

Full wwPDB X-ray Structure Validation Report (i)

Sep 11, 2023 – 02:23 PM EDT

PDB ID : 4MBO

Title : 1.65 Angstrom Crystal Structure of Serine-rich Repeat Adhesion Glycoprotein

(Srr1) from Streptococcus agalactiae

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Deposited on : 2013-08-19

Resolution : 1.65 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 (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 : 2.35.1

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)

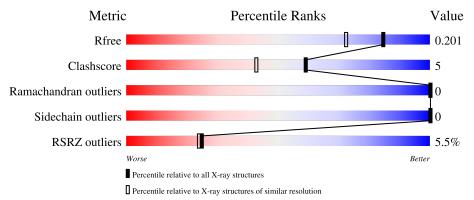
Validation Pipeline (wwPDB-VP) : 2.35.1

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 1.65 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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					
			5%					
1	A	340	82%	8% • 9%				



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3151 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 Serine-rich Repeat Adhesion Glycoprotein (Srr1).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	308	Total	С	N	О	S	0	29	0
1	A	300	2632	1634	441	549	8	0	29	

There are 9 discrepancies between the modelled and reference sequences:

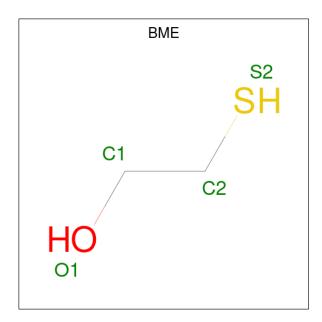
Chain	Residue	Modelled	Actual Comment		Reference
A	310	MET	- initiating methionine		UNP Q8E473
A	642	LEU	-	expression tag	UNP Q8E473
A	643	GLU	-	expression tag	UNP Q8E473
A	644	HIS	-	expression tag	UNP Q8E473
A	645	HIS	-	expression tag	UNP Q8E473
A	646	HIS	-	expression tag	UNP Q8E473
A	647	HIS	-	expression tag	UNP Q8E473
A	648	HIS	-	expression tag	UNP Q8E473
A	649	HIS	-	expression tag	UNP Q8E473

• Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total Ca 4 4	0	0

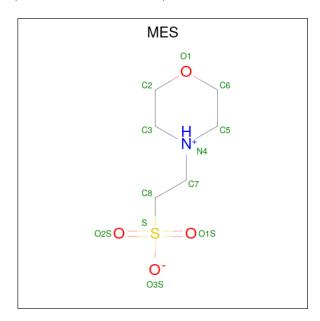
• Molecule 3 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C₂H₆OS).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 4	C 2	O 1	S 1	0	1

 \bullet Molecule 4 is 2-(N-MORPHOLINO)-ETHANE SULFONIC ACID (three-letter code: MES) (formula: $\rm C_6H_{13}NO_4S).$



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
4	A	1	Total 12	C 6	N 1	O 4	S 1	0	1

• Molecule 5 is water.



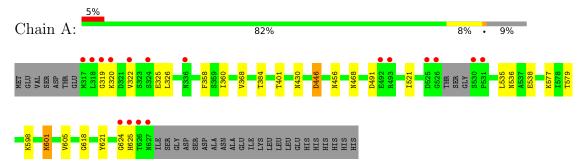
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	437	Total O 499 499	0	78



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: Serine-rich Repeat Adhesion Glycoprotein (Srr1)





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	73.36Å 83.23Å 57.06Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.93 - 1.65	Depositor
Resolution (A)	28.93 - 1.65	EDS
% Data completeness	99.5 (28.93-1.65)	Depositor
(in resolution range)	99.5 (28.93-1.65)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.50 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.8.0046	Depositor
D D.	0.154 , 0.187	Depositor
R, R_{free}	0.165 , 0.201	DCC
R_{free} test set	2141 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32, 52.2	EDS
L-test for twinning ²	$ < L > = 0.49, < L^2> = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3151	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.81% 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: CA, MES, MLZ, BME

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	Bond	lengths	Bo	nd angles
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.60	0/2659	0.79	1/3604 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
1	A	446	ASP	CB-CG-OD1	6.65	124.29	118.30

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	2632	0	2578	25	0
2	A	4	0	0	0	0
3	A	4	0	5	0	0
4	A	12	0	13	1	0
5	A	499	0	0	7	0
All	All	3151	0	2596	25	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 (25) 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:446:ASP:HB2	5:A:1077[A]:HOH:O	1.95	0.65
1:A:322:VAL:HG22	1:A:322:VAL:O	2.00	0.61
1:A:618:GLY:O	4:A:706[A]:MES:H81	2.03	0.58
1:A:326:LEU:HD23	1:A:360:ILE:HD12	1.85	0.58
1:A:430[A]:ASN:ND2	5:A:1006:HOH:O	2.39	0.54
1:A:319:GLY:O	1:A:320:LYS:HB2	2.08	0.53
1:A:468[B]:ASN:HB3	1:A:621[B]:TYR:OH	2.08	0.53
1:A:535[B]:LEU:O	1:A:536[B]:ASN:HB2	2.12	0.49
1:A:430[B]:ASN:O	1:A:430[B]:ASN:ND2	2.46	0.48
1:A:322:VAL:HG12	1:A:368:VAL:O	2.13	0.48
1:A:320:LYS:HE3	5:A:968:HOH:O	2.14	0.48
1:A:456:ASN:ND2	5:A:1112[A]:HOH:O	2.31	0.47
1:A:521:ILE:CD1	1:A:579:THR:HG23	2.44	0.47
1:A:601:MLZ:HE2	1:A:625:HIS:CE1	2.50	0.46
1:A:605:VAL:HA	1:A:624:GLY:HA2	1.97	0.45
1:A:535[B]:LEU:O	1:A:536[B]:ASN:CB	2.64	0.45
1:A:577[A]:LYS:HG3	5:A:1174[A]:HOH:O	2.16	0.45
1:A:384:THR:OG1	1:A:491[A]:ASP:OD1	2.35	0.44
1:A:430[B]:ASN:ND2	1:A:430[B]:ASN:C	2.71	0.44
1:A:326:LEU:HD22	1:A:358:PHE:CE1	2.52	0.44
1:A:322:VAL:HG21	1:A:360:ILE:HD11	2.00	0.43
1:A:401:THR:OG1	5:A:1098[A]:HOH:O	2.21	0.42
1:A:325:GLU:O	1:A:325:GLU:HG2	2.19	0.42
1:A:430[A]:ASN:ND2	5:A:1149:HOH:O	2.53	0.42
1:A:538[B]:GLU:OE2	1:A:598[B]:LYS:HD2	2.21	0.41

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	332/340 (98%)	322 (97%)	10 (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	Percentiles	
1	A	309/309 (100%)	309 (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 (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	389	GLN
1	A	397	GLN
1	A	456	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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		
IVIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
1	MLZ	A	601	1	8,9,10	0.59	0	4,9,11	1.85	1 (25%)

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
1	MLZ	A	601	1	-	0/7/8/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	601	MLZ	CM-NZ-CE	3.16	121.08	111.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	601	MLZ	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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).



Mal	Trino	Chain	Dog	Timle	Bond lengths			В	ond ang	les
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BME	A	705[A]	-	3,3,3	0.48	0	1,2,2	0.28	0
4	MES	A	706[A]	-	12,12,12	2.08	1 (8%)	14,16,16	1.31	2 (14%)

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	BME	A	705[A]	-	-	0/1/1/1	-
4	MES	A	706[A]	-	-	0/6/14/14	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(ext{\AA})$
4	A	706[A]	MES	C8-S	-6.87	1.67	1.77

All (2) bond angle outliers are listed below:

N	Iol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
	4	A	706[A]	MES	O1S-S-C8	3.05	110.59	106.92
	4	A	706[A]	MES	O3S-S-C8	2.55	109.89	105.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes	
4	A	706[A]	MES	1	0	

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	sed <rsrz></rsrz>		Z>2	$OWAB(A^2)$	Q < 0.9
1	A	307/340 (90%)	0.03	17 (5%) 2	25 24	13, 21, 40, 87	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	318	LEU	11.8
1	A	530	SER	7.6
1	A	626	THR	6.7
1	A	317	MET	6.1
1	A	322	VAL	5.8
1	A	526	GLY	5.4
1	A	627	ASN	4.3
1	A	324	SER	4.1
1	A	319	GLY	3.9
1	A	492	GLU	3.6
1	A	531	PRO	3.6
1	A	525	ASP	3.6
1	A	493	ARG	3.3
1	A	624	GLY	2.8
1	A	625	HIS	2.5
1	A	320	LYS	2.1
1	A	336[A]	ASN	2.1

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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	MLZ	A	601	10/11	0.93	0.10	21,25,41,41	0

6.3 Carbohydrates (i)

There are no monosaccharides 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^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	MES	A	706[A]	12/12	0.57	0.27	46,49,60,63	12
2	CA	A	704	1/1	0.90	0.25	49,49,49,49	0
3	BME	A	705[A]	4/4	0.93	0.17	17,18,21,23	4
2	CA	A	702	1/1	0.93	0.05	32,32,32,32	1
2	CA	A	703	1/1	0.96	0.11	42,42,42,42	1
2	CA	A	701	1/1	0.96	0.05	22,22,22,22	0

6.5 Other polymers (i)

There are no such residues in this entry.

