



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 12:01 PM UTC

PDB ID : 4N8Y / pdb_00004n8y
Title : Crystal structure of a trap periplasmic solute binding protein from bradyrhizobium sp. btail b (bbta_0128), target EFI-510056 (bbta_0128), complex with alpha/beta-d-galacturonate
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Deposited on : 2013-10-18
Resolution : 1.50 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)

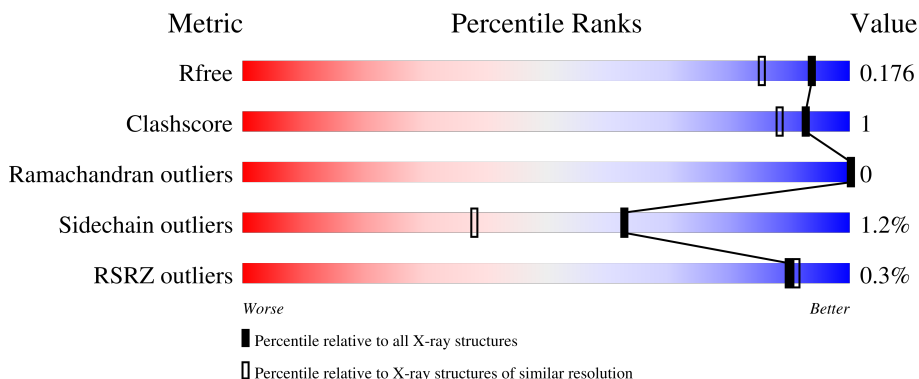
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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}	180053	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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 $\leq 5\%$. 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	330	
1	B	330	

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10325 atoms, of which 4704 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.

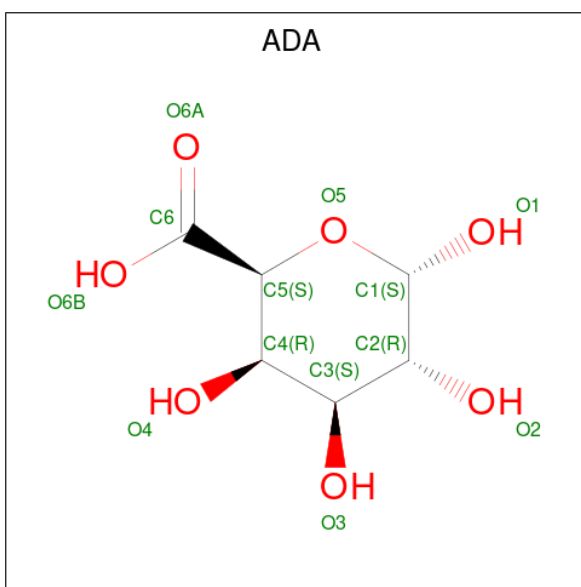
- Molecule 1 is a protein called Putative TRAP-type C4-dicarboxylate transport system, binding periplasmic protein (DctP subunit).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	300	Total	C	H	N	O	S	0	2	0
			4662	1474	2340	398	439	11			
1	B	300	Total	C	H	N	O	S	0	6	0
			4703	1485	2364	401	441	12			

There are 14 discrepancies between the modelled and reference sequences:

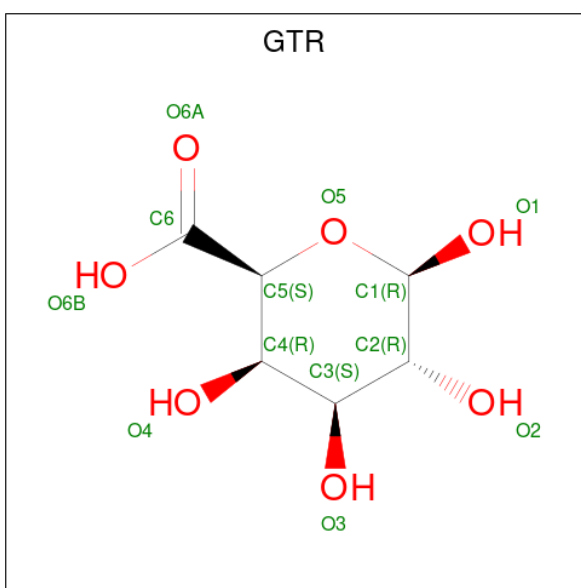
Chain	Residue	Modelled	Actual	Comment	Reference
A	324	ALA	-	expression tag	UNP A5E8D2
A	325	GLU	-	expression tag	UNP A5E8D2
A	326	ASN	-	expression tag	UNP A5E8D2
A	327	LEU	-	expression tag	UNP A5E8D2
A	328	TYR	-	expression tag	UNP A5E8D2
A	329	PHE	-	expression tag	UNP A5E8D2
A	330	GLN	-	expression tag	UNP A5E8D2
B	324	ALA	-	expression tag	UNP A5E8D2
B	325	GLU	-	expression tag	UNP A5E8D2
B	326	ASN	-	expression tag	UNP A5E8D2
B	327	LEU	-	expression tag	UNP A5E8D2
B	328	TYR	-	expression tag	UNP A5E8D2
B	329	PHE	-	expression tag	UNP A5E8D2
B	330	GLN	-	expression tag	UNP A5E8D2

- Molecule 2 is alpha-D-galactopyranuronic acid (CCD ID: ADA) (formula: C₆H₁₀O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	1
			13	6	7		
2	B	1	Total	C	O	0	1
			13	6	7		

- Molecule 3 is beta-D-galactopyranuronic acid (CCD ID: GTR) (formula: $C_6H_{10}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			13	6	7		
3	B	1	Total	C	O	0	1
			13	6	7		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	440	Total 440	O 440	0	0
4	B	468	Total 468	O 468	0	0

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.

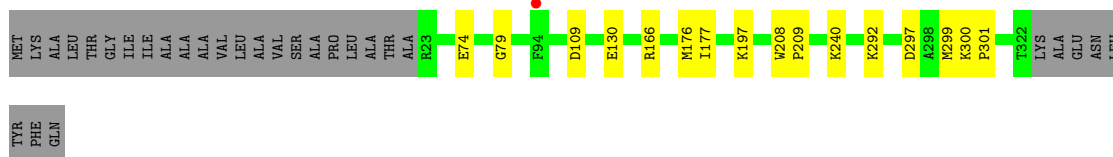
- Molecule 1: Putative TRAP-type C4-dicarboxylate transport system, binding periplasmic protein (DctP subunit)

Chain A:  88% 9%



- Molecule 1: Putative TRAP-type C4-dicarboxylate transport system, binding periplasmic protein (DctP subunit)

Chain B:  86% 5% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.80Å 85.70Å 86.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.09 – 1.50 29.09 – 1.50	Depositor EDS
% Data completeness (in resolution range)	95.8 (29.09-1.50) 95.8 (29.09-1.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.27 (at 1.50Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.144 , 0.175 0.145 , 0.176	Depositor DCC
R_{free} test set	4508 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.646	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.025 for -h,l,k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10325	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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 30.18 % 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 1.3654e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: ADA, GTR

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	0.57	0/2374	0.73	0/3214
1	B	0.60	0/2407	0.74	0/3257
All	All	0.59	0/4781	0.73	0/6471

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	2322	2340	2338	4	0
1	B	2339	2364	2338	10	0
2	A	13	0	9	0	0
2	B	13	0	9	0	0
3	A	13	0	9	0	0
3	B	13	0	9	0	0
4	A	440	0	0	2	1
4	B	468	0	0	3	1
All	All	5621	4704	4712	14	1

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 1.

The worst 5 of 14 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:155:LYS:NZ	4:A:909:HOH:O	2.19	0.74
1:A:266:LYS:NZ	4:A:928:HOH:O	2.37	0.57
1:B:130:GLU:OE1	4:B:827:HOH:O	2.16	0.56
1:B:79:GLY:HA3	1:B:240[A]:LYS:HD3	1.94	0.49
1:A:166:ARG:HD3	1:A:166:ARG:C	2.43	0.44

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:916:HOH:O	4:B:935:HOH:O[2_564]	2.16	0.04

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	300/330 (91%)	295 (98%)	5 (2%)	0	100	100
1	B	304/330 (92%)	301 (99%)	3 (1%)	0	100	100
All	All	604/660 (92%)	596 (99%)	8 (1%)	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	250/268 (93%)	247 (99%)	3 (1%)	63	38
1	B	254/268 (95%)	250 (98%)	4 (2%)	55	28
All	All	504/536 (94%)	497 (99%)	7 (1%)	63	33

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

Mol	Chain	Res	Type
1	B	177	ILE
1	B	292	LYS
1	B	300[B]	LYS
1	B	300[A]	LYS
1	A	304	GLN

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

Mol	Chain	Res	Type
1	A	183	ASN
1	B	92	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are 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		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ADA	A	401[A]	-	13,13,13	1.05	0	18,19,19	1.01	0
3	GTR	B	402[B]	-	13,13,13	1.04	0	18,19,19	0.94	1 (5%)
3	GTR	A	402[B]	-	13,13,13	1.17	1 (7%)	18,19,19	1.20	1 (5%)
2	ADA	B	401[A]	-	13,13,13	0.98	0	18,19,19	0.90	0

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
2	ADA	A	401[A]	-	-	0/4/24/24	0/1/1/1
3	GTR	B	402[B]	-	-	0/4/24/24	0/1/1/1
3	GTR	A	402[B]	-	-	2/4/24/24	0/1/1/1
2	ADA	B	401[A]	-	-	1/4/24/24	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402[B]	GTR	C1-C2	-2.03	1.47	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402[B]	GTR	O3-C3-C2	-2.09	105.44	110.38
3	A	402[B]	GTR	O3-C3-C2	-2.09	105.45	110.38

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401[A]	ADA	C4-C5-C6-O6A
3	A	402[B]	GTR	C4-C5-C6-O6A
3	A	402[B]	GTR	C4-C5-C6-O6B

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, 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	300/330 (90%)	-0.53	1 (0%) 90 91	6, 19, 48, 82	1 (0%)
1	B	300/330 (90%)	-0.64	1 (0%) 90 91	7, 17, 42, 62	3 (1%)
All	All	600/660 (90%)	-0.59	2 (0%) 90 91	6, 18, 45, 82	4 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	290	ALA	4.2
1	B	94	PHE	2.7

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 oligosaccharides 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, 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	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ADA	A	401[A]	13/13	0.97	0.05	6,9,12,17	13
2	ADA	B	401[A]	13/13	0.98	0.04	6,9,10,12	13
3	GTR	A	402[B]	13/13	0.98	0.04	8,10,12,12	13

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GTR	B	402[B]	13/13	0.98	0.04	6,9,11,12	13

6.5 Other polymers [i](#)

There are no such residues in this entry.