



wwPDB NMR Structure Validation Summary Report ⓘ

May 7, 2024 – 03:48 pm BST

PDB ID : 2XDF
Title : Solution Structure of the Enzyme I Dimer Complexed with HPr Using Residual Dipolar Couplings and Small Angle X-Ray Scattering
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Deposited on : 2010-04-30

This is a wwPDB NMR 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

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

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	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

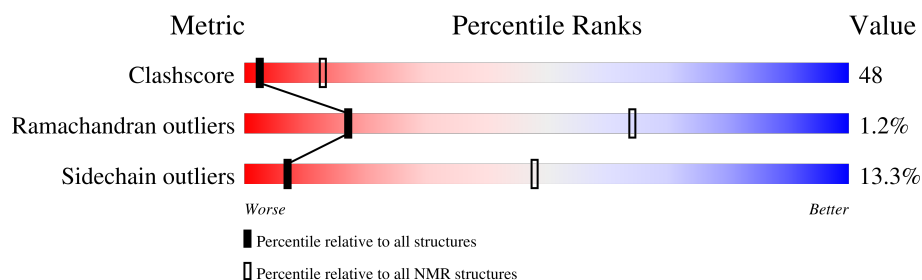
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR, SOLUTION SCATTERING

The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	573	63% 24% 13%
1	B	573	62% 26% 12%
2	C	85	81% 16% ..
2	D	85	82% 15% ..

2 Ensemble composition and analysis ⓘ

This entry contains 2 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 20502 atoms, of which 10338 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called PHOSPHOENOLPYRUVATE-PROTEIN PHOSPHOTRANSFERASE.

Mol	Chain	Residues	Atoms						Trace
1	A	573	Total	C	H	N	O	S	0
			8956	2790	4514	757	874	21	
1	B	573	Total	C	H	N	O	S	0
			8956	2790	4514	757	874	21	

- Molecule 2 is a protein called PHOSPHOCARRIER PROTEIN HPR.

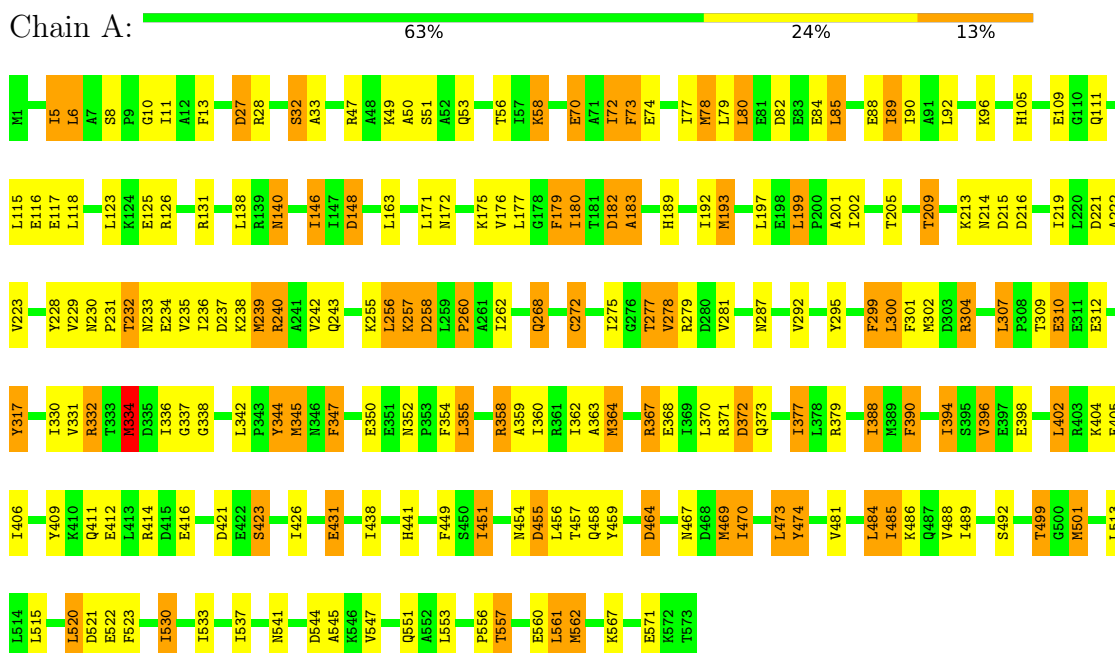
Mol	Chain	Residues	Atoms						Trace
2	C	85	Total	C	H	N	O	S	0
			1295	401	655	107	130	2	
2	D	85	Total	C	H	N	O	S	0
			1295	401	655	107	130	2	

4 Residue-property plots

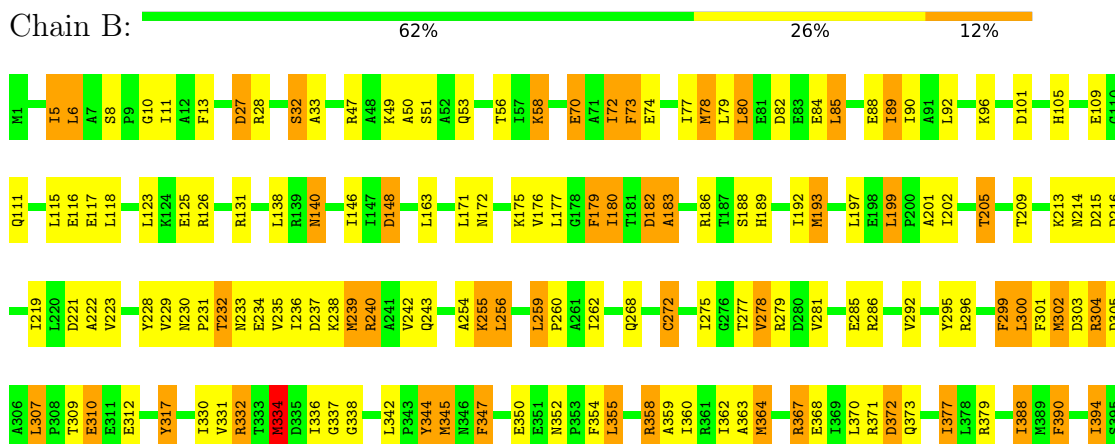
4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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 outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

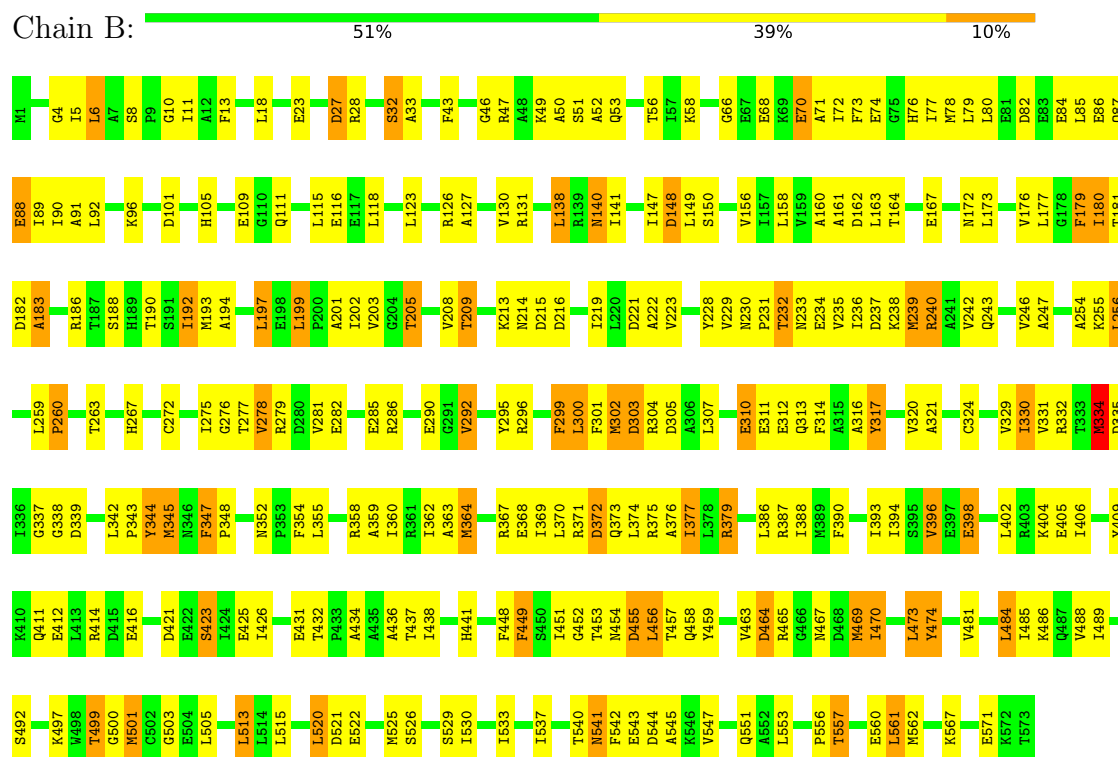
• Molecule 1: PHOSPHOENOLPYRUVATE-PROTEIN PHOSPHOTRANSFERASE



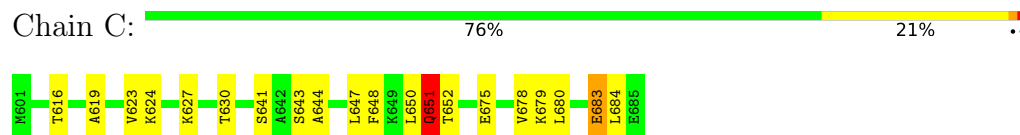
• Molecule 1: PHOSPHOENOLPYRUVATE-PROTEIN PHOSPHOTRANSFERASE



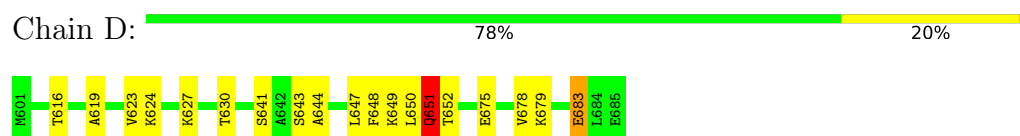
• Molecule 1: PHOSPHOENOLPYRUVATE-PROTEIN PHOSPHOTRANSFERASE



• Molecule 2: PHOSPHOCARRIER PROTEIN HPR



• Molecule 2: PHOSPHOCARRIER PROTEIN HPR



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: *simulated annealing*.

Of the 120 calculated structures, 2 were deposited, based on the following criterion: *BEST EXPERIMENT FIT, AND THEN LOWEST ENERGY*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
Xplor-NIH	refinement	2.25
Xplor-NIH	structure solution	

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

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 (average) 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	1.04±0.00	10±0/4495 (0.2± 0.0%)	0.88±0.00	4±0/6061 (0.1± 0.0%)
1	B	1.04±0.00	10±0/4495 (0.2± 0.0%)	0.88±0.00	4±0/6061 (0.1± 0.0%)
2	C	1.12±0.00	0±0/647 (0.0± 0.0%)	0.82±0.00	0±0/871 (0.0± 0.0%)
2	D	1.12±0.00	0±0/647 (0.0± 0.0%)	0.82±0.00	0±0/871 (0.0± 0.0%)
All	All	1.05	40/20568 (0.2%)	0.87	16/27728 (0.1%)

5 of 20 unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	562	MET	CG-SD	7.10	1.99	1.81	2	2
1	B	562	MET	CG-SD	7.08	1.99	1.81	1	2
1	A	345	MET	SD-CE	6.56	2.14	1.77	2	2
1	B	345	MET	SD-CE	6.55	2.14	1.77	1	2
1	A	469	MET	SD-CE	6.46	2.13	1.77	2	2

5 of 8 unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	561	LEU	CA-CB-CG	10.52	139.50	115.30	1	2
1	B	561	LEU	CA-CB-CG	10.52	139.49	115.30	1	2
1	A	300	LEU	CA-CB-CG	5.33	127.55	115.30	1	2
1	B	300	LEU	CA-CB-CG	5.33	127.55	115.30	1	2
1	A	414	ARG	NE-CZ-NH2	-5.31	117.64	120.30	1	2

There are no chirality outliers.

There are no planarity outliers.

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	4442	4514	4506	488±6
1	B	4442	4514	4506	487±4
2	C	640	655	650	38±5
2	D	640	655	650	38±4
All	All	20328	20676	20624	1953

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 48.

5 of 1730 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:126:ARG:HH22	2:C:619:ALA:CB	1.61	1.04	1	1
1:B:126:ARG:HH22	2:D:619:ALA:CB	1.55	1.04	1	1
1:B:314:PHE:CE2	1:B:376:ALA:HA	1.49	1.41	1	1
1:A:314:PHE:CE2	1:A:376:ALA:HA	1.48	1.40	1	1
1:B:562:MET:SD	1:B:562:MET:CE	1.48	2.01	2	1

6.3 Torsion angles [i](#)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	571/573 (100%)	538±0 (94±0%)	26±0 (5±0%)	7±0 (1±0%)	17	64
1	B	571/573 (100%)	538±2 (94±0%)	26±2 (5±0%)	7±0 (1±0%)	17	64
2	C	83/85 (98%)	76±0 (92±0%)	6±0 (7±0%)	1±0 (1±0%)	17	64
2	D	83/85 (98%)	76±0 (92±0%)	6±0 (7±0%)	1±0 (1±0%)	17	64
All	All	2616/2632 (99%)	2455 (94%)	129 (5%)	32 (1%)	17	64

5 of 19 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	6	LEU	2
1	A	148	ASP	2
1	A	183	ALA	2
1	A	232	THR	2
1	A	260	PRO	2

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	474/474 (100%)	409±4 (86±1%)	65±4 (14±1%)	7	47
1	B	474/474 (100%)	408±4 (86±1%)	66±4 (14±1%)	7	47
2	C	70/70 (100%)	63±2 (90±3%)	7±2 (10±3%)	11	56
2	D	70/70 (100%)	64±2 (91±2%)	6±2 (9±2%)	12	59
All	All	2176/2176 (100%)	1887 (87%)	289 (13%)	7	48

5 of 228 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	27	ASP	2
1	A	28	ARG	2
1	A	32	SER	2
1	A	58	LYS	2
1	A	80	LEU	2

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided