

wwPDB X-ray Structure Validation Summary Report (i)

Nov 6, 2023 – 12:14 AM EST

PDB ID : 1PKN

Title : STRUCTURE OF RABBIT MUSCLE PYRUVATE KINASE COMPLEXED

WITH MN2+, K+, AND PYRUVATE

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Deposited on : 1994-03-25

Resolution : 2.90 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)

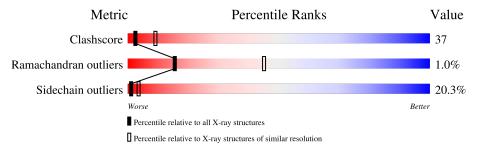
Validation Pipeline (wwPDB-VP) : 2.36

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 2.90 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}(\mathring{\rm A})) \end{array}$
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)

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	Λ	530	32%	44%	18%			



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3937 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 PYRUVATE KINASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Λ	514	Total	С	N	О	S	0	0	0
1	A	514	3929	2469	698	734	28	0	U	

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	149	ALA	GLU	conflict	UNP P11974
A	150	ALA	LYS	conflict	UNP P11974
A	168	GLU	ASP	conflict	UNP P11974
A	233	ASP	GLU	conflict	UNP P11974
A	234	GLU	GLN	conflict	UNP P11974
A	499	ALA	ARG	conflict	UNP P11974

• Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

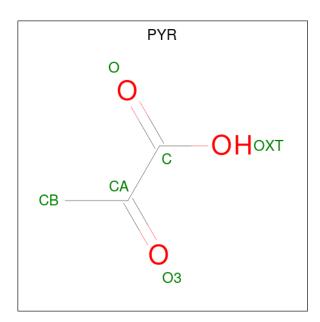
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total K 1 1	0	0

• Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mn 1 1	0	0

• Molecule 4 is PYRUVIC ACID (three-letter code: PYR) (formula: C₃H₄O₃).





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

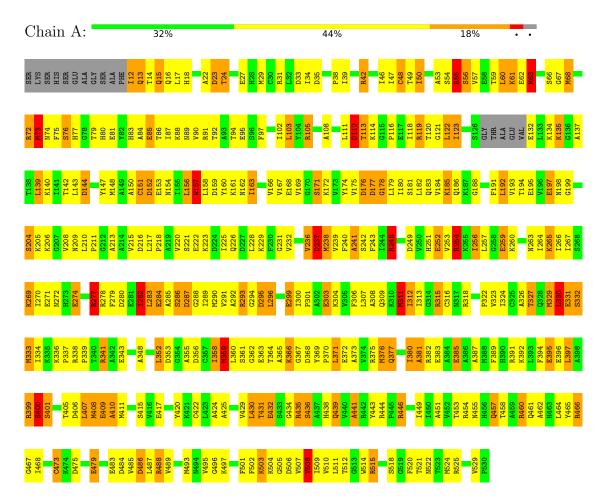


3 Residue-property plots (i)

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: PYRUVATE KINASE





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	Depositor	
Cell constants	83.60Å 109.90Å 146.80Å	Depositor	
a, b, c, α , β , γ	94.90° 93.60° 112.30°	Depositor	
Resolution (Å)	30.00 - 2.90	Depositor	
% Data completeness	(Not available) (30.00-2.90)	Depositor	
(in resolution range)	, , , , , , , , , , , , , , , , , , , ,		
R_{merge}	(Not available)	Depositor	
R_{sym}	(Not available)	Depositor	
Refinement program	TNT	Depositor	
R, R_{free}	0.191 , (Not available)	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3937	wwPDB-VP	
Average B, all atoms (Å ²)	37.0	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: MN, PYR, K

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).

Mal	Chain	Bo	nd lengths	Bond angles		
MIOI	Mol Chain		# Z > 5	RMSZ $\# Z > 5$		
1	A	1.02	19/3991~(0.5%)	1.92	$147/5384 \ (2.7\%)$	

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\text{\AA})$
1	A	432	GLU	CD-OE1	8.72	1.35	1.25
1	A	259	GLU	CD-OE1	8.47	1.34	1.25
1	A	331	GLU	CD-OE1	8.29	1.34	1.25
1	A	95	GLU	CD-OE1	7.50	1.33	1.25
1	A	479	GLU	CD-OE1	7.38	1.33	1.25

The worst 5 of 147 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$Ideal(^{o})$
1	A	369	TYR	CB-CG-CD1	-17.72	110.37	121.00
1	A	55	ARG	NE-CZ-NH2	-16.21	112.20	120.30
1	A	369	TYR	CB-CG-CD2	16.16	130.69	121.00
1	A	119	ARG	NE-CZ-NH1	15.81	128.21	120.30
1	A	55	ARG	NE-CZ-NH1	12.89	126.75	120.30

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	3929	0	4001	294	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	6	0	0	3	0
All	All	3937	0	4001	294	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 37.

The worst 5 of 294 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:279:PHE:CE2	1:A:283:LEU:HD12	1.65	1.29
1:A:122:LEU:O	1:A:151:CYS:HB2	1.29	1.26
1:A:134:LYS:NZ	1:A:139:LEU:HD13	1.56	1.20
1:A:139:LEU:HD12	1:A:153:GLU:O	1.44	1.16
1:A:72:ARG:HD2	1:A:359:MET:HE3	1.32	1.07

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	510/530 (96%)	460 (90%)	45 (9%)	5 (1%)	15 45

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	367	GLY
1	A	178	GLY
1	A	259	GLU
1	A	327	THR

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Mol	Chain	Res	Type
1	A	67	GLY

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	419/431 (97%)	334 (80%)	85 (20%)	1 3

5 of 85 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	332	SER
1	A	415	SER
1	A	352	LEU
1	A	392	LYS
1	A	439	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	463	HIS
1	A	478	GLN
1	A	494	ASN
1	A	226	GLN
1	A	377	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)

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	B	ond leng	$_{ m gths}$	В	ond ang	gles
WIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PYR	A	533	3	5,5,5	2.15	2 (40%)	3,6,6	4.48	2 (66%)

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
4	PYR	A	533	3	-	0/4/4/4	-

All (2) bond length outliers are listed below:

\mathbf{Mol}	Chain	Res	Type	Atoms	${f Z}$	Observed(A)	$\operatorname{Ideal}(ext{\AA})$
4	A	533	PYR	O-C	3.92	1.33	1.22
4	A	533	PYR	OXT-C	-2.04	1.24	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
4	A	533	PYR	O3-CA-CB	5.68	132.33	119.73
4	A	533	PYR	OXT-C-O	5.09	135.25	123.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.



1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	533	PYR	3	0

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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

