



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2018 – 10:12 AM EDT

PDB ID : 6D0M
Title : Polymerase Eta post-insertion binary complex with cytarabine (AraC)
Authors : Rechkoblit, O.; Aggarwal, A.K.
Deposited on : 2018-04-10
Resolution : 1.83 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : rb-20031172
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20031172

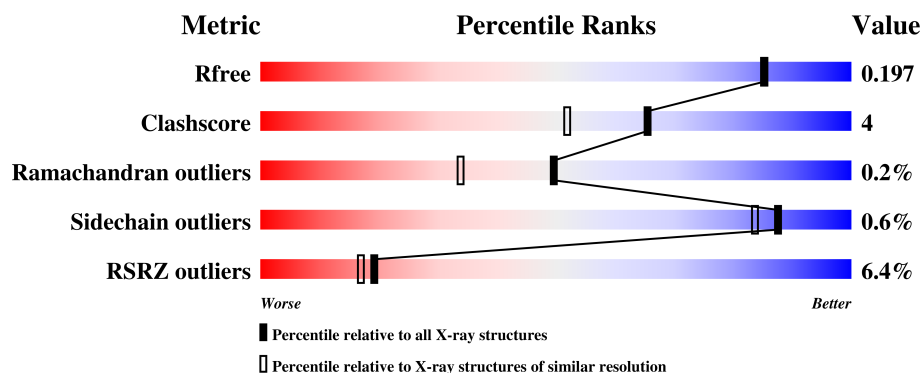
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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}	111664	3313 (1.86-1.82)
Clashscore	122126	3530 (1.86-1.82)
Ramachandran outliers	120053	3495 (1.86-1.82)
Sidechain outliers	120020	3496 (1.86-1.82)
RSRZ outliers	108989	3265 (1.86-1.82)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	435	<div> <div>7%</div> <div>90%</div> <div>8% ..</div> </div>
2	T	12	<div> <div>92%</div> <div>8%</div> </div>
3	P	9	<div> <div>67%</div> <div>33%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7956 atoms, of which 3670 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 DNA polymerase eta.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	430	Total	C	H	N	O	S	0	21	0
			6851	2163	3419	607	632	30			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9Y253
A	-1	PRO	-	expression tag	UNP Q9Y253
A	0	HIS	-	expression tag	UNP Q9Y253
A	406	MET	CYS	engineered mutation	UNP Q9Y253

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*AP*TP*GP*AP*CP*AP*GP*TP*GP*CP*T)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	T	12	Total	C	H	N	O	P	0	0	1
			353	108	125	42	67	11			

- Molecule 3 is a DNA chain called DNA/RNA (5'-D(*AP*GP*CP*AP*CP*TP*GP*T)-R(P*(CAR))-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	P	9	Total	C	H	N	O	P	0	2	0
			318	97	118	33	61	9			

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			14	3	8	3		

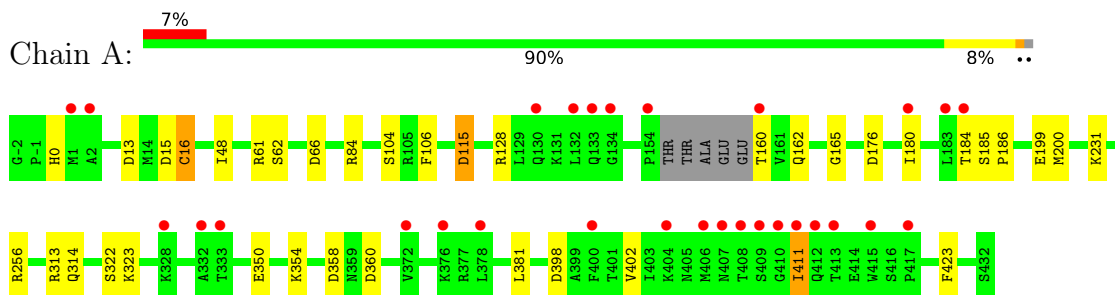
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	355	Total	O	0	8
			363	363		
5	T	34	Total	O	0	0
			34	34		
5	P	23	Total	O	0	0
			23	23		

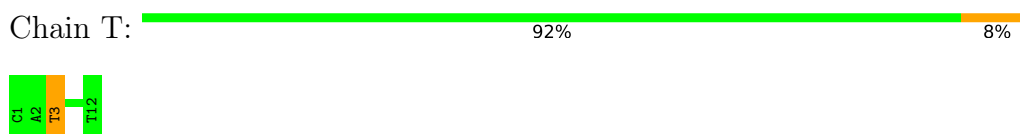
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

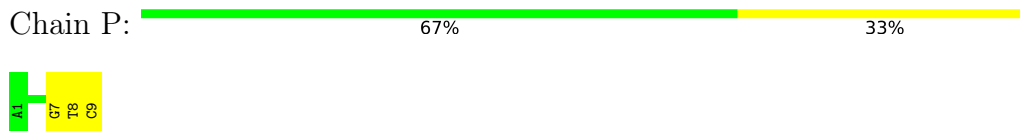
- Molecule 1: DNA polymerase eta



- Molecule 2: DNA (5'-D(*CP*AP*TP*GP*AP*CP*AP*GP*TP*GP*CP*T)-3')



- Molecule 3: DNA/RNA (5'-D(*AP*GP*CP*AP*CP*TP*GP*T)-R(P*(CAR))-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	98.57Å 98.57Å 81.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.68 – 1.83 42.68 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.68-1.83) 99.6 (42.68-1.83)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 1.83Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.160 , 0.196 0.162 , 0.197	Depositor DCC
R_{free} test set	1977 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 52.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.042 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7956	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

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: GOL, CAR

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	0.35	0/3541	0.52	0/4786
2	T	0.82	0/255	1.10	2/393 (0.5%)
3	P	0.84	0/226	1.03	1/348 (0.3%)
All	All	0.43	0/4022	0.62	3/5527 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	3	DT	O4'-C1'-N1	7.54	113.28	108.00
3	P	7	DG	O4'-C4'-C3'	-6.85	101.76	104.50
2	T	3	DT	C1'-O4'-C4'	-6.12	103.98	110.10

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	3432	3419	3472	31	1
2	T	228	125	125	1	0
3	P	200	118	110	4	0
4	A	6	8	8	0	0
5	A	363	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	P	23	0	0	1	0
5	T	34	0	0	0	0
All	All	4286	3670	3715	33	1

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 4.

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:256:ARG:NH1	5:A:601:HOH:O	1.96	0.97
1:A:199:GLU:OE1	5:A:602:HOH:O	2.07	0.72
1:A:66:ASP:OD1	5:A:603:HOH:O	2.08	0.71
3:P:9:CAR:OP1	5:P:101:HOH:O	2.12	0.67
1:A:128[B]:ARG:NH1	5:A:836[B]:HOH:O	2.26	0.67
1:A:160:THR:HG22	1:A:162:GLN:H	1.59	0.66
1:A:176:ASP:O	1:A:180[A]:ILE:HG13	1.99	0.61
1:A:381:LEU:HD23	1:A:381:LEU:C	2.22	0.59
1:A:61:ARG:HG2	2:T:3:DT:O4	2.06	0.56
1:A:48:ILE:O	3:P:9:CAR:O2'	2.23	0.55
1:A:256:ARG:NH2	5:A:611:HOH:O	2.41	0.54
1:A:313:ARG:NE	5:A:607:HOH:O	2.34	0.53
1:A:106:PHE:CG	1:A:200[B]:MET:HG2	2.42	0.53
1:A:104:SER:OG	1:A:314:GLN:OE1	2.22	0.53
1:A:185[A]:SER:OG	1:A:186:PRO:HD3	2.10	0.51
1:A:84:ARG:NH2	1:A:350:GLU:OE1	2.44	0.51
1:A:354:LYS:NZ	5:A:746[B]:HOH:O	2.36	0.51
1:A:354:LYS:NZ	1:A:358:ASP:OD2	2.42	0.48
1:A:160:THR:HG21	1:A:165:GLY:HA3	1.94	0.48
1:A:180[B]:ILE:HB	5:A:843:HOH:O	2.14	0.47
1:A:61:ARG:O	1:A:62:SER:CB	2.63	0.46
1:A:84:ARG:NE	5:A:614:HOH:O	2.43	0.45
1:A:398:ASP:O	1:A:402:VAL:HG13	2.19	0.43
1:A:115:ASP:OD1	3:P:9:CAR:H5'	2.18	0.42
1:A:322[B]:SER:HB2	1:A:423[B]:PHE:HD1	1.84	0.42
3:P:8:DT:H2'	3:P:9:CAR:C6	2.49	0.42
1:A:15:ASP:O	1:A:16:CYS:C	2.57	0.42
1:A:180[A]:ILE:HG22	1:A:180[A]:ILE:O	2.19	0.41
1:A:184[A]:THR:HB	1:A:186:PRO:HD2	2.02	0.41
1:A:411:ILE:H	1:A:411:ILE:HD12	1.86	0.41
1:A:13:ASP:OD2	1:A:231:LYS:NZ	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:ARG:NH1	5:A:629:HOH:O	2.55	0.40
1:A:323:LYS:NZ	5:A:616:HOH:O	2.43	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:0:HIS:HE2	1:A:360:ASP:OD2[5_454]	1.49	0.11

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	448/435 (103%)	437 (98%)	10 (2%)	1 (0%)	49 34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	CYS

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	378/372 (102%)	376 (100%)	2 (0%)	90 87

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	ASP
1	A	411	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	CAR	P	9	3,2	14,21,22	0.65	0	15,30,33	0.78	0

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	CAR	P	9	3,2	-	0/3/25/26	0/2/2/2

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.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	9	CAR	4	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
4	GOL	A	501	-	5,5,5	0.38	0	5,5,5	0.36	0

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	GOL	A	501	-	-	0/4/4/4	0/0/0/0

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

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, 95th 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>2	OWAB(Å ²)	Q<0.9
1	A	430/435 (98%)	-0.12	29 (6%) 18 16	9, 20, 53, 86	2 (0%)
2	T	12/12 (100%)	-0.23	0 100 100	18, 26, 56, 60	0
3	P	8/9 (88%)	-0.58	0 100 100	19, 27, 34, 35	0
All	All	450/456 (98%)	-0.13	29 (6%) 19 17	9, 20, 53, 86	2 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	132	LEU	4.6
1	A	406	MET	4.1
1	A	378	LEU	4.0
1	A	410	GLY	3.9
1	A	408	THR	3.8
1	A	134	GLY	3.8
1	A	412	GLN	3.6
1	A	376	LYS	3.2
1	A	409	SER	3.0
1	A	333	THR	3.0
1	A	180[A]	ILE	3.0
1	A	411	ILE	3.0
1	A	332	ALA	3.0
1	A	417	PRO	2.9
1	A	1	MET	2.9
1	A	413	THR	2.7
1	A	160	THR	2.6
1	A	133	GLN	2.6
1	A	372	VAL	2.6
1	A	130	GLN	2.5
1	A	400	PHE	2.5
1	A	415	TRP	2.5
1	A	2	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	407	ASN	2.4
1	A	154	PRO	2.3
1	A	404	LYS	2.3
1	A	183[A]	LEU	2.3
1	A	328	LYS	2.1
1	A	184[A]	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 95th 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	B-factors(Å ²)	Q<0.9
3	CAR	P	9	20/21	0.84	0.17	19,38,48,58	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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, 95th 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	B-factors(Å ²)	Q<0.9
4	GOL	A	501	6/6	0.96	0.12	13,16,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.