



Full wwPDB X-ray Structure Validation Report

May 29, 2014 – 12:32 AM EDT

PDB ID : 4L1G
Title : Crystal structure of the Bc1960 peptidoglycan N-acetylglucosamine deacetylase from *Bacillus cereus*
Authors : Tsalafouta, A.; Fadouloglou, V.E.; Kokkinidis, M.
Deposited on : 2013-06-03
Resolution : 2.34 Å(reported)

This is a full wwPDB validation 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/ValidationPDFNotes.html>

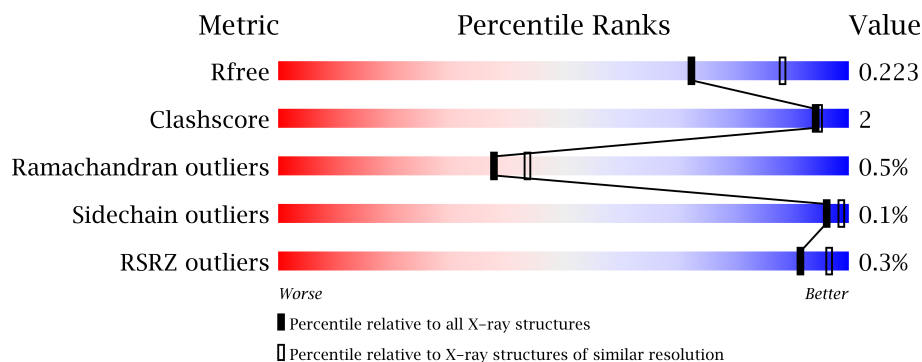
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : **FAILED**
Xtriage (Phenix) : dev-1439
EDS : stable23161
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23161

1 Overall quality at a glance

The reported resolution of this entry is 2.34 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}	66092	4049 (2.38-2.30)
Clashscore	79885	1094 (2.36-2.32)
Ramachandran outliers	78287	1080 (2.36-2.32)
Sidechain outliers	78261	1081 (2.36-2.32)
RSRZ outliers	66119	4050 (2.38-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	273	
1	B	273	
1	C	273	
1	D	273	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	ACT	A	302	-	X
3	ACT	C	303	-	X

2 Entry composition i

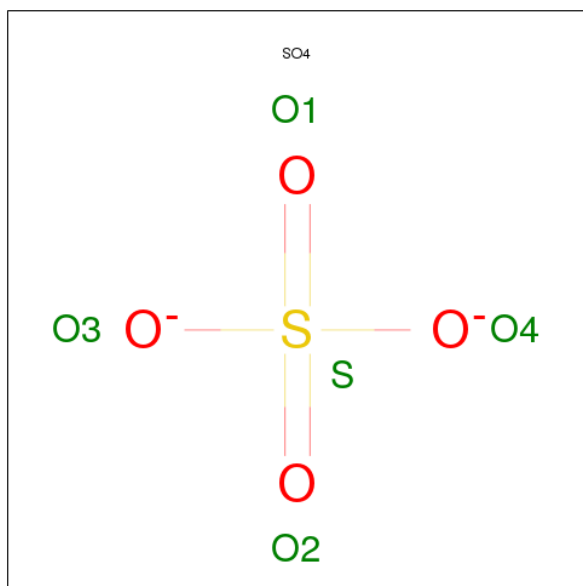
There are 4 unique types of molecules in this entry. The entry contains 7688 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Peptidoglycan N-acetylglucosamine deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	1	0
			1721	1109	290	320	2			
1	B	214	Total	C	N	O	S	0	1	0
			1723	1110	292	319	2			
1	C	214	Total	C	N	O	S	0	2	0
			1730	1114	292	322	2			
1	D	213	Total	C	N	O	S	0	0	0
			1699	1093	286	318	2			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



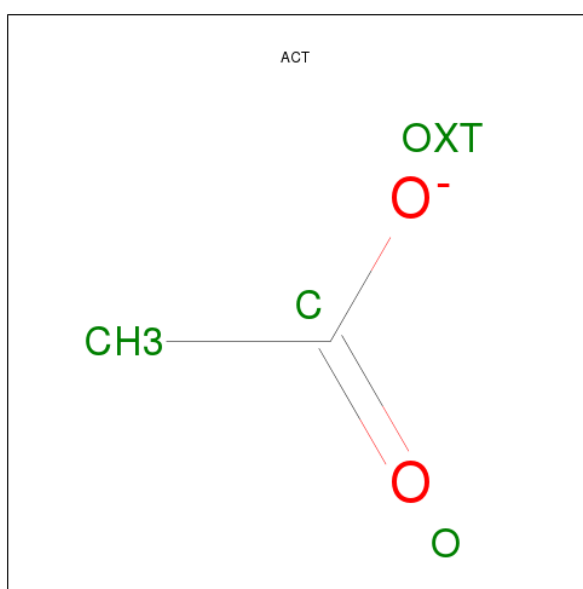
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	224	Total	O	0	0
			224	224		

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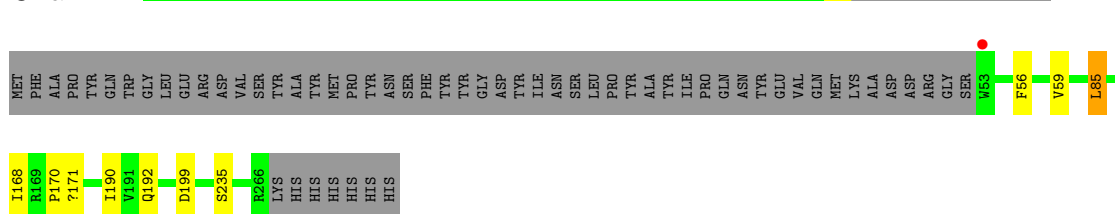
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	199	Total 199	O 199	0	0
4	C	189	Total 189	O 189	0	0
4	D	157	Total 157	O 157	0	0

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 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: Peptidoglycan N-acetylglucosamine deacetylase

Chain A:



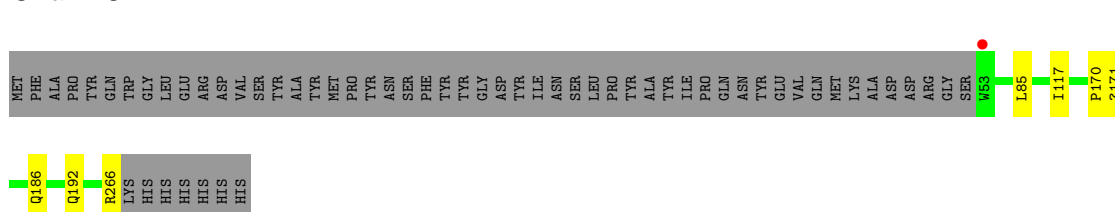
- Molecule 1: Peptidoglycan N-acetylglucosamine deacetylase

Chain B:



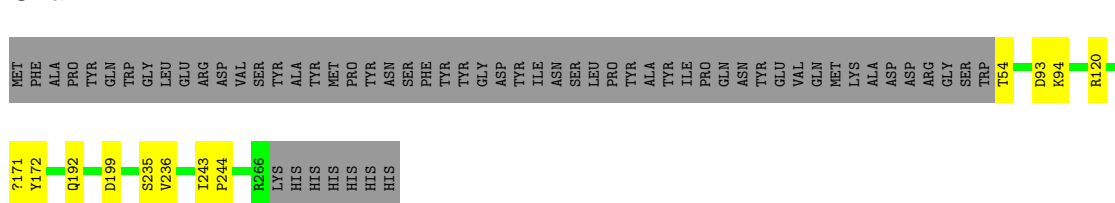
- Molecule 1: Peptidoglycan N-acetylglucosamine deacetylase

Chain C:



- Molecule 1: Peptidoglycan N-acetylglucosamine deacetylase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	92.66Å 92.66Å 242.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.64 – 2.34 86.57 – 2.34	Depositor EDS
% Data completeness (in resolution range)	98.9 (30.64-2.34) 98.9 (86.57-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.36 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1069)	Depositor
R, R_{free}	0.179 , 0.223 0.179 , 0.223	Depositor DCC
R_{free} test set	2283 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.470	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 21.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 45341 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7688	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 31.39 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.1159e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, PXU, ACT

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	0.24	0/1761	0.40	0/2390
1	B	0.24	0/1763	0.40	0/2393
1	C	0.23	0/1770	0.40	0/2403
1	D	0.23	0/1734	0.40	0/2354
All	All	0.23	0/7028	0.40	0/9540

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	D	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	171	PXU	Mainchain
1	D	171	PXU	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1721	0	1694	5	0
1	B	1723	0	1696	7	0
1	C	1730	0	1694	3	0
1	D	1699	0	1666	6	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	A	4	0	3	0	0
3	B	4	0	3	1	0
3	C	4	0	3	0	0
3	D	4	0	3	1	0
4	A	224	0	0	0	0
4	B	199	0	0	2	0
4	C	189	0	0	0	0
4	D	157	0	0	1	0
All	All	7688	0	6762	21	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

All (21) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:54:THR:N	4:D:528:HOH:O	2.31	0.63
1:C:186:GLN:O	1:C:266:ARG:HD2	2.08	0.53
1:B:199:ASP:HB2	1:B:235:SER:HB2	1.95	0.48
1:B:172:TYR:H	3:B:302:ACT:H3	1.79	0.48
1:D:199:ASP:HB2	1:D:235:SER:HB2	1.97	0.47
1:A:168:ILE:HB	1:A:190:ILE:HG12	1.96	0.47
1:D:172:TYR:H	3:D:303:ACT:H3	1.80	0.47
1:B:72:LYS:NZ	4:B:581:HOH:O	2.35	0.46
1:B:56:PHE:O	1:B:59:VAL:HG22	2.16	0.44
1:B:116:ASN:HB2	4:B:556:HOH:O	2.16	0.44
1:C:85:LEU:HD23	1:C:117:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:93:ASP:OD2	1:D:120:ARG:NH2	2.41	0.43
1:C:170:PRO:HA	1:C:171:PXU:H8	1.85	0.43
1:B:263:SER:OG	1:B:265:GLU:OE1	2.37	0.42
1:A:199:ASP:HB2	1:A:235:SER:HB2	2.00	0.42
1:A:170:PRO:HA	1:A:171:PXU:H8	1.62	0.42
1:B:170:PRO:HA	1:B:171:PXU:H8	1.78	0.42
1:A:56:PHE:O	1:A:59:VAL:HG22	2.21	0.40
1:A:85:LEU:HA	1:A:85:LEU:HD12	1.83	0.40
1:D:94:LYS:HE2	1:D:236:VAL:HG12	2.03	0.40
1:D:243:ILE:HB	1:D:244:PRO:HD3	2.04	0.40

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	212/273 (78%)	206 (97%)	5 (2%)	1 (0%)	38	44
1	B	212/273 (78%)	206 (97%)	5 (2%)	1 (0%)	38	44
1	C	213/273 (78%)	209 (98%)	3 (1%)	1 (0%)	38	44
1	D	210/273 (77%)	204 (97%)	5 (2%)	1 (0%)	38	44
All	All	847/1092 (78%)	825 (97%)	18 (2%)	4 (0%)	38	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	GLN
1	C	192	GLN
1	D	192	GLN
1	B	192	GLN

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	186/237 (78%)	185 (100%)	1 (0%)	94	97
1	B	185/237 (78%)	185 (100%)	0	100	100
1	C	185/237 (78%)	185 (100%)	0	100	100
1	D	182/237 (77%)	182 (100%)	0	100	100
All	All	738/948 (78%)	737 (100%)	1 (0%)	96	99

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	LEU

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 chains in this entry.

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

Mogul failed to run properly - this section will therefore be empty.

5.5 Carbohydrates ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.6 Ligand geometry ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.7 Other polymers ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.8 Polymer linkage issues

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	213/273 (78%)	-0.26	1 (0%) 88 95	8, 16, 29, 42	0
1	B	213/273 (78%)	-0.26	1 (0%) 88 95	6, 15, 27, 46	0
1	C	213/273 (78%)	-0.27	1 (0%) 88 95	6, 16, 31, 82	0
1	D	212/273 (77%)	-0.17	0 100 100	10, 21, 39, 54	0
All	All	851/1092 (77%)	-0.24	3 (0%) 91 96	6, 17, 34, 82	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	53	TRP	3.3
1	B	53	TRP	3.2
1	A	53	TRP	2.4

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

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	PXU	C	171	8/9	0.12	-0.17	14,15,16,16	0
1	PXU	A	171	8/9	0.12	-0.32	14,15,15,15	0
1	PXU	D	171	8/9	0.10	-0.90	16,16,17,18	0
1	PXU	B	171	8/9	0.10	-0.94	13,14,16,17	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ACT	A	302	4/4	0.18	3.35	28,31,32,45	0
3	ACT	C	303	4/4	0.17	2.53	30,35,40,43	0
3	ACT	B	302	4/4	0.13	1.71	31,32,36,42	0
2	SO4	D	301	5/5	0.14	1.67	18,18,18,18	0
2	SO4	D	302	5/5	0.17	1.52	61,63,63,64	0
2	SO4	C	301	5/5	0.13	1.49	12,14,14,15	0
3	ACT	D	303	4/4	0.14	0.94	32,33,34,39	0
2	SO4	C	302	5/5	0.13	-0.26	67,69,69,70	0
2	SO4	B	301	5/5	0.09	-0.70	58,60,60,61	0
2	SO4	A	301	5/5	0.09	-1.28	49,50,51,51	0

6.5 Other polymers ⓘ

There are no such residues in this entry.