wwPDB X-ray Structure Validation Summary Report (i)

Feb 3, 2024 – 11:46 PM EST

PDB ID	:	1PYI
Title	:	CRYSTAL STRUCTURE OF A PPR1-DNA COMPLEX: DNA RECOGNITION BY PROTEINS CONTAINING A ZN2CYS6 BINUCLEAR CLUSTER
Authors	:	Marmorstein, R.; Harrison, S.C.
Deposited on	:	1995-01-04
Resolution	:	3.20 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:

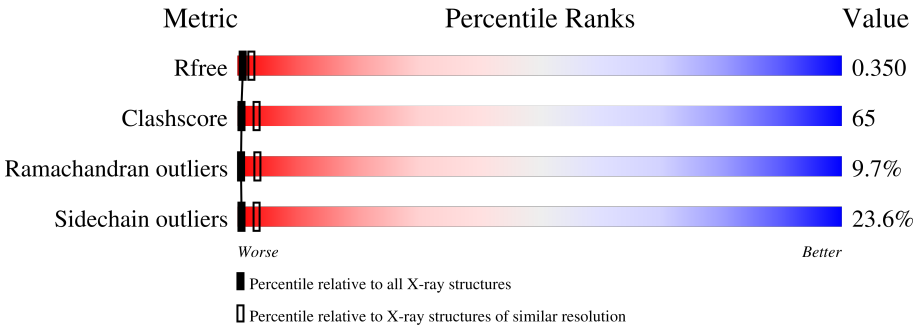
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 3.20 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))
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)

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%

Mol	Chain	Length	Quality of chain
1	D	14	21% 64% 14%
1	E	14	7% 14% 79%
2	A	96	15% 45% 21% 11% 8%
2	B	96	20% 23% 26% 27%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 1826 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 DNA chain called DNA (5'-D(*TP*CP*GP*GP*CP*AP*AP*TP*TP*GP*CP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	14	Total	C	N	O	P	0	0	0
			284	136	53	82	13			
1	E	14	Total	C	N	O	P	0	0	0
			284	136	53	82	13			

- Molecule 2 is a protein called PROTEIN (PYRIMIDINE PATHWAY REGULATOR 1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	88	Total	C	N	O	S	0	0	0
			697	438	124	127	8			
2	B	70	Total	C	N	O	S	0	0	0
			556	350	100	98	8			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	1	Total	O	0	0
			1	1		

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.

• Molecule 1: DNA (5'-D(*TP*CP*GP*GP*CP*AP*AP*TP*TP*GP*CP*CP*GP*A)-3')

Chain D:

• Molecule 1: DNA (5'-D(*TP*CP*GP*GP*CP*AP*AP*TP*TP*GP*CP*CP*GP*A)-3')

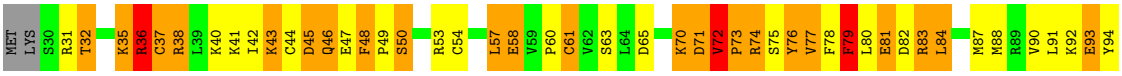
Chain E:

• Molecule 2: PROTEIN (PYRIMIDINE PATHWAY REGULATOR 1)

Chain A:

• Molecule 2: PROTEIN (PYRIMIDINE PATHWAY REGULATOR 1)

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	53.80Å 92.60Å 174.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 3.20 10.04 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-3.20) 98.3 (10.04-3.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.91 (at 2.99Å)	Xtriage
Refinement program		Depositor
R, R_{free}	0.245 , 0.330 0.286 , 0.350	Depositor DCC
R_{free} test set	830 reflections (9.61%)	wwPDB-VP
Wilson B-factor (Å^2)	63.2	Xtriage
Anisotropy	0.480	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.24 , 86.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.87	EDS
Total number of atoms	1826	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 6.99% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	D	1.98	10/318 (3.1%)	3.47	41/489 (8.4%)
1	E	1.58	3/318 (0.9%)	2.65	27/489 (5.5%)
2	A	1.11	5/709 (0.7%)	2.12	33/952 (3.5%)
2	B	1.25	5/565 (0.9%)	2.17	27/757 (3.6%)
All	All	1.41	23/1910 (1.2%)	2.53	128/2687 (4.8%)

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 mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	2
2	A	0	2
All	All	0	4

The worst 5 of 23 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	73	PRO	N-CD	9.44	1.61	1.47
1	D	9	DT	C1'-N1	9.20	1.61	1.49
1	D	9	DT	C4-C5	-8.62	1.37	1.45
1	E	23	DT	C5-C7	8.61	1.55	1.50
1	D	9	DT	C2-N3	7.68	1.43	1.37

The worst 5 of 128 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	9	DT	C4-C5-C7	-28.72	101.77	119.00
1	D	9	DT	C6-C5-C7	27.81	139.59	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	9	DT	C5-C4-O4	-22.06	109.45	124.90
1	D	9	DT	N3-C4-O4	15.63	129.28	119.90
1	D	14	DA	O4'-C1'-N9	13.77	117.64	108.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	104	ASN	Sidechain
2	A	111	ASP	Sidechain
1	D	3	DG	Sidechain
1	D	9	DT	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	D	284	0	159	25	0
1	E	284	0	159	32	0
2	A	697	0	714	112	0
2	B	556	0	575	96	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	E	1	0	0	0	0
All	All	1826	0	1607	223	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 65.

The worst 5 of 223 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)
2:A:113:PRO:HB2	2:A:116:LEU:HD11	1.29	1.14
2:A:97:ASP:HB2	2:A:98:PRO:HD2	1.31	1.05
1:D:1:DT:H6	1:D:1:DT:HO5'	1.05	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:101:ILE:HG21	2:B:83:ARG:HE	1.32	0.95
1:E:21:DA:H2"	1:E:22:DT:H5"	1.48	0.93

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
2	A	86/96 (90%)	59 (69%)	16 (19%)	11 (13%)	0 1
2	B	68/96 (71%)	54 (79%)	10 (15%)	4 (6%)	1 12
All	All	154/192 (80%)	113 (73%)	26 (17%)	15 (10%)	0 3

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	46	GLN
2	A	97	ASP
2	A	113	PRO
2	B	97	ASP
2	A	67	ALA

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
2	A	80/88 (91%)	61 (76%)	19 (24%)	0 3
2	B	64/88 (73%)	49 (77%)	15 (23%)	1 3
All	All	144/176 (82%)	110 (76%)	34 (24%)	1 3

5 of 34 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	76	TYR
2	B	77	VAL
2	B	93	GLU
2	A	98	PRO
2	A	96	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	46	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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.