

Full wwPDB X-ray Structure Validation Report (i)

Feb 15, 2024 – 08:49 AM EST

PDB ID : 3PNP

Title : THE HIGH RESOLUTION CRYSTAL STRUCTURE OF BOVINE SPLEEN

PURINE NUCLEOSIDE PHOSPHORYLASE IN COMPLEX FORMS WITH

PHOSPHATE AND 9-DEAZAINOSINE

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Deposited on : 1998-03-27

Resolution : 1.60 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 (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 (1)) were used in the production of this report:

Mol Probity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

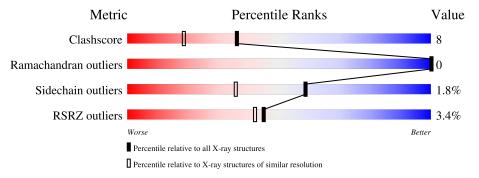
Validation Pipeline (wwPDB-VP) : 2.36

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 1.60 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	Similar resolution
Wiediic	$(\# {\rm Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain				
			3%				
1	A	289	77%	13%	9%		



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2172 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 PURINE NUCLEOSIDE PHOSPHORYLASE.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	263	Total	С	N	О	S	0	2	0
1	Α	203	2067	1312	364	374	17	0	3	

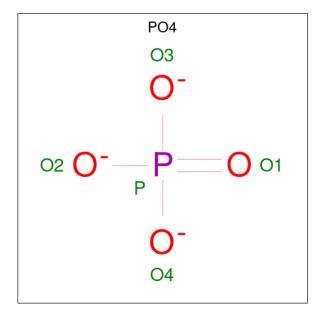
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	GLN	conflict	UNP P55859
A	249	THR	TYR	conflict	UNP P55859

• Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

• Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).





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

• Molecule 4 is water.

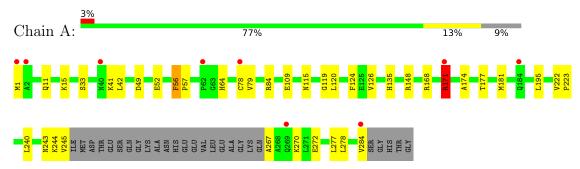
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	99	Total O 99 99	0	0



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 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.

• Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants	93.00Å 93.00Å 93.00Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 - 1.60	Depositor
Resolution (A)	14.70 - 1.55	EDS
% Data completeness	96.0 (8.00-1.60)	Depositor
(in resolution range)	84.3 (14.70-1.55)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) > 1$	4.09 (at 1.55Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.208 , 0.232	Depositor
it, it free	0.203 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	14.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.46,65.3	EDS
L-test for twinning ²	$< L >=0.49, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.032 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2172	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.27% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: PO4, MG

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	Bond	lengths	Bond angles		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.38	0/2130	0.68	$2/2882 \ (0.1\%)$	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	A	171[A]	ARG	NE-CZ-NH2	8.40	124.50	120.30
1	A	171[B]	ARG	NE-CZ-NH2	8.40	124.50	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	171[A]	ARG	Sidechain
1	A	171[B]	ARG	Sidechain

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	2067	0	2038	33	0
2	A	1	0	0	0	0
3	A	5	0	0	0	0
4	A	99	0	0	1	0
All	All	2172	0	2038	33	0

The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is 8.

All (33) 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:171[B]:ARG:HB2	1:A:171[B]:ARG:NH1	1.53	1.23
1:A:171[B]:ARG:NH1	1:A:171[B]:ARG:CB	2.29	0.94
1:A:171[B]:ARG:CG	1:A:171[B]:ARG:HH11	1.81	0.94
1:A:171[B]:ARG:CB	1:A:171[B]:ARG:HH11	1.85	0.90
1:A:1:MET:H2	1:A:148:ARG:HD3	1.37	0.89
1:A:171[B]:ARG:HB2	1:A:171[B]:ARG:HH11	1.37	0.84
1:A:42:LEU:HD21	1:A:78[B]:CYS:SG	2.21	0.80
1:A:171[B]:ARG:HH11	1:A:171[B]:ARG:HG3	1.44	0.80
1:A:64:HIS:HD2	1:A:84:ARG:H	1.31	0.74
1:A:120:LEU:HD23	1:A:245:VAL:HG21	1.70	0.72
1:A:126:VAL:HG22	1:A:243:ASN:HA	1.76	0.66
1:A:171[B]:ARG:NH1	1:A:171[B]:ARG:CG	2.49	0.61
1:A:42:LEU:CD2	1:A:78[B]:CYS:SG	2.88	0.60
1:A:11:GLN:O	1:A:15:LYS:HG2	2.01	0.60
1:A:33:SER:OG	1:A:64:HIS:HE1	1.89	0.56
1:A:49:ASP:O	1:A:52:GLU:HG2	2.06	0.56
1:A:78[C]:CYS:SG	1:A:79:VAL:N	2.81	0.54
1:A:177:THR:O	1:A:181:MET:HG3	2.09	0.51
1:A:1:MET:N	4:A:380:HOH:O	2.45	0.50
1:A:174:ALA:HA	1:A:278:LEU:HD21	1.95	0.47
1:A:181:MET:HE3	1:A:277:LEU:HD12	1.96	0.47
1:A:119:GLY:O	1:A:245:VAL:HG22	2.15	0.46
1:A:171[B]:ARG:HB2	1:A:171[B]:ARG:CZ	2.36	0.45
1:A:240:LEU:HD11	1:A:267:ALA:HB3	1.97	0.45
1:A:41:LYS:HE3	1:A:272:GLU:HB3	1.99	0.45
1:A:171[B]:ARG:CB	1:A:171[B]:ARG:CZ	2.93	0.45
1:A:124:PHE:O	1:A:244:LYS:HE3	2.19	0.42
1:A:270:LYS:HA	1:A:270:LYS:HD2	1.90	0.42

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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	$\operatorname{distance}\ (ext{Å})$	overlap (Å)
1:A:168:ARG:O	1:A:171[B]:ARG:HB3	2.20	0.42
1:A:1:MET:N	1:A:148:ARG:HD3	2.19	0.41
1:A:222:VAL:N	1:A:223:PRO:HD2	2.36	0.41
1:A:56:PHE:HA	1:A:57:PRO:HD3	1.91	0.41
1:A:109:GLU:OE1	1:A:284:VAL:HG21	2.21	0.40

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	$263/289 \ (91\%)$	257 (98%)	6 (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.

M	ol	Chain	Analysed	Rotameric	Outliers	Percentiles
1		A	222/238 (93%)	218 (98%)	4 (2%)	59 36

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	PHE

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Mol	Chain	Res	Type
1	A	115	ASN
1	A	135	HIS
1	A	195	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	115	ASN
1	A	211	ASN

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 2 ligands modelled in this entry, 1 is 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	Pog	Link	\mathbf{B}	ond leng	$_{ m gths}$	В	ond ang	${ m gles}$
WIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	PO4	A	290	-	4,4,4	2.07	1 (25%)	6,6,6	0.42	0



All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
3	A	290	PO4	P-O1	3.82	1.59	1.50

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 (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	$\langle { m RSRZ} \rangle$	# RSRZ > 2	$OWAB(A^2)$	Q<0.9
1	A	263/289 (91%)	0.02	9 (3%) 45 42	8, 14, 36, 63	6 (2%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	9.5
1	A	284	VAL	6.6
1	A	62	PRO	4.3
1	A	269	GLN	3.8
1	A	78[A]	CYS	3.0
1	A	40	ASN	2.7
1	A	171[A]	ARG	2.1
1	A	2	ALA	2.1
1	A	184	GLN	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.



Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	MG	A	291	1/1	0.99	0.04	17,17,17,17	0
3	PO4	A	290	5/5	0.99	0.06	9,9,10,10	0

6.5 Other polymers (i)

There are no such residues in this entry.

