

wwPDB X-ray Structure Validation Summary Report (i)

Oct 15, 2023 – 02:28 AM EDT

PDB ID : 8D2M

Title : Covalent Schiff base complex of YedK C2A and abasic DNA

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Deposited on : 2022-05-30

Resolution : 1.82 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 (1)) were used in the production of this report:

MolProbity : 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) roteins) : Engh & Huber (2001)

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

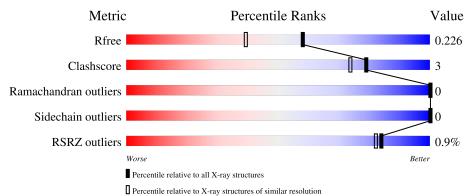
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 1.82 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar \ resolution} \\ (\#{\rm Entries, \ resolution \ range(\AA)}) \end{array}$		
R_{free}	130704	7484 (1.84-1.80)		
Clashscore	141614	8401 (1.84-1.80)		
Ramachandran outliers	138981	8290 (1.84-1.80)		
Sidechain outliers	138945	8290 (1.84-1.80)		
RSRZ outliers	127900	7371 (1.84-1.80)		

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% 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		
1	A	227	93%		6% •
1	В	227	91%		7% •
2	С	7	71%	299	%
2	D	7	71%	14%	14%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7418 atoms, of which 3490 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 Abasic site processing protein YedK.

Mol	Chain	Residues		Atoms				ZeroOcc	AltConf	Trace	
1	A	224	Total 3420	C 1126	H 1667	N 304	O 318	S 5	2	2	0
1	В	222	20000	C 1121		N 305	O 316	S 5	2	3	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	CYS	engineered mutation	UNP P76318
A	223	LEU	-	expression tag	UNP P76318
A	224	GLU	-	expression tag	UNP P76318
A	225	VAL	-	expression tag	UNP P76318
A	226	LEU	-	expression tag	UNP P76318
A	227	PHE	-	expression tag	UNP P76318
A	228	GLN	-	expression tag	UNP P76318
В	2	ALA	CYS	engineered mutation	UNP P76318
В	223	LEU	-	expression tag	UNP P76318
В	224	GLU	-	expression tag	UNP P76318
В	225	VAL	-	expression tag	UNP P76318
В	226	LEU	-	expression tag	UNP P76318
В	227	PHE	-	expression tag	UNP P76318
В	228	GLN	-	expression tag	UNP P76318

• Molecule 2 is a DNA chain called DNA (5'-D(*GP*TP*CP*(PED)P*GP*GP*A)-3').

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace		
2	С	7	Total	С	Н	N	О	Р	0	0	0
		1	211	64	77	25	39	6	U	U	
2	D	7	Total	С	Н	N	О	Р	0	0	0
	ש	1	211	64	77	25	39	6	U		U

• Molecule 3 is water.



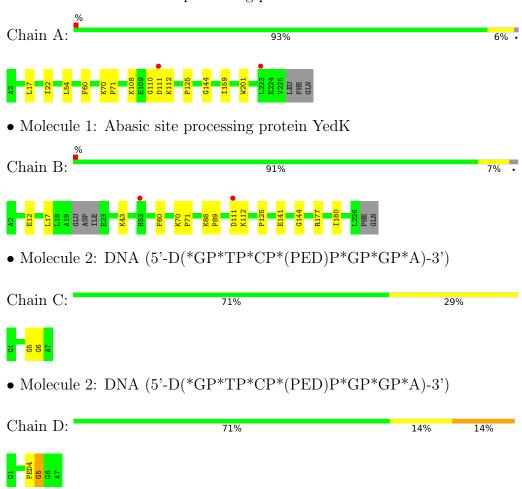
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	70	Total O 70 70	0	0
3	С	6	Total O 6 6	0	0
3	В	77	Total O 77 77	0	0
3	D	7	Total O 7 7	0	0



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: Abasic site processing protein YedK





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	60.94Å 41.47Å 82.58Å	Donositor
a, b, c, α , β , γ	90.00° 95.65° 90.00°	Depositor
Resolution (Å)	37.02 - 1.82	Depositor
rtesolution (A)	37.03 - 1.82	EDS
% Data completeness	99.5 (37.02-1.82)	Depositor
(in resolution range)	94.1 (37.03-1.82)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) > 1$	2.73 (at 1.82Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
P. P.	0.178 , 0.226	Depositor
R, R_{free}	0.178 , 0.226	DCC
R_{free} test set	1852 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.43, 44.6	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7418	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.10% 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: PED

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		
Mol	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.54	0/1815	0.64	0/2472	
1	В	0.62	2/1817 (0.1%)	0.69	0/2473	
2	С	0.98	0/137	1.01	0/208	
2	D	1.28	0/137	1.18	1/208~(0.5%)	
All	All	0.64	2/3906 (0.1%)	0.71	1/5361 (0.0%)	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	${f Z}$	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$\operatorname{Ideal}(ext{\AA})$
1	В	12[A]	GLU	CB-CG	5.57	1.62	1.52
1	В	12[B]	GLU	CB-CG	5.57	1.62	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	D	5	DG	O4'-C1'-N9	7.31	113.12	108.00

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	1753	1667	1669	9	0

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Continued	trom	mmoninonic	maaa
COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	В	1747	1669	1664	9	0
2	С	134	77	78	1	0
2	D	134	77	78	1	0
3	A	70	0	0	0	0
3	В	77	0	0	0	0
3	С	6	0	0	0	0
3	D	7	0	0	0	0
All	All	3928	3490	3489	20	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 3.

The worst 5 of 20 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:B:43:LYS:HE2	1:B:141:GLU:OE1	1.97	0.62
1:A:22:ILE:HG23	1:A:54:LEU:CD1	2.36	0.55
1:A:70:LYS:H	1:A:70:LYS:HD2	1.73	0.54
1:B:111:ASP:OD2	1:B:112:LYS:HG2	2.08	0.52
1:A:70:LYS:HB3	1:A:71:PRO:CD	2.42	0.49

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	Perce	$_{ m ntiles}$
1	A	$224/227 \ (99\%)$	217 (97%)	7 (3%)	0	100	100
1	В	$221/227 \ (97\%)$	216 (98%)	5 (2%)	0	100	100
All	All	445/454 (98%)	433 (97%)	12 (3%)	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	Percen	ntiles
1	A	174/184 (95%)	174 (100%)	0	100	100
1	В	176/184 (96%)	176 (100%)	0	100	100
All	All	350/368 (95%)	350 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	198	GLN
1	В	55	HIS

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)

There are no ligands in this entry.

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	<rsrz></rsrz>	# RSRZ > 2	$OWAB(A^2)$	Q<0.9
1	A	224/227 (98%)	-0.06	2 (0%) 84 82	15, 25, 52, 60	0
1	В	222/227 (97%)	-0.16	2 (0%) 84 82	13, 23, 41, 58	0
2	С	6/7 (85%)	-0.36	0 100 100	27, 28, 39, 40	0
2	D	6/7 (85%)	-0.32	0 100 100	23, 26, 32, 32	0
All	All	458/468 (97%)	-0.11	4 (0%) 84 82	13, 24, 47, 60	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	LEU	3.7
1	В	53	HIS	2.8
1	A	111	ASP	2.7
1	В	111	ASP	2.7

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)

There are no ligands in this entry.



6.5 Other polymers (i)

There are no such residues in this entry.

