



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 08:45 am BST

PDB ID : 5WC8  
Title : Structure of a bacterial polysialyltransferase at 2.75 Angstrom resolution  
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Deposited on : 2017-06-29  
Resolution : 2.75 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](mailto: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.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

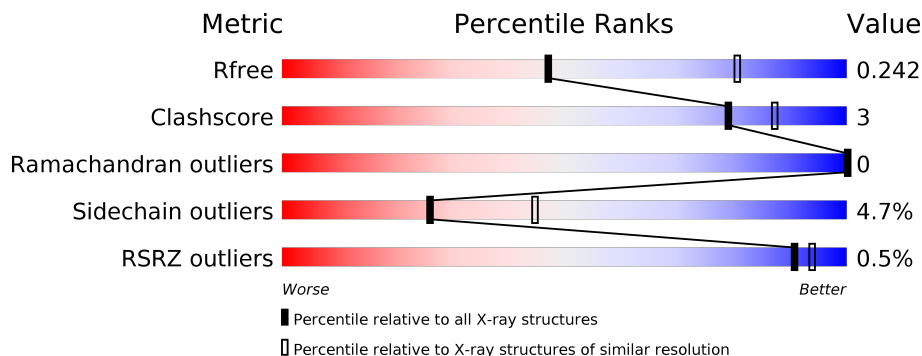
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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 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  $\leq 5\%$ . 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	382	
1	M	382	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BR	A	516	-	-	-	X
2	BR	A	517	-	-	-	X
2	BR	M	513	-	-	X	-
2	BR	M	516	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6451 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 SiaD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	M	381	3182	2075	523	569	15	0	0	0
1	A	381	3182	2075	523	569	15	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	20	ALA	-	expression tag	UNP G4RIN4
M	68	ALA	LYS	conflict	UNP G4RIN4
M	69	ALA	LYS	conflict	UNP G4RIN4
A	20	ALA	-	expression tag	UNP G4RIN4
A	68	ALA	LYS	conflict	UNP G4RIN4
A	69	ALA	LYS	conflict	UNP G4RIN4

- Molecule 2 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	20	Total	Br	0	0
			20	20		
2	M	16	Total	Br	0	0
			16	16		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	M	18	Total	O	0	0
			18	18		
3	A	33	Total	O	0	0
			33	33		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.08Å 78.08Å 303.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.76 – 2.75 61.76 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (61.76-2.75) 99.9 (61.76-2.75)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.71 (at 2.77Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.190 , 0.242 0.197 , 0.242	Depositor DCC
$R_{free}$ test set	1431 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	60.8	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 32.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6451	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.65	0/3247	0.82	4/4353 (0.1%)
1	M	0.65	0/3247	0.81	5/4353 (0.1%)
All	All	0.65	0/6494	0.82	9/8706 (0.1%)

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	109	MET	CG-SD-CE	-6.17	90.33	100.20
1	A	26	LEU	CA-CB-CG	5.82	128.69	115.30
1	M	293	LYS	CB-CG-CD	5.53	125.99	111.60
1	A	123	LYS	CB-CG-CD	5.51	125.92	111.60
1	A	55	LEU	N-CA-C	5.46	125.74	111.00

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	3182	0	3308	20	0
1	M	3182	0	3308	18	0
2	A	20	0	0	5	0
2	M	16	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	33	0	0	3	0
3	M	18	0	0	1	0
All	All	6451	0	6616	37	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 37 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:118:LYS:NZ	2:A:501:BR:BR	2.28	1.22
1:M:118:LYS:NZ	2:M:501:BR:BR	2.51	0.99
2:M:509:BR:BR	3:M:609:HOH:O	2.42	0.91
1:M:391:GLY:HA2	2:M:516:BR:BR	2.50	0.65
1:M:183:LYS:HG3	2:M:513:BR:BR	2.52	0.65

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	379/382 (99%)	371 (98%)	8 (2%)	0	100	100
1	M	379/382 (99%)	372 (98%)	7 (2%)	0	100	100
All	All	758/764 (99%)	743 (98%)	15 (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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	361/361 (100%)	345 (96%)	16 (4%)	28	47
1	M	361/361 (100%)	343 (95%)	18 (5%)	24	42
All	All	722/722 (100%)	688 (95%)	34 (5%)	26	45

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

Mol	Chain	Res	Type
1	M	350	LYS
1	A	32	LYS
1	A	365	VAL
1	M	390	ASP
1	M	123	LYS

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

Mol	Chain	Res	Type
1	A	133	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 36 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 [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<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	381/382 (99%)	0.10	2 (0%) 91 94	36, 58, 90, 121	0
1	M	381/382 (99%)	0.15	2 (0%) 91 94	36, 60, 94, 153	0
All	All	762/764 (99%)	0.12	4 (0%) 91 94	36, 59, 92, 153	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	30	TYR	5.7
1	A	25	HIS	2.9
1	A	30	TYR	2.8
1	M	232	SER	2.3

### 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 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	BR	A	517	1/1	0.20	0.51	133,133,133,133	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	BR	A	518	1/1	0.44	0.30	80,80,80,80	1
2	BR	M	514	1/1	0.53	0.35	72,72,72,72	1
2	BR	M	516	1/1	0.64	0.52	57,57,57,57	1
2	BR	A	516	1/1	0.65	0.46	89,89,89,89	1
2	BR	A	512	1/1	0.73	0.27	82,82,82,82	1
2	BR	M	515	1/1	0.81	0.29	53,53,53,53	1
2	BR	A	513	1/1	0.83	0.26	99,99,99,99	1
2	BR	A	508	1/1	0.83	0.18	80,80,80,80	1
2	BR	A	519	1/1	0.83	0.34	96,96,96,96	1
2	BR	A	514	1/1	0.85	0.21	78,78,78,78	1
2	BR	M	509	1/1	0.85	0.76	73,73,73,73	1
2	BR	M	512	1/1	0.86	0.26	60,60,60,60	1
2	BR	A	510	1/1	0.88	0.17	63,63,63,63	1
2	BR	A	520	1/1	0.88	0.75	63,63,63,63	1
2	BR	A	515	1/1	0.88	0.43	60,60,60,60	1
2	BR	A	506	1/1	0.89	0.14	66,66,66,66	1
2	BR	M	513	1/1	0.90	0.23	76,76,76,76	1
2	BR	A	511	1/1	0.90	0.35	94,94,94,94	1
2	BR	M	511	1/1	0.90	0.15	73,73,73,73	1
2	BR	M	504	1/1	0.90	0.14	67,67,67,67	1
2	BR	M	510	1/1	0.92	0.34	95,95,95,95	1
2	BR	M	508	1/1	0.93	0.14	49,49,49,49	1
2	BR	A	509	1/1	0.94	0.17	73,73,73,73	1
2	BR	M	501	1/1	0.94	0.08	68,68,68,68	1
2	BR	A	504	1/1	0.94	0.29	62,62,62,62	1
2	BR	M	502	1/1	0.94	0.13	66,66,66,66	1
2	BR	A	503	1/1	0.95	0.09	53,53,53,53	1
2	BR	A	505	1/1	0.95	0.11	60,60,60,60	1
2	BR	A	501	1/1	0.95	0.07	59,59,59,59	1
2	BR	M	503	1/1	0.95	0.13	53,53,53,53	1
2	BR	M	507	1/1	0.96	0.18	57,57,57,57	1
2	BR	A	507	1/1	0.97	0.14	37,37,37,37	1
2	BR	M	505	1/1	0.97	0.28	59,59,59,59	1
2	BR	A	502	1/1	0.98	0.09	45,45,45,45	1
2	BR	M	506	1/1	0.99	0.22	57,57,57,57	1

## 6.5 Other polymers

There are no such residues in this entry.