



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 5, 2016 – 01:35 PM EDT

PDB ID : 4ZRL  
Title : Structure of the non canonical Poly(A) polymerase complex GLD-2 - GLD-3  
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Deposited on : 2015-05-12  
Resolution : 2.28 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027790  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027790

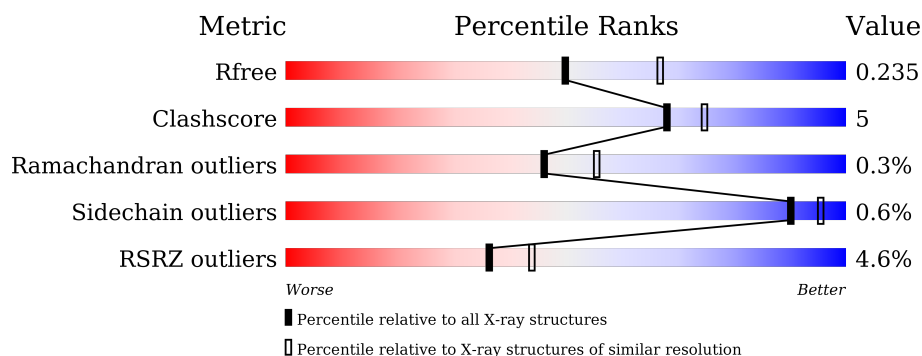
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.28 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))
$R_{free}$	91344	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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  $\geq 3$ , 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	364	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>10%</div> <div>14%</div> </div> </div>
2	B	77	<div> <div>9%</div> <div> <div></div> <div>78%</div> <div>5%</div> <div>17%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3081 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 Poly(A) RNA polymerase gld-2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	Se	0	0	0
			2458	1566	411	466	6	9			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	525	GLY	-	expression tag	UNP O17087
A	526	ALA	-	expression tag	UNP O17087
A	527	MSE	ARG	conflict	UNP O17087
A	?	-	ARG	deletion	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	LEU	deletion	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	VAL	deletion	UNP O17087
A	?	-	ARG	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	LYS	deletion	UNP O17087
A	?	-	ILE	deletion	UNP O17087
A	?	-	HIS	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	ASN	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	GLU	deletion	UNP O17087
A	?	-	GLY	deletion	UNP O17087
A	?	-	ASP	deletion	UNP O17087
A	?	-	LYS	deletion	UNP O17087
A	?	-	GLU	deletion	UNP O17087
A	?	-	THR	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	SER	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	THR	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	ILE	deletion	UNP O17087
A	?	-	HIS	deletion	UNP O17087
A	?	-	ASN	deletion	UNP O17087
A	?	-	GLY	deletion	UNP O17087

- Molecule 2 is a protein called Defective in germ line development protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	64	Total	C	N	O	Se	0	0	0
			507	308	94	103	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	12	MSE	-	initiating methionine	UNP Q95ZK7

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

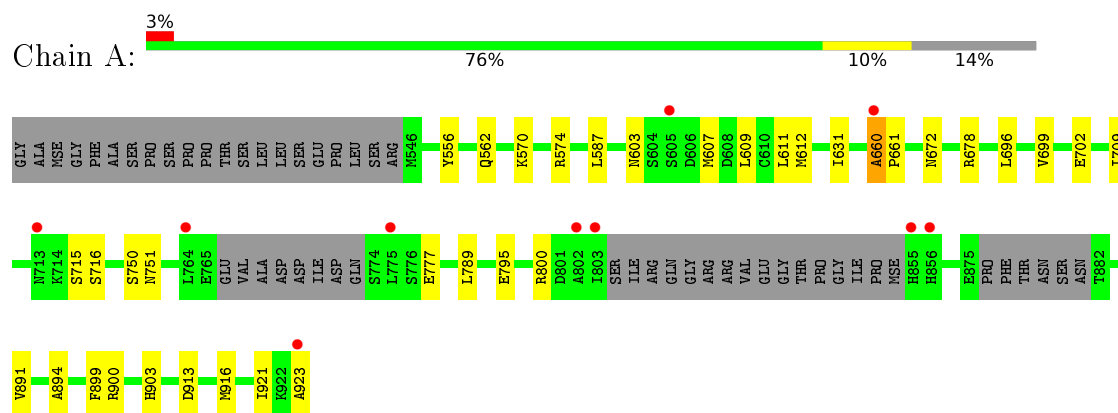
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	86	Total	O	0	0
			86	86		
4	B	29	Total	O	0	0
			29	29		

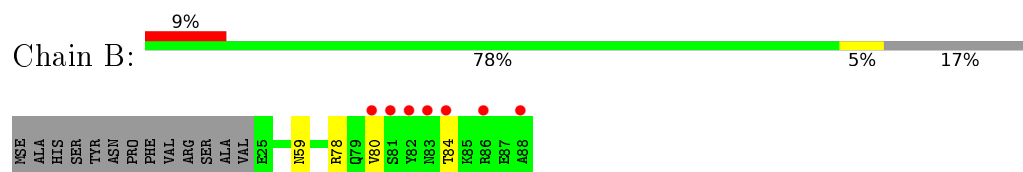
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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: Poly(A) RNA polymerase gld-2



- Molecule 2: Defective in germ line development protein 3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.66 Å 89.37 Å 94.67 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.30 – 2.28 47.33 – 2.28	Depositor EDS
% Data completeness (in resolution range)	98.4 (47.30-2.28) 98.4 (47.33-2.28)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.29 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1615)	Depositor
R, $R_{free}$	0.187 , 0.231 0.192 , 0.235	Depositor DCC
$R_{free}$ test set	987 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.4	Xtriage
Anisotropy	0.818	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3081	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

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

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	0.23	0/2502	0.41	0/3386
2	B	0.24	0/512	0.38	0/685
All	All	0.23	0/3014	0.40	0/4071

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	2458	0	2352	26	0
2	B	507	0	467	3	0
3	A	1	0	0	0	0
4	A	86	0	0	10	0
4	B	29	0	0	0	0
All	All	3081	0	2819	27	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 5.

All (27) 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:607:MSE:HE2	1:A:609:LEU:HD21	1.67	0.75
1:A:612:MSE:SE	1:A:672:ASN:HD22	2.22	0.72
1:A:699:VAL:HG13	1:A:916:MSE:HE1	1.72	0.71
1:A:562:GLN:O	2:B:78:ARG:NH2	2.23	0.71
1:A:899:PHE:O	4:A:1122:HOH:O	2.09	0.70
1:A:696:LEU:O	4:A:1123:HOH:O	2.10	0.68
1:A:562:GLN:NE2	4:A:1125:HOH:O	2.24	0.67
1:A:570:LYS:NZ	4:A:1126:HOH:O	2.29	0.65
1:A:660:ALA:HB1	1:A:661:PRO:HD2	1.84	0.59
1:A:900:ARG:NH2	4:A:1101:HOH:O	2.23	0.55
1:A:678:ARG:HD2	4:A:1121:HOH:O	2.05	0.55
1:A:903:HIS:N	4:A:1122:HOH:O	2.43	0.50
1:A:556:TYR:CE2	1:A:702:GLU:HG2	2.47	0.49
2:B:80:VAL:O	2:B:84:THR:HB	2.15	0.47
1:A:715:SER:O	1:A:891:VAL:HG11	2.14	0.46
1:A:750:SER:OG	1:A:751:ASN:N	2.47	0.46
1:A:789:LEU:HB3	4:A:1122:HOH:O	2.17	0.45
1:A:660:ALA:HB1	1:A:661:PRO:CD	2.46	0.44
1:A:587:LEU:HB3	2:B:59:ASN:HA	1.99	0.44
1:A:611:LEU:HD22	1:A:631:ILE:HD11	1.98	0.43
1:A:795:GLU:HG2	1:A:800:ARG:HD3	2.01	0.43
1:A:699:VAL:N	4:A:1123:HOH:O	2.52	0.43
1:A:913:ASP:OD1	4:A:1124:HOH:O	2.21	0.43
1:A:891:VAL:HG22	1:A:923:ALA:HB1	2.00	0.42
1:A:894:ALA:O	1:A:921:ILE:HD11	2.20	0.41
1:A:709:ILE:O	1:A:716:SER:HB2	2.20	0.41
1:A:574:ARG:HB2	1:A:607:MSE:HE1	2.03	0.41

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	305/364 (84%)	291 (95%)	13 (4%)	1 (0%)	46	55
2	B	62/77 (80%)	61 (98%)	1 (2%)	0	100	100
All	All	367/441 (83%)	352 (96%)	14 (4%)	1 (0%)	46	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	660	ALA

### 5.3.2 Protein sidechains ⓘ

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	270/319 (85%)	268 (99%)	2 (1%)	88	94
2	B	52/62 (84%)	52 (100%)	0	100	100
All	All	322/381 (84%)	320 (99%)	2 (1%)	90	95

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

Mol	Chain	Res	Type
1	A	603	ASN
1	A	777	GLU

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

Mol	Chain	Res	Type
1	A	672	ASN
1	A	743	GLN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 ⓘ

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	304/364 (83%)	0.04	10 (3%) 50 58	22, 39, 75, 98	0
2	B	62/77 (80%)	0.57	7 (11%) 7 9	27, 39, 97, 124	0
All	All	366/441 (82%)	0.13	17 (4%) 36 44	22, 39, 81, 124	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	82	TYR	8.3
1	A	855	HIS	5.9
1	A	923	ALA	5.5
2	B	88	ALA	4.3
2	B	83	ASN	3.3
1	A	803	ILE	3.1
1	A	660	ALA	3.1
2	B	80	VAL	2.9
1	A	605	SER	2.7
1	A	802	ALA	2.7
2	B	81	SER	2.6
2	B	86	ARG	2.6
2	B	84	THR	2.5
1	A	856	HIS	2.5
1	A	775	LEU	2.3
1	A	764	LEU	2.2
1	A	713	ASN	2.0

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

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

### 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	CL	A	1001	1/1	0.89	0.11	-0.18	55,55,55,55	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.