



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:53 PM UTC

PDB ID : 1PC3 / pdb_00001pc3
Title : Crystal structure of the extracellular phosphate ABC transport receptor (PstS-1) and immunodominant antigen of M. tuberculosis.
Authors : Vyas, N.K.; Vyas, M.N.; Quiocho, F.A.
Deposited on : 2003-05-15
Resolution : 2.16 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)	:	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

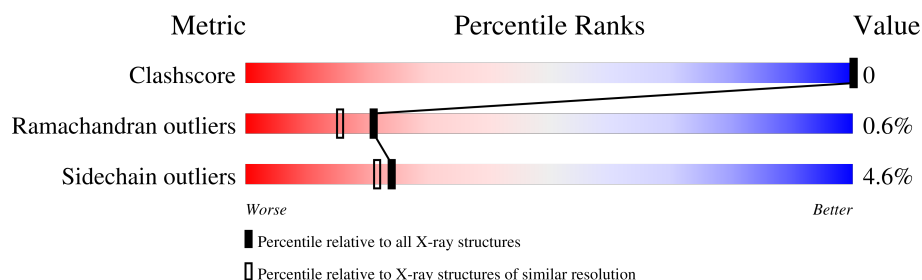
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 Å.

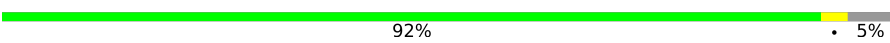
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)

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	1-A	350	 88% 6% • 5%
1	1-B	350	 92% • 5%
1	2-A	350	 88% 6% • 5%
1	2-B	350	 92% • 5%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10196 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 Phosphate-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1-A	333	Total	C	N	O	S	0	0	0
			2424	1532	408	477	7			
1	2-A	333	Total	C	N	O	S	0	0	0
			2424	1532	408	477	7			
1	1-B	333	Total	C	N	O	S	0	0	0
			2424	1532	408	477	7			
1	2-B	333	Total	C	N	O	S	0	0	0
			2424	1532	408	477	7			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	1-A	225	Total	O	0	0
			225	225		
3	2-A	222	Total	O	0	0
			222	222		
3	1-B	15	Total	O	0	0
			15	15		
3	2-B	18	Total	O	0	0
			18	18		

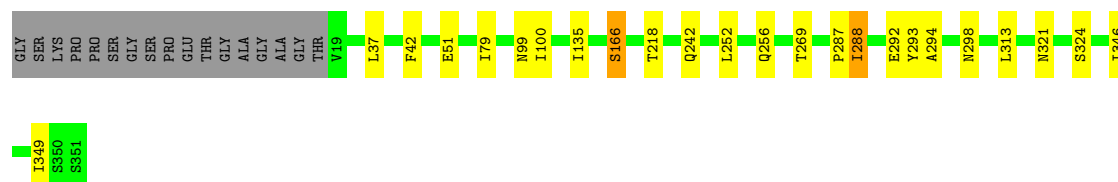
3 Residue-property plots [i](#)

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. 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.

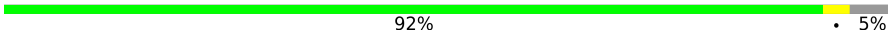
Note EDS was not executed.

- Molecule 1: Phosphate-binding protein 1

Chain 1-A:  88% 6% • 5%



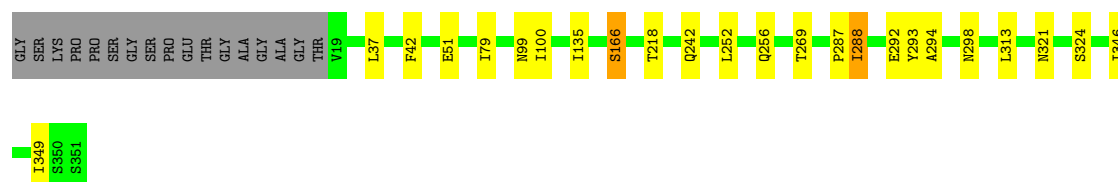
- Molecule 1: Phosphate-binding protein 1

Chain 1-B:  92% • 5%

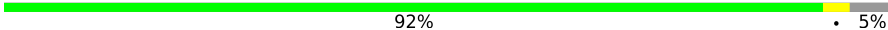


- Molecule 1: Phosphate-binding protein 1

Chain 2-A:  88% 6% • 5%



- Molecule 1: Phosphate-binding protein 1

Chain 2-B:  92% • 5%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	125.40Å 72.30Å 73.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	18.64 – 2.16	Depositor
% Data completeness (in resolution range)	86.0 (18.64-2.16)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.220 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10196	wwPDB-VP
Average B, all atoms (Å ²)	10.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	1-A	0.47	0/2482	0.99	10/3394 (0.3%)
1	1-B	0.25	0/2482	0.66	1/3394 (0.0%)
1	2-A	0.47	0/2482	0.99	10/3394 (0.3%)
1	2-B	0.25	0/2482	0.66	1/3394 (0.0%)
All	All	0.38	0/9928	0.84	22/13576 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1-A	242	GLN	N-CA-C	-6.26	98.53	108.73
1	2-A	242	GLN	N-CA-C	-6.26	98.53	108.73
1	1-A	166	SER	N-CA-C	6.23	119.65	109.06
1	2-A	166	SER	N-CA-C	6.23	119.65	109.06
1	1-A	294	ALA	N-CA-C	-5.47	99.82	108.73
1	2-A	294	ALA	N-CA-C	-5.47	99.82	108.73
1	1-A	288	ILE	CB-CA-C	-5.39	105.33	111.59
1	2-A	288	ILE	CB-CA-C	-5.39	105.33	111.59
1	1-B	83	ASP	N-CA-C	-5.33	106.77	113.28
1	2-B	83	ASP	N-CA-C	-5.33	106.77	113.28
1	1-A	287	PRO	N-CA-C	5.33	120.52	113.86
1	2-A	287	PRO	N-CA-C	5.33	120.52	113.86
1	1-A	269	THR	CA-C-N	5.20	125.17	120.03
1	1-A	269	THR	C-N-CA	5.20	125.17	120.03
1	2-A	269	THR	CA-C-N	5.20	125.17	120.03
1	2-A	269	THR	C-N-CA	5.20	125.17	120.03
1	1-A	292	GLU	N-CA-C	-5.10	100.99	109.46
1	2-A	292	GLU	N-CA-C	-5.10	100.99	109.46
1	1-A	293	TYR	N-CA-C	5.06	117.73	109.59
1	2-A	293	TYR	N-CA-C	5.06	117.73	109.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1-A	346	ILE	CB-CA-C	-5.00	105.49	112.04
1	2-A	346	ILE	CB-CA-C	-5.00	105.49	112.04

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	1-A	2424	0	2356	0	0
1	1-B	2424	0	2356	0	0
1	2-A	2424	0	2356	0	0
1	2-B	2424	0	2356	0	0
2	1-A	5	0	0	0	0
2	1-B	5	0	0	0	0
2	2-A	5	0	0	0	0
2	2-B	5	0	0	0	0
3	1-A	225	0	0	0	0
3	1-B	15	0	0	0	0
3	2-A	222	0	0	0	0
3	2-B	18	0	0	0	0
All	All	10196	0	9424	0	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 0.

There are no clashes within the asymmetric unit.

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	1-A	331/350 (95%)	323 (98%)	8 (2%)	0	100	100
1	1-B	331/350 (95%)	298 (90%)	28 (8%)	5 (2%)	8	4
1	2-A	331/350 (95%)	323 (98%)	8 (2%)	0	100	100
1	2-B	331/350 (95%)	310 (94%)	18 (5%)	3 (1%)	14	9
All	All	1324/1400 (95%)	1254 (95%)	62 (5%)	8 (1%)	21	15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1-B	76	THR
1	2-B	116	VAL
1	1-B	35	SER
1	1-B	75	GLY
1	1-B	151	ASN
1	2-B	35	SER
1	2-B	301	LYS
1	1-B	74	ALA

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	1-A	249/259 (96%)	232 (93%)	17 (7%)	14	10
1	1-B	249/259 (96%)	243 (98%)	6 (2%)	43	47
1	2-A	249/259 (96%)	232 (93%)	17 (7%)	14	10
1	2-B	249/259 (96%)	243 (98%)	6 (2%)	43	47
All	All	996/1036 (96%)	950 (95%)	46 (5%)	24	22

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1-A	37	LEU
1	1-A	42	PHE
1	1-A	51	GLU
1	1-A	79	ILE
1	1-A	99	ASN
1	1-A	100	ILE
1	1-A	135	ILE
1	1-A	166	SER
1	1-A	218	THR
1	1-A	252	LEU
1	1-A	256	GLN
1	1-A	288	ILE
1	1-A	298	ASN
1	1-A	313	LEU
1	1-A	321	ASN
1	1-A	324	SER
1	1-A	349	ILE
1	1-B	99	ASN
1	1-B	100	ILE
1	1-B	258	ILE
1	1-B	269	THR
1	1-B	272	ASN
1	1-B	284	ASP
1	2-A	37	LEU
1	2-A	42	PHE
1	2-A	51	GLU
1	2-A	79	ILE
1	2-A	99	ASN
1	2-A	100	ILE
1	2-A	135	ILE
1	2-A	166	SER
1	2-A	218	THR
1	2-A	252	LEU
1	2-A	256	GLN
1	2-A	288	ILE
1	2-A	298	ASN
1	2-A	313	LEU
1	2-A	321	ASN
1	2-A	324	SER
1	2-A	349	ILE
1	2-B	99	ASN
1	2-B	100	ILE
1	2-B	258	ILE

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Mol	Chain	Res	Type
1	2-B	269	THR
1	2-B	272	ASN
1	2-B	284	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	1-A	50	HIS
1	1-A	71	GLN
1	1-A	119	HIS
1	1-A	142	GLN
1	1-A	242	GLN
1	1-A	259	GLN
1	1-A	298	ASN
1	1-A	330	HIS
1	1-A	332	GLN
1	1-B	71	GLN
1	1-B	108	GLN
1	1-B	174	GLN
1	1-B	208	ASN
1	1-B	259	GLN
1	1-B	272	ASN
1	1-B	298	ASN
1	1-B	332	GLN
1	2-A	50	HIS
1	2-A	71	GLN
1	2-A	119	HIS
1	2-A	142	GLN
1	2-A	242	GLN
1	2-A	259	GLN
1	2-A	298	ASN
1	2-A	330	HIS
1	2-A	332	GLN
1	2-B	43	ASN
1	2-B	71	GLN
1	2-B	142	GLN
1	2-B	206	ASN
1	2-B	235	GLN
1	2-B	259	GLN
1	2-B	272	ASN
1	2-B	298	ASN
1	2-B	321	ASN

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Mol	Chain	Res	Type
1	2-B	328	GLN
1	2-B	332	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 oligosaccharides 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	1-B	2352	-	4,4,4	1.73	0	6,6,6	0.47	0
2	PO4	2-A	1352	-	4,4,4	1.49	0	6,6,6	0.53	0
2	PO4	2-B	2352	-	4,4,4	1.76	0	6,6,6	0.51	0
2	PO4	1-A	1352	-	4,4,4	1.49	0	6,6,6	0.53	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

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.