



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 04:16 PM GMT

PDB ID : 4A0R  
Title : Structure of bifunctional DAPA aminotransferase-DTB synthetase from *Arabidopsis thaliana* bound to dethiobiotin (DTB).  
Authors : Cobessi, D.; Dumas, R.; Pautre, V.; Meinguet, C.; Ferrer, J.L.; Alban, C.  
Deposited on : 2011-09-12  
Resolution : 2.68 Å(reported)

This is a full wwPDB validation 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 <http://wwpdb.org/ValidationPDFNotes.html>

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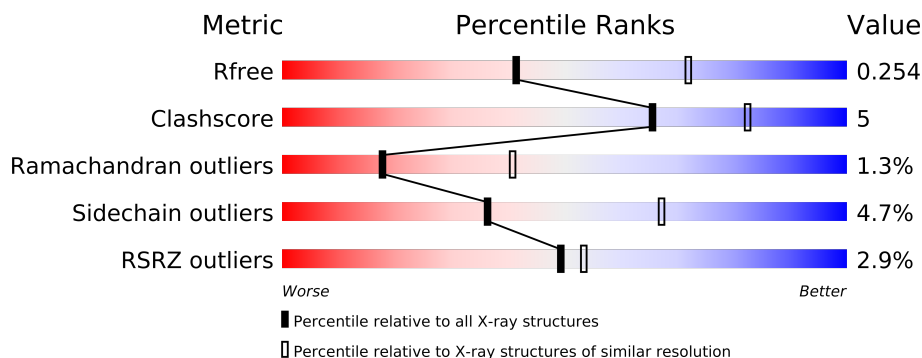
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
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) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.68 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	2010 (2.70-2.66)
Clashscore	79885	2450 (2.70-2.66)
Ramachandran outliers	78287	2410 (2.70-2.66)
Sidechain outliers	78261	2410 (2.70-2.66)
RSRZ outliers	66119	2013 (2.70-2.66)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	831	
1	B	831	

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

Mol	Type	Chain	Res	Geometry	Electron density
4	DTB	A	1811	-	X
4	DTB	B	1808	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11693 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 ADENOSYLMETHIONINE-8-AMINO-7-OXONONANOAT EAMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	755	Total	C	N	O	S	0	2	0
			5788	3703	974	1079	32			
1	B	748	Total	C	N	O	S	0	2	0
			5704	3652	961	1059	32			

There are 40 discrepancies between the modelled and reference sequences:

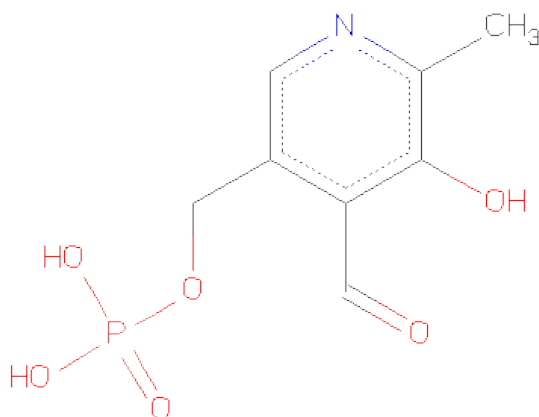
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	GLY	-	EXPRESSION TAG	UNP B0F481
A	-18	SER	-	EXPRESSION TAG	UNP B0F481
A	-17	SER	-	EXPRESSION TAG	UNP B0F481
A	-16	HIS	-	EXPRESSION TAG	UNP B0F481
A	-15	HIS	-	EXPRESSION TAG	UNP B0F481
A	-14	HIS	-	EXPRESSION TAG	UNP B0F481
A	-13	HIS	-	EXPRESSION TAG	UNP B0F481
A	-12	HIS	-	EXPRESSION TAG	UNP B0F481
A	-11	HIS	-	EXPRESSION TAG	UNP B0F481
A	-10	SER	-	EXPRESSION TAG	UNP B0F481
A	-9	SER	-	EXPRESSION TAG	UNP B0F481
A	-8	GLY	-	EXPRESSION TAG	UNP B0F481
A	-7	LEU	-	EXPRESSION TAG	UNP B0F481
A	-6	VAL	-	EXPRESSION TAG	UNP B0F481
A	-5	PRO	-	EXPRESSION TAG	UNP B0F481
A	-4	ARG	-	EXPRESSION TAG	UNP B0F481
A	-3	GLY	-	EXPRESSION TAG	UNP B0F481
A	-2	SER	-	EXPRESSION TAG	UNP B0F481
A	-1	HIS	-	EXPRESSION TAG	UNP B0F481
A	0	MET	-	EXPRESSION TAG	UNP B0F481
B	-19	GLY	-	EXPRESSION TAG	UNP B0F481
B	-18	SER	-	EXPRESSION TAG	UNP B0F481
B	-17	SER	-	EXPRESSION TAG	UNP B0F481
B	-16	HIS	-	EXPRESSION TAG	UNP B0F481

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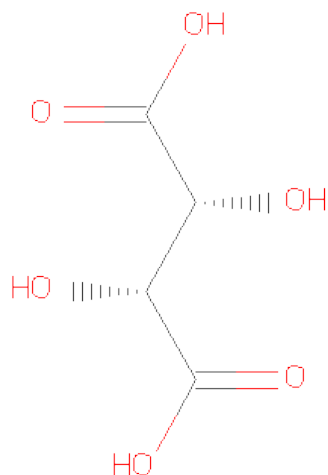
Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	EXPRESSION TAG	UNP B0F481
B	-14	HIS	-	EXPRESSION TAG	UNP B0F481
B	-13	HIS	-	EXPRESSION TAG	UNP B0F481
B	-12	HIS	-	EXPRESSION TAG	UNP B0F481
B	-11	HIS	-	EXPRESSION TAG	UNP B0F481
B	-10	SER	-	EXPRESSION TAG	UNP B0F481
B	-9	SER	-	EXPRESSION TAG	UNP B0F481
B	-8	GLY	-	EXPRESSION TAG	UNP B0F481
B	-7	LEU	-	EXPRESSION TAG	UNP B0F481
B	-6	VAL	-	EXPRESSION TAG	UNP B0F481
B	-5	PRO	-	EXPRESSION TAG	UNP B0F481
B	-4	ARG	-	EXPRESSION TAG	UNP B0F481
B	-3	GLY	-	EXPRESSION TAG	UNP B0F481
B	-2	SER	-	EXPRESSION TAG	UNP B0F481
B	-1	HIS	-	EXPRESSION TAG	UNP B0F481
B	0	MET	-	EXPRESSION TAG	UNP B0F481

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula:  $C_8H_{10}NO_6P$ ).



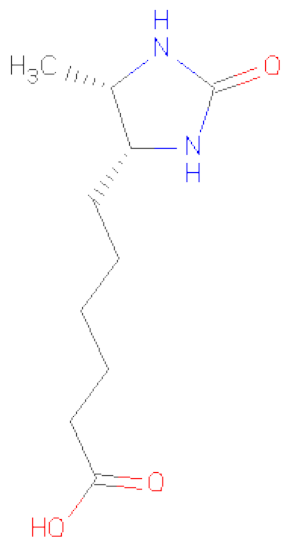
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula:  $C_4H_6O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	4	6		
3	B	1	Total	C	O	0	0
			10	4	6		

- Molecule 4 is 6-(5-METHYL-2-OXO-IMIDAZOLIDIN-4-YL)-HEXANOICACID (three-letter code: DTB) (formula: C<sub>10</sub>H<sub>18</sub>N<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			15	10	2	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			15	10	2	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	70	Total	O	0	0
			70	70		
5	B	51	Total	O	0	0
			51	51		







## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	246.67Å 76.63Å 79.84Å 90.00° 108.02° 90.00°	Depositor
Resolution (Å)	40.44 – 2.68 40.72 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.44-2.68) 99.5 (40.72-2.68)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.90 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.184 , 0.259 0.179 , 0.254	Depositor DCC
$R_{free}$ test set	2000 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.6	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 37.3	EDS
Estimated twinning fraction	0.027 for -h+k-l,-l,-k 0.005 for -h-k-l,l,k 0.023 for -h-2*l,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 39935 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11693	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: DTB, TLA, PLP

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	0.48	1/5922 (0.0%)	0.64	3/8050 (0.0%)
1	B	0.48	2/5836 (0.0%)	0.65	3/7938 (0.0%)
All	All	0.48	3/11758 (0.0%)	0.64	6/15988 (0.0%)

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	307	ARG	CZ-NH1	-11.03	1.18	1.33
1	A	307	ARG	CZ-NH1	-9.79	1.20	1.33
1	B	307	ARG	CZ-NH2	-5.56	1.25	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	307	ARG	NE-CZ-NH2	17.58	129.09	120.30
1	B	307	ARG	NE-CZ-NH2	16.14	128.37	120.30
1	B	307	ARG	NH1-CZ-NH2	-8.37	110.19	119.40
1	A	44	PRO	N-CA-CB	6.33	110.90	103.30
1	A	307	ARG	NH1-CZ-NH2	-5.93	112.87	119.40
1	B	44	PRO	N-CA-CB	5.83	110.29	103.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	45	SER	Peptide

## 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	5788	0	0	27	0
1	B	5704	0	0	33	0
2	A	15	0	0	0	0
2	B	15	0	0	0	0
3	A	10	0	3	2	0
3	B	10	0	4	4	0
4	A	15	0	17	1	0
4	B	15	0	17	4	0
5	A	70	0	0	0	0
5	B	51	0	0	2	0
All	All	11693	0	41	60	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 5.

All (60) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:25:SER:N	3:B:1809:TLA:HA	1.79	0.79
1:A:223:GLY:CA	4:B:1808:DTB:HCA1	2.17	0.73
1:B:768:ARG:NH2	1:B:773:PHE:CE2	2.59	0.70
1:B:28:LYS:NZ	3:B:1809:TLA:H2	2.09	0.67
1:B:490:THR:CG2	1:B:490:THR:O	2.45	0.65
1:B:547:ASP:OD1	1:B:599:ARG:NH1	2.29	0.65
1:A:490:THR:O	1:A:490:THR:CG2	2.46	0.64
1:B:522:PRO:CG	1:B:525:PHE:CD2	2.80	0.64
1:B:535:PHE:CD1	1:B:535:PHE:N	2.66	0.64
1:A:223:GLY:N	4:B:1808:DTB:HCA1	2.15	0.62
1:B:730:SER:O	1:B:747:LYS:NZ	2.36	0.58
1:A:104:LEU:O	1:A:105:ASN:C	2.43	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:547:ASP:OD1	1:A:599:ARG:NH1	2.38	0.56
1:B:588:GLY:N	5:B:2040:HOH:O	2.40	0.55
1:A:60:GLY:O	1:A:64:ASP:N	2.40	0.55
1:A:676:LEU:N	1:B:494:GLN:OE1	2.40	0.54
1:A:768:ARG:NH2	1:A:773:PHE:CE2	2.75	0.54
1:B:42:GLN:OE1	1:B:42:GLN:CA	2.57	0.52
1:A:28:LYS:NZ	3:A:1810:TLA:H3	2.26	0.51
1:B:370:TRP:N	1:B:370:TRP:CD1	2.78	0.51
1:A:706:ASN:O	1:A:715:ARG:N	2.43	0.51
1:B:417:VAL:CG1	1:B:642:PHE:CZ	2.94	0.51
1:A:659:THR:CG2	1:A:660:ASP:N	2.73	0.51
1:A:725:GLN:O	1:A:728:SER:N	2.44	0.51
1:A:805:LEU:O	1:A:808:PHE:N	2.45	0.49
1:B:307:ARG:NH2	5:B:2018:HOH:O	2.45	0.49
1:B:659:THR:CG2	1:B:660:ASP:N	2.75	0.49
1:B:191:GLY:N	3:B:1809:TLA:O11	2.46	0.49
1:A:370:TRP:N	1:A:370:TRP:CD1	2.80	0.48
1:B:60:GLY:O	1:B:64:ASP:N	2.46	0.48
1:A:417:VAL:CG1	1:A:642:PHE:CZ	2.97	0.48
1:B:44:PRO:O	1:B:45:SER:CB	2.63	0.46
1:B:143:SER:OG	1:B:145:HIS:ND1	2.49	0.45
4:B:1808:DTB:HCS	3:B:1809:TLA:C1	2.46	0.45
1:B:59:THR:N	4:B:1808:DTB:O	2.49	0.45
1:A:562:HIS:ND1	1:A:562:HIS:N	2.64	0.45
1:B:558:TYR:O	1:B:562:HIS:ND1	2.50	0.44
1:A:248:GLU:OE2	1:A:274:VAL:N	2.51	0.44
1:A:708:THR:N	1:A:713:THR:O	2.51	0.44
1:B:354:SER:O	1:B:355:ASP:O	2.36	0.44
1:B:538:ARG:NH1	1:B:737:VAL:O	2.52	0.43
1:A:497:TRP:CD1	1:B:501:ARG:CD	3.02	0.43
1:A:353:ALA:C	1:A:355:ASP:N	2.71	0.43
1:A:808:PHE:N	1:A:808:PHE:CD1	2.86	0.43
1:A:700:ASP:C	1:A:700:ASP:OD1	2.57	0.42
1:A:444:ARG:NH1	1:A:673:LYS:O	2.52	0.42
1:B:582:PRO:O	1:B:584:ILE:N	2.52	0.42
1:B:562:HIS:ND1	1:B:562:HIS:N	2.67	0.42
1:A:708:THR:CG2	1:A:711:GLY:N	2.83	0.42
1:B:805:LEU:O	1:B:807:GLU:N	2.53	0.42
3:A:1810:TLA:O2	3:A:1810:TLA:O41	2.38	0.41
1:B:78:SER:OG	1:B:83:ILE:O	2.38	0.41
1:B:224:GLY:O	1:B:228:THR:OG1	2.38	0.41
1:A:281[A]:ASP:O	1:A:282:LEU:CB	2.68	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:208:ARG:N	1:B:209:PRO:CD	2.84	0.41
1:B:593:VAL:CG1	1:B:594:ASP:N	2.84	0.41
4:A:1811:DTB:HCB1	1:B:222:LEU:CD1	2.50	0.41
1:B:204:CYS:O	1:B:240:TYR:OH	2.40	0.40
1:A:445:LYS:CE	1:A:449:ASP:OD2	2.69	0.40
1:A:615:ASP:C	1:A:615:ASP:OD1	2.58	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	745/831 (90%)	687 (92%)	46 (6%)	12 (2%)	14	33
1	B	736/831 (89%)	696 (95%)	33 (4%)	7 (1%)	22	49
All	All	1481/1662 (89%)	1383 (93%)	79 (5%)	19 (1%)	18	41

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	490	THR
1	B	355	ASP
1	B	490	THR
1	A	373	GLY
1	A	561	LYS
1	B	45	SER
1	B	373	GLY
1	A	105	ASN
1	A	355	ASP
1	A	500	GLY
1	A	643	ALA
1	A	644	LYS
1	B	643	ALA
1	B	644	LYS

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Mol	Chain	Res	Type
1	A	421	TRP
1	A	725	GLN
1	B	421	TRP
1	A	708	THR
1	A	103	GLY

### 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	622/704 (88%)	592 (95%)	30 (5%)	35	66
1	B	605/704 (86%)	575 (95%)	30 (5%)	34	64
All	All	1227/1408 (87%)	1167 (95%)	60 (5%)	36	65

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	13	ASN
1	A	63	SER
1	A	67	SER
1	A	74	LEU
1	A	109	SER
1	A	110	GLU
1	A	117	ASN
1	A	120	ASP
1	A	146	LEU
1	A	154	THR
1	A	193	VAL
1	A	253	VAL
1	A	281[A]	ASP
1	A	281[B]	ASP
1	A	308	LEU
1	A	317	LEU
1	A	321	VAL
1	A	370	TRP
1	A	384	ARG

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Mol	Chain	Res	Type
1	A	417	VAL
1	A	428	SER
1	A	466	VAL
1	A	562	HIS
1	A	628	THR
1	A	655	VAL
1	A	664	ASP
1	A	713	THR
1	A	727	SER
1	A	764	LEU
1	B	10	LEU
1	B	63	SER
1	B	67	SER
1	B	74	LEU
1	B	120	ASP
1	B	146	LEU
1	B	154	THR
1	B	193	VAL
1	B	281[A]	ASP
1	B	281[B]	ASP
1	B	307	ARG
1	B	317	LEU
1	B	321	VAL
1	B	370	TRP
1	B	408	LYS
1	B	417	VAL
1	B	428	SER
1	B	466	VAL
1	B	501	ARG
1	B	535	PHE
1	B	539	ASP
1	B	562	HIS
1	B	624	LEU
1	B	628	THR
1	B	655	VAL
1	B	664	ASP
1	B	708	THR
1	B	709	SER
1	B	723	VAL
1	B	763	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 ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

6 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	PLP	A	1644	1	14,15,16	1.91	3 (21%)	20,22,23	1.59	3 (15%)
3	TLA	A	1810	-	9,9,9	2.00	2 (22%)	12,12,12	2.09	6 (50%)
4	DTB	A	1811	-	15,15,15	0.98	1 (6%)	19,19,19	1.06	2 (10%)
2	PLP	B	1644	1	14,15,16	1.90	3 (21%)	20,22,23	1.67	1 (5%)
4	DTB	B	1808	-	15,15,15	1.00	1 (6%)	19,19,19	1.63	4 (21%)
3	TLA	B	1809	-	9,9,9	2.02	3 (33%)	12,12,12	2.01	6 (50%)

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	PLP	A	1644	1	-	0/6/6/8	0/1/1/1
3	TLA	A	1810	-	-	0/12/12/12	0/0/0/0
4	DTB	A	1811	-	-	0/8/20/20	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PLP	B	1644	1	-	0/6/6/8	0/1/1/1
4	DTB	B	1808	-	-	0/8/20/20	0/1/1/1
3	TLA	B	1809	-	-	0/12/12/12	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1644	PLP	O3-C3	-5.60	1.23	1.37
2	A	1644	PLP	O3-C3	-5.50	1.23	1.37
3	A	1810	TLA	C3-C4	-3.97	1.47	1.52
3	B	1809	TLA	C2-C1	-3.62	1.47	1.52
3	A	1810	TLA	O3-C3	-2.98	1.36	1.42
2	A	1644	PLP	C2-N1	2.61	1.38	1.33
2	B	1644	PLP	C2-N1	2.28	1.38	1.33
2	A	1644	PLP	C6-N1	2.28	1.39	1.34
3	B	1809	TLA	O2-C2	-2.24	1.37	1.42
4	A	1811	DTB	CN-N1	-2.21	1.32	1.35
3	B	1809	TLA	C3-C4	-2.07	1.50	1.52
2	B	1644	PLP	C3-C2	-2.05	1.39	1.40
4	B	1808	DTB	CN-N2	-2.03	1.32	1.35

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1644	PLP	O4P-C5A-C5	5.91	121.27	109.26
2	A	1644	PLP	O4P-C5A-C5	5.21	119.86	109.26
4	B	1808	DTB	CD-CE-CR	-4.27	106.26	113.69
3	A	1810	TLA	C3-C2-C1	-3.56	100.54	109.90
3	B	1809	TLA	O1-C1-C2	-2.81	112.94	121.15
3	B	1809	TLA	C2-C3-C4	-2.79	102.56	109.90
3	B	1809	TLA	C3-C2-C1	-2.71	102.78	109.90
3	A	1810	TLA	O4-C4-C3	-2.70	113.25	121.15
4	B	1808	DTB	CR-N2-CN	-2.64	109.31	112.35
3	B	1809	TLA	O4-C4-C3	-2.63	113.44	121.15
3	A	1810	TLA	O1-C1-C2	-2.59	113.58	121.15
3	A	1810	TLA	O2-C2-C1	2.58	116.27	110.73
3	A	1810	TLA	O11-C1-C2	2.46	121.06	113.89
2	A	1644	PLP	O3P-P-O4P	-2.40	100.01	106.65
3	B	1809	TLA	O41-C4-C3	2.39	120.86	113.89
3	B	1809	TLA	O11-C1-C2	2.27	120.50	113.89
4	A	1811	DTB	CS-CR-N2	-2.20	99.63	102.46
3	A	1810	TLA	O41-C4-C3	2.20	120.31	113.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1808	DTB	CS-N1-CN	-2.13	110.77	112.51
4	A	1811	DTB	CT-CS-CR	-2.12	111.10	115.73
2	A	1644	PLP	O3P-P-O1P	2.09	117.26	110.44
4	B	1808	DTB	CT-CS-CR	-2.03	111.29	115.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 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, 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	755/831 (90%)	-0.08	21 (2%)	50 54	15, 31, 68, 94	0
1	B	748/831 (90%)	-0.08	23 (3%)	47 50	14, 30, 68, 94	0
All	All	1503/1662 (90%)	-0.08	44 (2%)	49 53	14, 31, 68, 94	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	492	PHE	4.7
1	B	488	PRO	4.6
1	B	563	LEU	4.2
1	A	451	ASN	3.9
1	A	758	LEU	3.7
1	A	333	HIS	3.7
1	A	493	LEU	3.5
1	A	710	GLN	3.4
1	B	661	ALA	3.3
1	B	806	GLY	3.2
1	B	99	ALA	3.1
1	A	708	THR	3.1
1	A	709	SER	3.0
1	B	559	LEU	2.8
1	A	334	GLN	2.8
1	B	463	VAL	2.8
1	B	96	LEU	2.7
1	B	446	PHE	2.6
1	B	761	LYS	2.6
1	B	759	TYR	2.6
1	B	607	ASN	2.5
1	A	330	LYS	2.5
1	B	807	GLU	2.5
1	A	759	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	354	SER	2.4
1	B	668	GLY	2.3
1	B	730	SER	2.3
1	A	668	GLY	2.3
1	A	450	HIS	2.2
1	B	492	PHE	2.2
1	A	521	LEU	2.2
1	B	575	VAL	2.2
1	B	609	LYS	2.2
1	A	609	LYS	2.2
1	B	802	TYR	2.1
1	B	669	ASP	2.1
1	B	489	TYR	2.1
1	A	431	GLY	2.1
1	B	800	LYS	2.1
1	A	45	SER	2.0
1	A	335	GLU	2.0
1	A	608	ARG	2.0
1	A	332	VAL	2.0
1	B	760	ALA	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 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, 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	DTB	A	1811	15/15	0.19	3.45	5,33,64,71	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	DTB	B	1808	15/15	0.20	2.13	2,23,45,51	0
3	TLA	A	1810	10/10	0.15	1.18	7,28,39,45	0
3	TLA	B	1809	10/10	0.16	0.73	22,37,47,56	0
2	PLP	A	1644	15/16	0.20	0.14	16,31,39,40	0
2	PLP	B	1644	15/16	0.17	-0.31	15,31,36,40	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.