



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 4UG3
Title : B. subtilis GpsB N-terminal Domain
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Deposited on : 2015-03-21
Resolution : 2.80 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

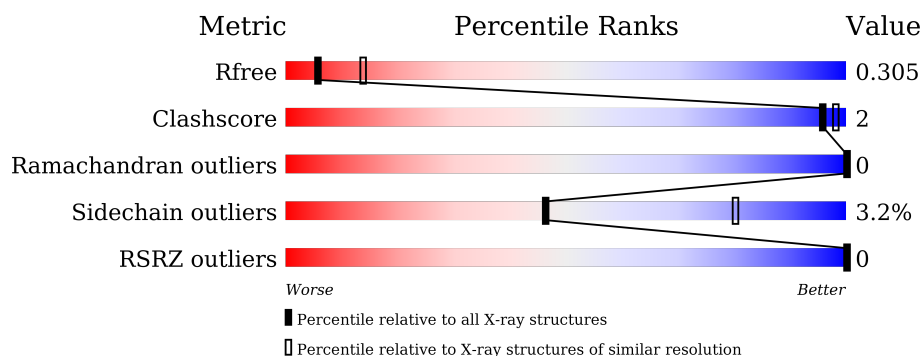
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107





i

X-RAY DIFFRACTION

A.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

Mol	Chain	Length	Quality of chain
1	A	71	
1	B	71	
1	C	71	
1	D	71	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1946 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 CELL CYCLE PROTEIN GPSB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	66	Total	C	N	O	S	0	0	0
			556	353	90	111	2			
1	B	60	Total	C	N	O	S	0	0	0
			511	326	82	102	1			
1	C	57	Total	C	N	O	S	0	0	0
			479	305	76	96	2			
1	D	47	Total	C	N	O	S	0	0	0
			398	257	62	78	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P0CI74
A	-1	SER	-	EXPRESSION TAG	UNP P0CI74
A	0	HIS	-	EXPRESSION TAG	UNP P0CI74
B	-2	GLY	-	EXPRESSION TAG	UNP P0CI74
B	-1	SER	-	EXPRESSION TAG	UNP P0CI74
B	0	HIS	-	EXPRESSION TAG	UNP P0CI74
C	-2	GLY	-	EXPRESSION TAG	UNP P0CI74
C	-1	SER	-	EXPRESSION TAG	UNP P0CI74
C	0	HIS	-	EXPRESSION TAG	UNP P0CI74
D	-2	GLY	-	EXPRESSION TAG	UNP P0CI74
D	-1	SER	-	EXPRESSION TAG	UNP P0CI74
D	0	HIS	-	EXPRESSION TAG	UNP P0CI74


- Molecule 2 is water.

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

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 ($RSRZ > 2$). 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: CELL CYCLE PROTEIN GPSB

Chain A: 



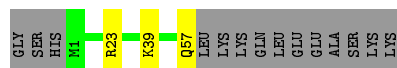
- Molecule 1: CELL CYCLE PROTEIN GPSB

Chain B: 



- Molecule 1: CELL CYCLE PROTEIN GPSB

Chain C: 



- Molecule 1: CELL CYCLE PROTEIN GPSB

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.67Å 39.89Å 77.28Å 90.00° 103.28° 90.00°	Depositor
Resolution (Å)	50.30 – 2.80 50.29 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.30-2.80) 100.0 (50.29-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.216 , 0.301 0.219 , 0.305	Depositor DCC
R_{free} test set	801 reflections (11.47%)	DCC
Wilson B-factor (Å ²)	51.0	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 23.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 7785 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	1946	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.53% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.55	0/562	0.76	0/747
1	B	0.59	0/516	0.70	0/685
1	C	0.52	0/484	0.70	0/644
1	D	0.63	0/403	0.73	0/535
All	All	0.57	0/1965	0.72	0/2611

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 [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	556	0	561	4	0
1	B	511	0	517	3	0
1	C	479	0	481	1	0
1	D	398	0	402	2	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0
All	All	1946	0	1961	6	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 2.

The worst 5 of 6 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:D:14:LEU:HD12	1:D:38:ILE:HD13	1.76	0.66
1:A:64:GLU:N	1:A:64:GLU:OE1	2.41	0.54
1:A:22:VAL:O	1:B:22:VAL:O	2.32	0.48
1:A:59:LYS:NZ	1:B:54:GLU:OE2	2.43	0.47
1:C:23:ARG:NE	1:D:20:THR:CG2	2.80	0.44

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	64/71 (90%)	62 (97%)	2 (3%)	0	100	100
1	B	58/71 (82%)	57 (98%)	1 (2%)	0	100	100
1	C	55/71 (78%)	53 (96%)	2 (4%)	0	100	100
1	D	45/71 (63%)	44 (98%)	1 (2%)	0	100	100
All	All	222/284 (78%)	216 (97%)	6 (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	62/65 (95%)	60 (97%)	2 (3%)	46	80
1	B	57/65 (88%)	56 (98%)	1 (2%)	66	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	53/65 (82%)	51 (96%)	2 (4%)	40	74
1	D	44/65 (68%)	42 (96%)	2 (4%)	34	68
All	All	216/260 (83%)	209 (97%)	7 (3%)	46	80

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	39	LYS
1	D	23	ARG
1	C	57	GLN
1	A	60	LYS
1	D	14	LEU

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

Mol	Chain	Res	Type
1	B	57	GLN
1	C	55	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](#)

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](#)

There are no ligands in this entry.

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, 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	66/71 (92%)	-0.42	0 100 100	27, 51, 71, 91	0
1	B	60/71 (84%)	-0.39	0 100 100	31, 45, 79, 89	0
1	C	57/71 (80%)	-0.38	0 100 100	33, 53, 91, 94	0
1	D	47/71 (66%)	-0.37	0 100 100	34, 49, 83, 94	0
All	All	230/284 (80%)	-0.39	0 100 100	27, 51, 85, 94	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.