wwPDB X-ray Structure Validation Summary Report (i)

Mar 4, 2024 – 10:32 PM EST

PDB ID	:	2CXO
Title	:	Crystal structure of mouse AMF / E4P complex
Authors	:	Tanaka, N.; Haga, A.; Naba, N.; Shiraiwa, K.; Kusakabe, Y.; Hashimoto, K.; Funasaka, T.; Nagase, H.; Raz, A.; Nakamura, K.T.
Deposited on	:	2005-06-30
Resolution	:	1.80 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

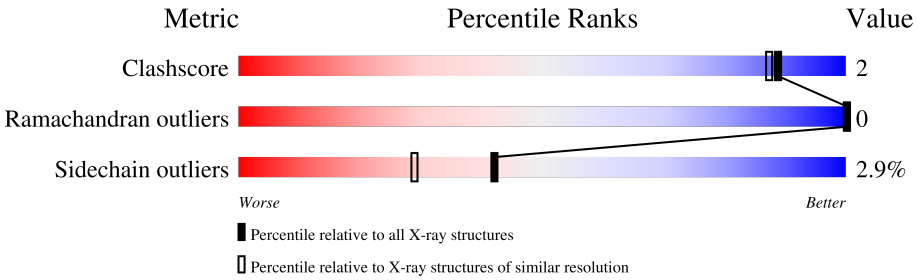
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	557	93% 7%
1	B	557	93% 7%

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	DER	A	1601	X	-	-	-
2	DER	B	2601	X	-	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 9899 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.

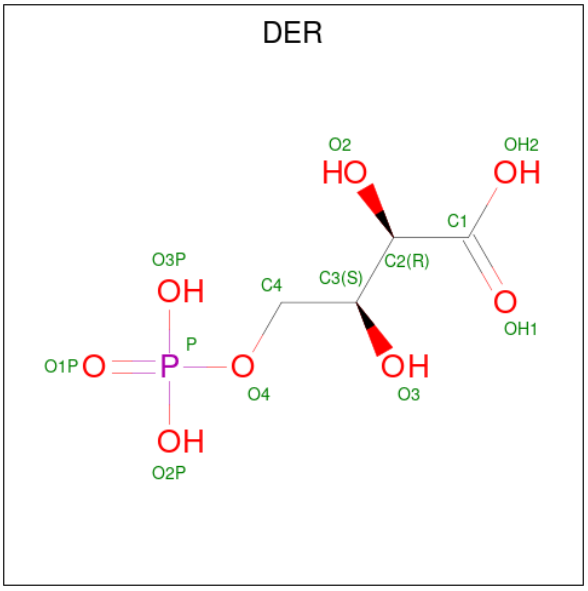
- Molecule 1 is a protein called Glucose-6-phosphate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4416	2826	764	807	19			
1	B	557	Total	C	N	O	S	0	0	0
			4416	2826	764	807	19			

There are 4 discrepancies between the modelled and reference sequences:

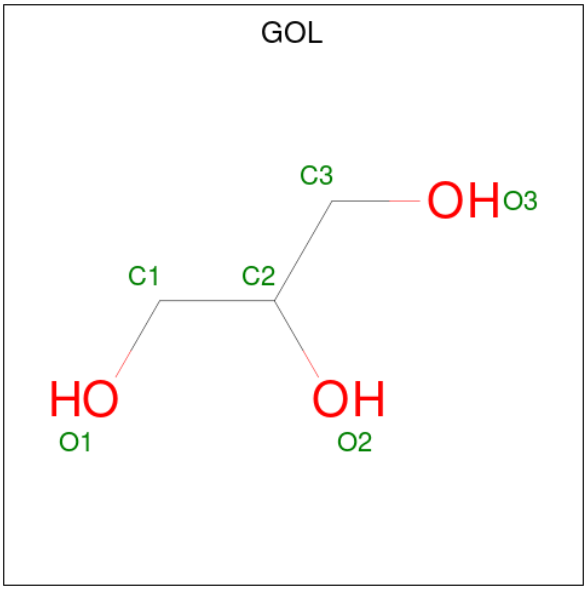
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P06745
A	61	SER	ASN	conflict	UNP P06745
B	1	MET	-	initiating methionine	UNP P06745
B	61	SER	ASN	conflict	UNP P06745

- Molecule 2 is D-4-PHOSPHOERYTHRONIC ACID (three-letter code: DER) (formula: $C_4H_9O_8P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O P 13 4 8 1	0	0
2	B	1	Total C O P 13 4 8 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	513	Total O 513 513	0	0
4	B	504	Total O 504 504	0	0

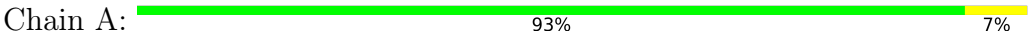
3 Residue-property plots (i)

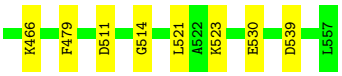
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: Glucose-6-phosphate isomerase

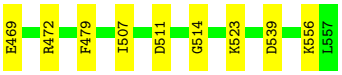
Chain A:





- Molecule 1: Glucose-6-phosphate isomerase

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	69.63Å 115.86Å 73.32Å 90.00° 101.64° 90.00°	Depositor
Resolution (Å)	40.00 – 1.80	Depositor
% Data completeness (in resolution range)	100.0 (40.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.159 , 0.189	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9899	wwPDB-VP
Average B, all atoms (Å^2)	16.0	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: DER, GOL

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	0.39	0/4525	0.64	6/6124 (0.1%)
1	B	0.39	0/4525	0.62	3/6124 (0.0%)
All	All	0.39	0/9050	0.63	9/12248 (0.1%)

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	511	ASP	CB-CG-OD2	6.49	124.14	118.30
1	A	188	ASP	CB-CG-OD2	6.39	124.05	118.30
1	A	259	ASP	CB-CG-OD2	6.35	124.02	118.30
1	B	511	ASP	CB-CG-OD2	5.75	123.47	118.30
1	B	188	ASP	CB-CG-OD2	5.70	123.43	118.30

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	A	4416	0	4398	15	0
1	B	4416	0	4398	23	0
2	A	13	0	7	0	0
2	B	13	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	12	0	16	0	0
3	B	12	0	16	0	0
4	A	513	0	0	3	0
4	B	504	0	0	2	0
All	All	9899	0	8842	34	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 2.

The worst 5 of 34 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:393:GLN:HE22	1:B:514:GLY:H	1.23	0.85
1:A:514:GLY:H	1:B:393:GLN:HE22	1.26	0.83
1:A:521:LEU:O	4:A:2060:HOH:O	2.04	0.76
1:A:388:GLN:HE22	1:A:428:ASN:HB3	1.55	0.71
1:B:390:ALA:C	4:B:3105:HOH:O	2.27	0.71

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	555/557 (100%)	542 (98%)	13 (2%)	0	100	100
1	B	555/557 (100%)	543 (98%)	12 (2%)	0	100	100
All	All	1110/1114 (100%)	1085 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	475/475 (100%)	461 (97%)	14 (3%)	42 29
1	B	475/475 (100%)	461 (97%)	14 (3%)	42 29
All	All	950/950 (100%)	922 (97%)	28 (3%)	42 29

5 of 28 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	104	ARG
1	B	556	LYS
1	B	213	PHE
1	B	466	LYS
1	B	198	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	397	GLN
1	B	475	ASN
1	B	465	HIS
1	B	47	ASN
1	B	393	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

6 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	$\#\|Z\| > 2$
3	GOL	B	1703	-	5,5,5	0.33	0	5,5,5	0.36	0
3	GOL	A	1704	-	5,5,5	0.31	0	5,5,5	0.30	0
3	GOL	A	1701	-	5,5,5	0.35	0	5,5,5	0.38	0
3	GOL	B	1702	-	5,5,5	0.32	0	5,5,5	0.40	0
2	DER	A	1601	-	11,12,12	2.39	2 (18%)	13,17,17	2.36	2 (15%)
2	DER	B	2601	-	11,12,12	2.43	2 (18%)	13,17,17	2.24	2 (15%)

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	GOL	B	1703	-	-	0/4/4/4	-
3	GOL	A	1704	-	-	0/4/4/4	-
3	GOL	A	1701	-	-	0/4/4/4	-
3	GOL	B	1702	-	-	2/4/4/4	-
2	DER	A	1601	-	1/1/4/4	3/14/14/14	-
2	DER	B	2601	-	1/1/4/4	3/14/14/14	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1601	DER	OH1-C1	6.84	1.43	1.22
2	B	2601	DER	OH1-C1	6.83	1.43	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2601	DER	OH2-C1	3.67	1.42	1.30
2	A	1601	DER	OH2-C1	3.41	1.41	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1601	DER	OH2-C1-OH1	-6.63	109.03	124.09
2	B	2601	DER	OH2-C1-OH1	-6.04	110.38	124.09
2	B	2601	DER	OH1-C1-C2	-4.09	110.87	121.63
2	A	1601	DER	OH1-C1-C2	-3.81	111.62	121.63

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1601	DER	C3
2	B	2601	DER	C3

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2601	DER	OH1-C1-C2-O2
2	A	1601	DER	OH1-C1-C2-O2
2	B	2601	DER	OH1-C1-C2-C3
2	A	1601	DER	OH1-C1-C2-C3
2	A	1601	DER	OH2-C1-C2-O2

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.