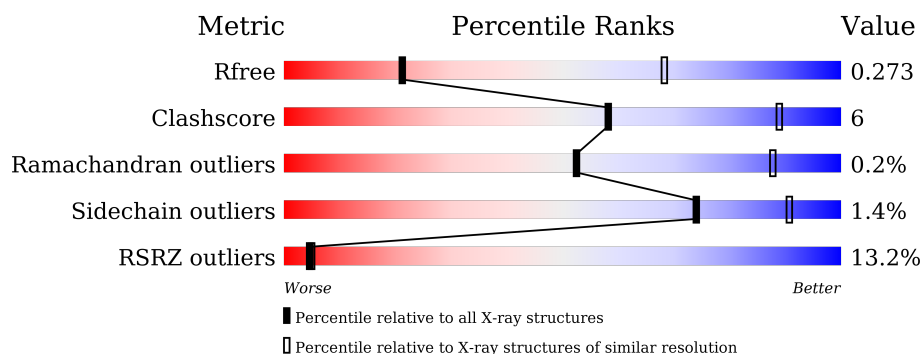
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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))
$R_{free}$	91344	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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  $\geq 3$ , 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	735	<div> <div>13%</div> <div>83%</div> <div>15%</div> <div>.</div> </div>
1	B	735	<div> <div>13%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MG	A	801	-	-	-	X
2	MG	A	802	-	-	-	X
2	MG	A	803	-	-	-	X
2	MG	B	801	-	-	-	X
2	MG	B	802	-	-	-	X
2	MG	B	805	-	-	-	X
3	2PN	A	806	-	-	-	X
3	2PN	A	809	-	-	-	X
3	2PN	B	806	-	-	-	X
4	NA	B	807	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10448 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 K(+)-stimulated pyrophosphate-energized sodium pump.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	725	Total	C	N	O	S	0	0	0
			5224	3423	816	958	27			
1	B	725	Total	C	N	O	S	0	0	0
			5170	3387	805	953	25			

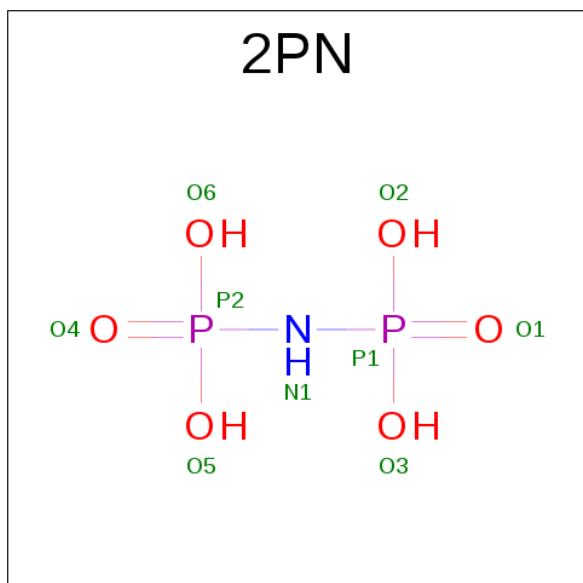
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MET	-	initiating methionine	UNP Q9S5X0
A	-7	ARG	-	expression tag	UNP Q9S5X0
A	-6	GLY	-	expression tag	UNP Q9S5X0
A	-5	SER	-	expression tag	UNP Q9S5X0
A	-4	HIS	-	expression tag	UNP Q9S5X0
A	-3	HIS	-	expression tag	UNP Q9S5X0
A	-2	HIS	-	expression tag	UNP Q9S5X0
A	-1	HIS	-	expression tag	UNP Q9S5X0
A	0	HIS	-	expression tag	UNP Q9S5X0
A	1	HIS	-	expression tag	UNP Q9S5X0
A	353	LEU	VAL	engineered mutation	UNP Q9S5X0
A	395	GLY	SER	engineered mutation	UNP Q9S5X0
B	-8	MET	-	initiating methionine	UNP Q9S5X0
B	-7	ARG	-	expression tag	UNP Q9S5X0
B	-6	GLY	-	expression tag	UNP Q9S5X0
B	-5	SER	-	expression tag	UNP Q9S5X0
B	-4	HIS	-	expression tag	UNP Q9S5X0
B	-3	HIS	-	expression tag	UNP Q9S5X0
B	-2	HIS	-	expression tag	UNP Q9S5X0
B	-1	HIS	-	expression tag	UNP Q9S5X0
B	0	HIS	-	expression tag	UNP Q9S5X0
B	1	HIS	-	expression tag	UNP Q9S5X0
B	353	LEU	VAL	engineered mutation	UNP Q9S5X0
B	395	GLY	SER	engineered mutation	UNP Q9S5X0

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	5	Total	Mg	0	0
			5	5		
2	A	5	Total	Mg	0	0
			5	5		

- Molecule 3 is IMIDODIPHOSPHORIC ACID (three-letter code: 2PN) (formula:  $\text{H}_5\text{NO}_6\text{P}_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	N	O	P	0	0
			9	1	6	2		
3	A	1	Total	N	O	P	0	0
			9	1	6	2		
3	A	1	Total	N	O	P	0	0
			9	1	6	2		
3	B	1	Total	N	O	P	0	0
			9	1	6	2		

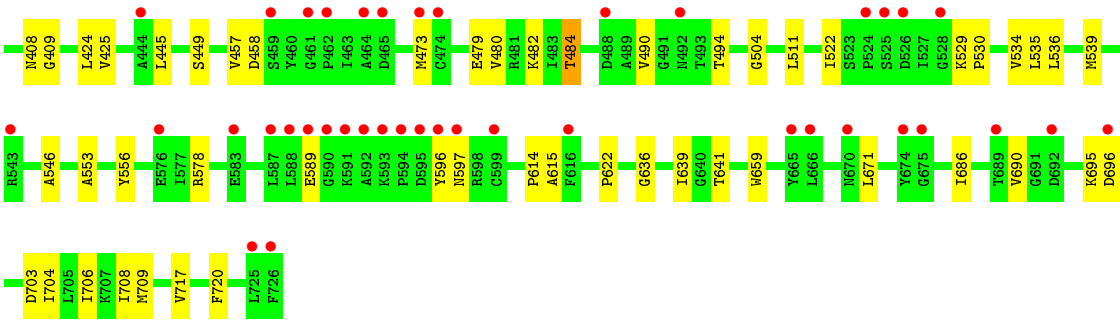
- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Na	0	0
			1	1		
4	A	1	Total	Na	0	0
			1	1		

- Molecule 5 is water.

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







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.42Å 106.76Å 161.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.75 – 3.50 29.75 – 3.49	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.75-3.50) 99.8 (29.75-3.49)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.18 (at 3.47Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.245 , 0.273 0.245 , 0.273	Depositor DCC
$R_{free}$ test set	1195 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	148.7	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 104.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.028 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	10448	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	136.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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: NA, MG, 2PN

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	0.23	0/5337	0.40	0/7295
1	B	0.23	0/5281	0.40	0/7229
All	All	0.23	0/10618	0.40	0/14524

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 [i](#)

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	5224	0	5166	67	0
1	B	5170	0	5076	63	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	27	0	2	0	0
3	B	9	0	1	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	2	0	0	0	0
5	B	4	0	0	0	0
All	All	10448	0	10245	121	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.

The worst 5 of 121 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:539:MET:HB2	1:B:539:MET:HE3	1.69	0.74
1:A:636:GLY:HA2	1:A:639:ILE:HD12	1.79	0.65
1:B:424:LEU:HD21	1:B:504:GLY:HA2	1.79	0.65
1:A:546:ALA:HB1	1:B:425:VAL:HG13	1.80	0.64
1:A:14:ALA:HB2	1:A:298:LEU:HD21	1.80	0.62

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	723/735 (98%)	696 (96%)	27 (4%)	0	100	100
1	B	723/735 (98%)	692 (96%)	28 (4%)	3 (0%)	39	81
All	All	1446/1470 (98%)	1388 (96%)	55 (4%)	3 (0%)	52	88

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	310	VAL
1	B	30	GLU
1	B	311	LYS

### 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	517/575 (90%)	511 (99%)	6 (1%)	78	92
1	B	507/575 (88%)	499 (98%)	8 (2%)	70	89
All	All	1024/1150 (89%)	1010 (99%)	14 (1%)	74	91

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	17	PHE
1	B	22	PHE
1	B	484	THR
1	A	703	ASP
1	B	261	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 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	2PN	A	806	2	8,8,8	2.90	2 (25%)	6,13,13	1.33	0
3	2PN	A	808	-	8,8,8	2.90	2 (25%)	6,13,13	1.35	0
3	2PN	A	809	-	8,8,8	2.88	2 (25%)	6,13,13	1.35	0
3	2PN	B	806	2	8,8,8	2.89	2 (25%)	6,13,13	1.31	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	2PN	A	806	2	-	0/2/6/6	0/0/0/0
3	2PN	A	808	-	-	0/2/6/6	0/0/0/0
3	2PN	A	809	-	-	0/2/6/6	0/0/0/0
3	2PN	B	806	2	-	0/2/6/6	0/0/0/0

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	809	2PN	P1-O1	4.93	1.51	1.46
3	B	806	2PN	P2-O4	4.94	1.51	1.46
3	A	808	2PN	P1-O1	4.96	1.51	1.46
3	A	806	2PN	P1-O1	4.98	1.51	1.46
3	A	809	2PN	P2-O4	4.99	1.51	1.46

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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	725/735 (98%)	0.63	98 (13%) 4 5	90, 130, 161, 182	0
1	B	725/735 (98%)	0.63	94 (12%) 5 5	94, 141, 180, 204	0
All	All	1450/1470 (98%)	0.63	192 (13%) 4 5	90, 135, 173, 204	0

The worst 5 of 192 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	465	ASP	6.2
1	A	670	ASN	6.1
1	B	726	PHE	6.1
1	B	202	ASP	6.0
1	B	594	PRO	5.9

### 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	MG	B	802	1/1	0.87	2.10	19.86	155,155,155,155	0
2	MG	A	802	1/1	0.96	1.61	8.38	126,126,126,126	0
3	2PN	A	809	9/9	0.76	0.50	3.74	146,161,184,192	0
4	NA	B	807	1/1	0.81	0.44	3.19	125,125,125,125	0
2	MG	A	801	1/1	0.95	0.67	1.98	127,127,127,127	0
3	2PN	A	806	9/9	0.93	0.67	1.33	142,145,156,164	0
2	MG	A	803	1/1	0.94	0.58	1.20	104,104,104,104	0
2	MG	B	805	1/1	0.95	0.72	0.86	140,140,140,140	0
3	2PN	A	808	9/9	0.67	0.34	0.54	182,189,195,202	0
2	MG	B	801	1/1	0.88	0.59	0.39	136,136,136,136	0
3	2PN	B	806	9/9	0.91	0.57	0.23	153,159,174,179	0
4	NA	A	807	1/1	0.82	0.22	-0.48	115,115,115,115	0
2	MG	A	804	1/1	0.91	0.75	-	130,130,130,130	0
2	MG	B	803	1/1	0.92	0.76	-	149,149,149,149	0
2	MG	A	805	1/1	0.97	0.39	-	130,130,130,130	0
2	MG	B	804	1/1	0.92	0.38	-	138,138,138,138	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.