



wwPDB X-ray Structure Validation Summary Report i

Feb 27, 2014 – 11:15 PM GMT

PDB ID : 1C47
Title : BINDING DRIVEN STRUCTURAL CHANGES IN CRYSTALLINE PHOSPHOGLUCOMUTASE ASSOCIATED WITH CHEMICAL REACTION
Authors : Baranidharan, S.; Ray Jr., W.J.
Deposited on : 1999-08-11
Resolution : 2.70 Å(reported)

This is a wwPDB 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/ValidationPDFNotes.html>

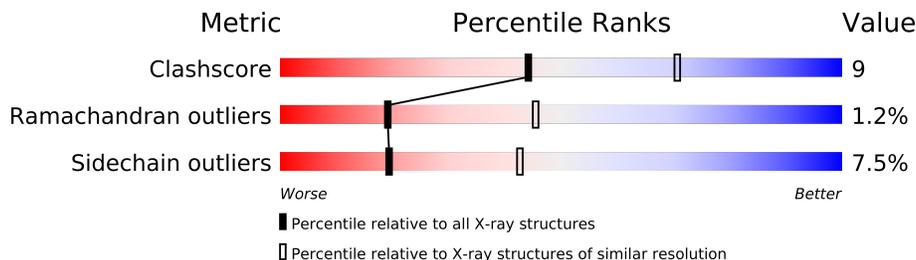
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	561	
1	B	561	

2 Entry composition (i)

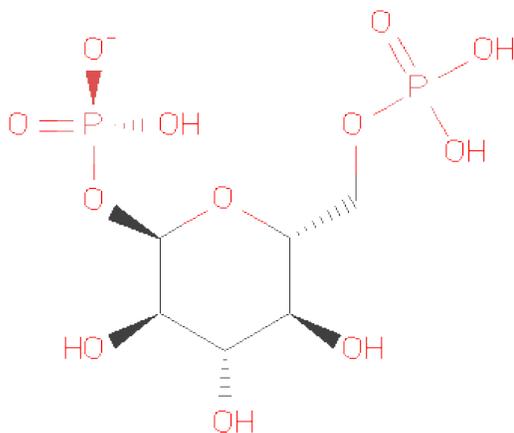
There are 4 unique types of molecules in this entry. The entry contains 8828 atoms, of which 0 are hydrogen and 0 are deuterium.

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 ALPHA-D-GLUCOSE 1,6-BISPHOSPHATE PHOSPHOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	561	4329	2753	743	817	16	0	0	0
1	B	561	4329	2753	743	817	16	0	0	0

- Molecule 2 is SUGAR (ALPHA-D-GLUCOSE 1,6-BISPHOSPHATE) (three-letter code: G16) (formula: $C_6H_{13}O_{12}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
2	A	1	20	6	12	2	0	0

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cd 1 1	0	0
3	A	1	Total Cd 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	59	Total O 59 59	0	0
4	B	89	Total O 89 89	0	0

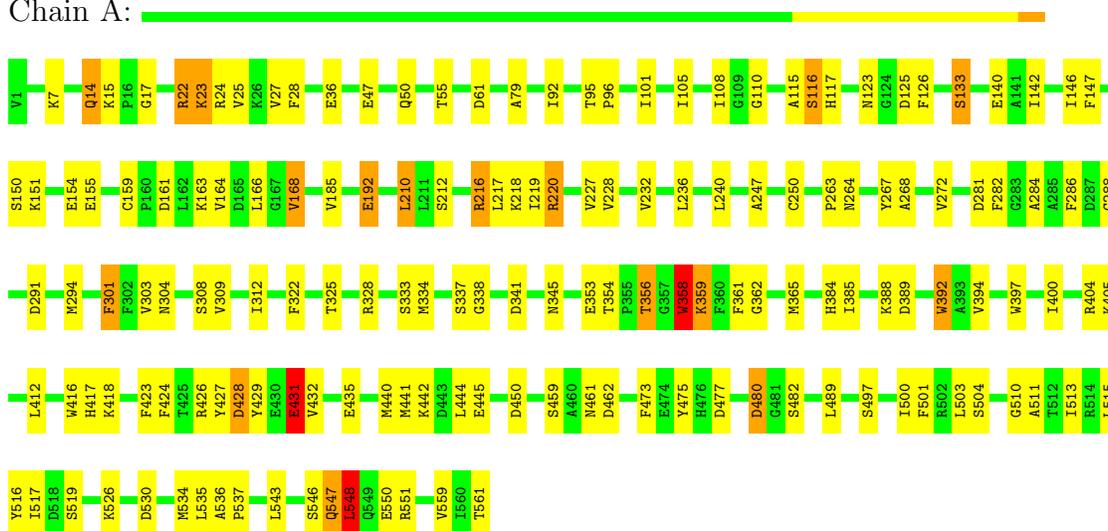
3 Residue-property plots

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.

Note EDS was not executed.

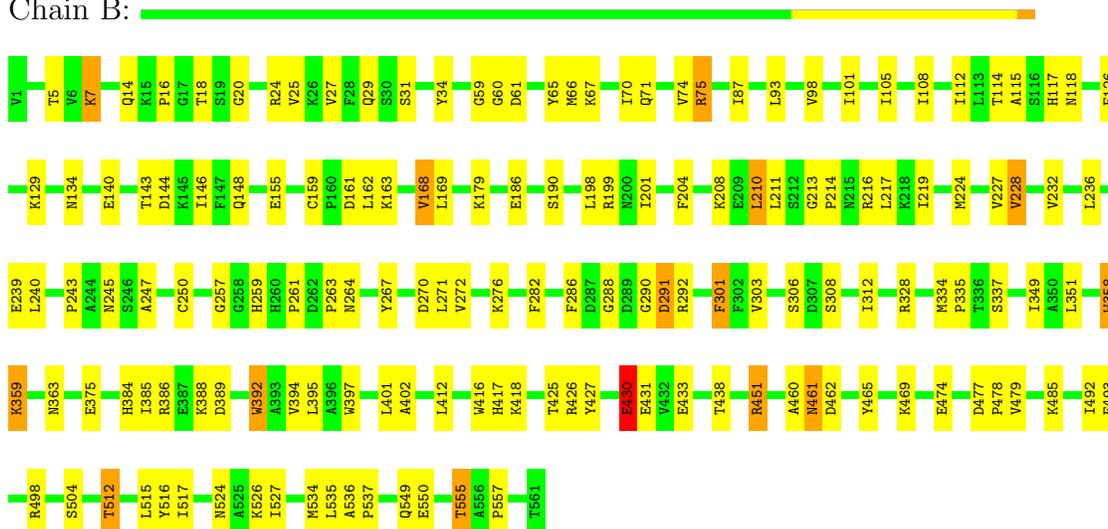
- Molecule 1: ALPHA-D-GLUCOSE 1,6-BISPHOSPHATE PHOSPHOTRANSFERASE

Chain A:



- Molecule 1: ALPHA-D-GLUCOSE 1,6-BISPHOSPHATE PHOSPHOTRANSFERASE

Chain B:



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	174.42Å 174.42Å 101.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.70	Depositor
% Data completeness (in resolution range)	96.5 (6.00-2.70)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.183 , 0.248	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	8828	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: G16, CD

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	1.47	15/4416 (0.3%)	0.94	12/5969 (0.2%)
1	B	1.54	21/4416 (0.5%)	1.00	17/5969 (0.3%)
All	All	1.51	36/8832 (0.4%)	0.97	29/11938 (0.2%)

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	550	GLU	CD-OE2	-8.27	1.16	1.25
1	B	308	SER	CB-OG	-8.17	1.31	1.42
1	A	36	GLU	CD-OE1	-7.92	1.17	1.25
1	B	312	ILE	C-O	-7.23	1.09	1.23
1	B	186	GLU	CG-CD	-7.22	1.41	1.51

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	392	TRP	CD1-CG-CD2	14.56	117.95	106.30
1	B	416	TRP	CD1-CG-CD2	14.15	117.62	106.30
1	B	397	TRP	CD1-CG-CD2	13.73	117.29	106.30
1	B	358	TRP	CD1-CG-CD2	13.63	117.21	106.30
1	A	416	TRP	CD1-CG-CD2	13.57	117.15	106.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4329	0	4332	71	1
1	B	4329	0	4332	80	0
2	A	20	0	9	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	59	0	0	2	1
4	B	89	0	0	7	0
All	All	8828	0	8673	150	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

The worst 5 of 150 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:428:ASP:HB2	1:A:559:VAL:HB	1.51	0.92
1:B:224:MET:SD	1:B:286:PHE:O	2.44	0.76
1:A:17:GLY:HA3	4:A:616:HOH:O	1.87	0.75
1:B:31:SER:HB2	1:B:34:TYR:HB2	1.71	0.72
1:A:547:GLN:HA	1:A:547:GLN:HE21	1.54	0.72

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	Distance(Å)	Clash(Å)
1:A:345:ASN:ND2	4:A:569:HOH:O[3_555]	2.19	0.01

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	559/561 (100%)	517 (92%)	34 (6%)	8 (1%)	16 41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	559/561 (100%)	517 (92%)	37 (7%)	5 (1%)	25	55
All	All	1118/1122 (100%)	1034 (92%)	71 (6%)	13 (1%)	19	45

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	430	GLU
1	B	461	ASN
1	A	358	TRP
1	B	301	PHE
1	A	192	GLU

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	462/462 (100%)	420 (91%)	42 (9%)	14	30
1	B	462/462 (100%)	435 (94%)	27 (6%)	28	57
All	All	924/924 (100%)	855 (92%)	69 (8%)	19	43

5 of 69 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	459	SER
1	A	547	GLN
1	B	418	LYS
1	A	462	ASP
1	A	504	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	422	ASN
1	A	547	GLN
1	B	37	ASN

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Mol	Chain	Res	Type
1	A	345	ASN
1	B	29	GLN

5.3.3 RNA [i](#)

There are no RNA chains 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](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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$
2	G16	A	563	3	20,20,20	1.69	4 (20%)	31,31,31	1.50	5 (16%)

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	G16	A	563	3	-	0/11/31/31	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	563	G16	P'-O1	-4.20	1.49	1.60
2	A	563	G16	O4-C4	-3.64	1.34	1.43
2	A	563	G16	O6-C6	-2.74	1.33	1.44
2	A	563	G16	C4-C5	2.12	1.57	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	563	G16	O5-C5-C6	3.43	113.51	106.61
2	A	563	G16	O3-C3-C2	-3.25	103.06	110.35
2	A	563	G16	O1-C1-C2	-2.90	103.07	108.38
2	A	563	G16	C1-O5-C5	2.67	118.91	113.73
2	A	563	G16	O1X-P'-O1	2.03	111.03	104.93

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.