



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2016 – 11:38 AM EST

PDB ID : 5F9S  
Title : Crystal structure of human Alanine:Glyoxylate Aminotransferase major allele (AGT-Ma) at 1.7 Angstrom; internal aldimine with PLP in the active site  
Authors : Giardina, G.; Cutruzzola, F.; Borri Voltattorni, C.; Cellini, B.; Montioli, R.  
Deposited on : 2015-12-10  
Resolution : 1.70 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](mailto: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.

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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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20028442

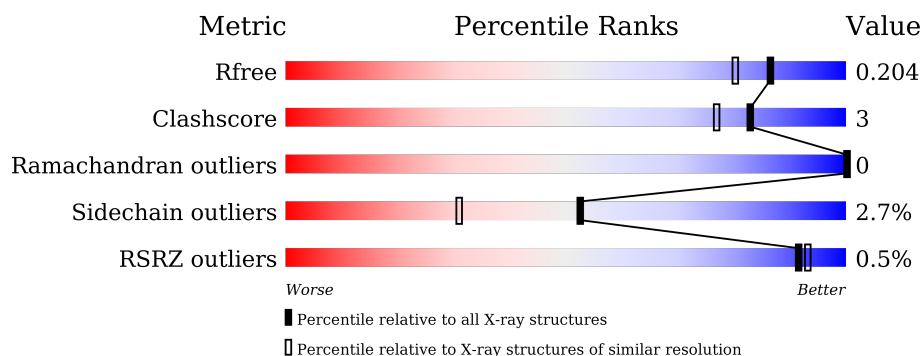
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
$R_{free}$	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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. 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	386	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>%</span> <span>94%</span> <span>5%</span> </div> </div>
1	B	386	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>%</span> <span>91%</span> <span>8% •</span> </div> </div>

## 2 Entry composition [i](#)

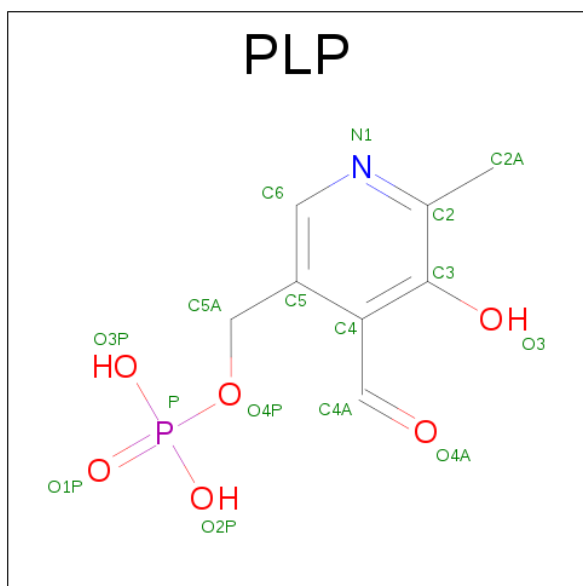
There are 3 unique types of molecules in this entry. The entry contains 13111 atoms, of which 6215 are hydrogens and 0 are deuteriums.

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–pyruvate aminotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	385	Total	C	H	N	O	S	95	13	0
			6123	1936	3096	527	546	18			
1	B	386	Total	C	H	N	O	S	110	14	0
			6138	1939	3101	531	547	20			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C<sub>8</sub>H<sub>10</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			24	8	9	1	5	1		
2	B	1	Total	C	H	N	O	P	0	0
			24	8	9	1	5	1		

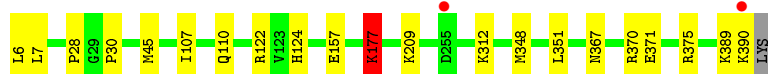
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	417	Total 417	O 417	0	0
3	B	385	Total 385	O 385	0	0

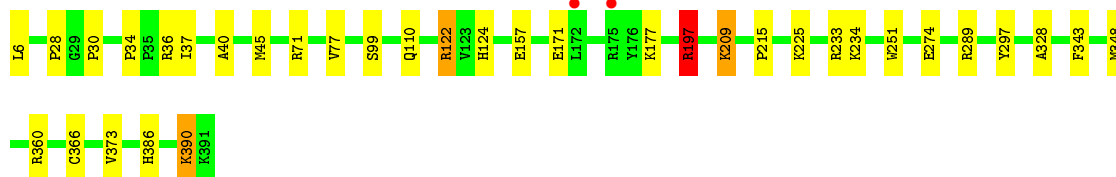
### 3 Residue-property plots [i](#)

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: Serine-pyruvate aminotransferase



- Molecule 1: Serine-pyruvate aminotransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.48Å 101.88Å 131.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.12 – 1.70 45.12 – 1.70	Depositor EDS
% Data completeness (in resolution range)	86.3 (45.12-1.70) 86.3 (45.12-1.70)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.81 (at 1.70Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.158 , 0.205 0.158 , 0.204	Depositor DCC
$R_{free}$ test set	3409 reflections (4.87%)	DCC
Wilson B-factor (Å <sup>2</sup> )	19.8	Xtriage
Anisotropy	0.191	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 45.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	13111	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.55	1/3146 (0.0%)	0.68	2/4267 (0.0%)
1	B	0.50	0/3165	0.67	3/4291 (0.1%)
All	All	0.53	1/6311 (0.0%)	0.67	5/8558 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	177	LYS	CE-NZ	-8.37	1.28	1.49

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	197	ARG	NE-CZ-NH1	7.62	124.11	120.30
1	A	375	ARG	NE-CZ-NH2	6.19	123.40	120.30
1	A	7	LEU	CB-CG-CD1	-5.69	101.33	111.00
1	B	289	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	B	390	LYS	CA-CB-CG	5.00	124.41	113.40

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	3027	3096	3058	14	2
1	B	3037	3101	3050	28	3
2	A	15	9	7	1	0
2	B	15	9	7	1	0
3	A	417	0	0	4	0
3	B	385	0	0	9	1
All	All	6896	6215	6122	38	3

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 3.

All (38) 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:B:348[B]:MET:SD	3:B:681:HOH:O	2.42	0.77
1:A:122[B]:ARG:NH2	3:A:505:HOH:O	2.18	0.77
1:B:328:ALA:O	3:B:501:HOH:O	2.04	0.76
1:A:122[B]:ARG:NH1	1:A:124:HIS:NE2	2.35	0.74
1:B:122[B]:ARG:NH2	3:B:502:HOH:O	2.05	0.71
1:A:6:LEU:N	3:A:506:HOH:O	2.23	0.70
1:B:110:GLN:OE1	3:B:503:HOH:O	2.08	0.70
1:B:386:HIS:O	3:B:505:HOH:O	2.12	0.67
1:A:122[B]:ARG:NH1	1:A:124:HIS:CE1	2.63	0.67
1:B:99:SER:HB2	1:B:122[A]:ARG:NH2	2.10	0.67
1:A:110:GLN:OE1	3:A:503:HOH:O	2.13	0.65
1:B:99:SER:HB2	1:B:122[A]:ARG:CZ	2.27	0.65
1:B:122[A]:ARG:HG2	1:B:122[A]:ARG:HH11	1.61	0.65
1:B:348[A]:MET:SD	3:B:635:HOH:O	2.55	0.63
1:B:197:ARG:HG2	1:B:197:ARG:HH11	1.62	0.63
1:B:348[B]:MET:HE2	3:B:635:HOH:O	1.99	0.62
1:A:30:PRO:HB2	1:B:45[B]:MET:HE2	1.84	0.60
1:B:122[A]:ARG:NH2	3:B:504:HOH:O	2.11	0.59
1:A:28:PRO:HD2	1:A:348:MET:HG3	1.84	0.58
1:A:209:LYS:NZ	2:A:401:PLP:O3	2.36	0.58
1:B:297:TYR:CD2	1:B:373:VAL:HG11	2.40	0.57
1:B:122[A]:ARG:HG3	1:B:122[A]:ARG:O	2.03	0.57
1:A:45[A]:MET:HE2	1:B:215:PRO:HD3	1.87	0.56
1:A:45[A]:MET:HE3	1:B:30:PRO:CG	2.36	0.56
1:A:30:PRO:CB	1:B:45[B]:MET:HE2	2.35	0.56
1:A:367[A]:ASN:OD1	3:A:504:HOH:O	2.15	0.55
1:B:209:LYS:NZ	2:B:401:PLP:O3	2.42	0.53
1:B:225:LYS:HG3	3:B:718:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122[A]:ARG:NH1	1:B:122[A]:ARG:HG2	2.25	0.51
1:A:107:ILE:HD13	1:A:351:LEU:HG	1.93	0.51
1:B:40:ALA:HB2	1:B:274:GLU:HG3	1.96	0.48
1:B:77:VAL:HB	1:B:251:TRP:CZ2	2.52	0.45
1:A:389:LYS:O	1:A:390:LYS:C	2.55	0.44
1:B:197:ARG:HG2	1:B:197:ARG:NH1	2.32	0.43
1:B:122[A]:ARG:HD2	1:B:124:HIS:NE2	2.34	0.43
1:B:28:PRO:HG3	1:B:360:ARG:CZ	2.49	0.42
1:B:34:PRO:HD2	1:B:37:ILE:HD12	2.01	0.42
1:B:28:PRO:HD2	1:B:348[A]:MET:SD	2.60	0.41

All (3) 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)
1:B:366[B]:CYS:HG	3:B:828:HOH:O[3_655]	1.12	0.48
1:A:177:LYS:NZ	1:B:343:PHE:O[3_545]	2.00	0.20
1:A:177:LYS:HZ1	1:B:343:PHE:O[3_545]	1.41	0.19

## 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	396/386 (103%)	390 (98%)	6 (2%)	0	100	100
1	B	398/386 (103%)	389 (98%)	9 (2%)	0	100	100
All	All	794/772 (103%)	779 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	332/321 (103%)	327 (98%)	5 (2%)	72	56
1	B	335/321 (104%)	321 (96%)	14 (4%)	36	14
All	All	667/642 (104%)	648 (97%)	19 (3%)	52	29

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	157	GLU
1	A	177	LYS
1	A	312	LYS
1	A	370	ARG
1	A	371	GLU
1	B	6	LEU
1	B	36[A]	ARG
1	B	36[B]	ARG
1	B	71	ARG
1	B	122[A]	ARG
1	B	122[B]	ARG
1	B	157	GLU
1	B	171	GLU
1	B	177	LYS
1	B	197	ARG
1	B	209	LYS
1	B	233	ARG
1	B	234	LYS
1	B	390	LYS

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 molecules 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 ⓘ

2 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	401	1	15,15,16	2.04	4 (26%)	21,22,23	2.29	8 (38%)
2	PLP	B	401	1	15,15,16	2.26	6 (40%)	21,22,23	1.63	4 (19%)

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	401	1	-	0/6/6/8	0/1/1/1
2	PLP	B	401	1	-	0/6/6/8	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	PLP	C5-C4	-2.69	1.37	1.40
2	A	401	PLP	O4P-C5A	-2.30	1.36	1.45
2	B	401	PLP	C5A-C5	2.08	1.56	1.51
2	A	401	PLP	C2A-C2	2.10	1.54	1.50
2	B	401	PLP	C6-N1	2.11	1.38	1.34
2	B	401	PLP	C3-C2	2.36	1.42	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	PLP	O3-C3	2.61	1.43	1.37
2	B	401	PLP	C2A-C2	2.69	1.55	1.50
2	A	401	PLP	C4A-C4	5.34	1.62	1.51
2	B	401	PLP	C4A-C4	5.92	1.63	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PLP	C2A-C2-C3	-3.51	117.34	120.90
2	A	401	PLP	C5-C6-N1	-2.77	119.02	123.86
2	A	401	PLP	C4A-C4-C3	-2.67	115.76	120.42
2	A	401	PLP	C5A-C5-C6	-2.48	114.69	119.33
2	B	401	PLP	C2A-C2-C3	-2.39	118.48	120.90
2	B	401	PLP	O4P-P-O1P	-2.22	101.50	107.08
2	A	401	PLP	O2P-P-O4P	2.18	113.08	106.72
2	A	401	PLP	C3-C2-N1	2.47	123.78	120.69
2	B	401	PLP	C6-C5-C4	2.47	120.23	118.12
2	B	401	PLP	C4A-C4-C5	4.18	125.20	120.90
2	A	401	PLP	C6-C5-C4	4.77	122.17	118.12
2	A	401	PLP	C4A-C4-C5	5.57	126.64	120.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PLP	1	0
2	B	401	PLP	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<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	385/386 (99%)	-0.22	2 (0%) 91 93	13, 19, 32, 66	12 (3%)
1	B	386/386 (100%)	-0.03	2 (0%) 91 93	15, 21, 33, 47	12 (3%)
All	All	771/772 (99%)	-0.12	4 (0%) 91 93	13, 20, 32, 66	24 (3%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	390	LYS	5.4
1	A	255	ASP	2.4
1	B	172	LEU	2.1
1	B	175	ARG	2.1

### 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](#)

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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	PLP	A	401	15/16	0.97	0.13	1.51	12,16,28,28	0
2	PLP	B	401	15/16	0.98	0.10	-0.00	16,22,28,28	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.