



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:45 PM GMT

PDB ID : 1QJ5  
Title : CRYSTAL STRUCTURE OF 7,8-DIAMINOPELARGONIC ACID SYNTHASE  
Authors : Kack, H.; Sandmark, J.; Gibson, K.J.; Lindqvist, Y.; Schneider, G.  
Deposited on : 1999-06-21  
Resolution : 1.80 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

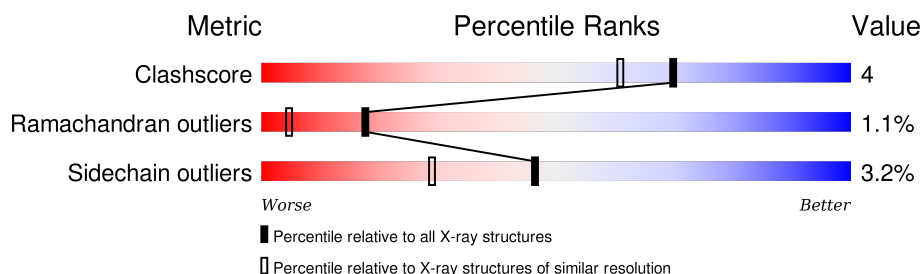
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	429	 84% 13% •
1	B	429	 86% 12% •

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7227 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 7,8-DIAMINOPELARGONIC ACID SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	6	0
			3324	2108	576	607	33			
1	B	429	Total	C	N	O	S	0	3	0
			3315	2100	577	605	33			

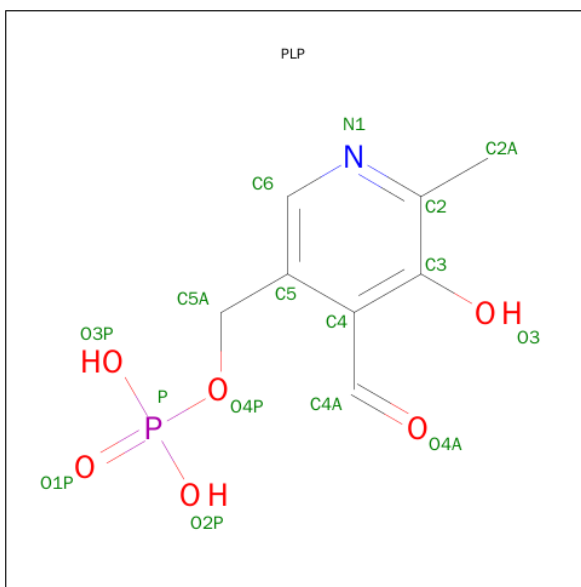
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	LEU	TRP	CONFLICT	UNP P12995
B	14	LEU	TRP	CONFLICT	UNP P12995

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	K	0	0
			1	1		
2	A	1	Total	K	0	0
			1	1		

- Molecule 3 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
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 4 is water.

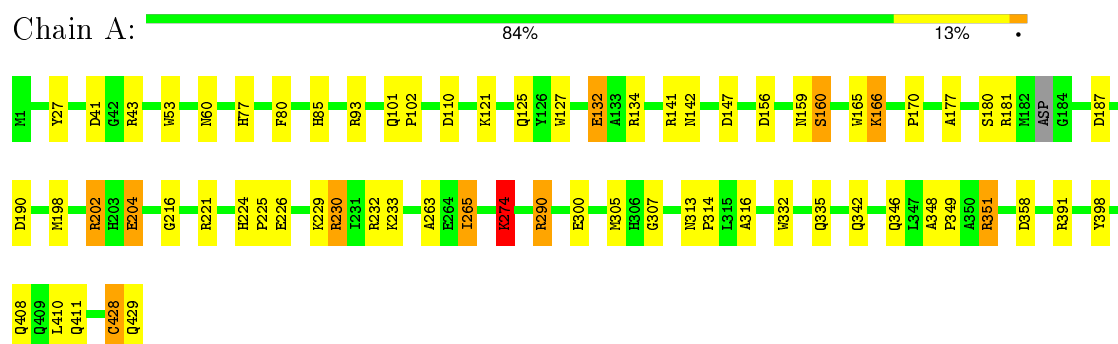
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	277	Total	O	0	0
			277	277		
4	B	279	Total	O	0	0
			279	279		

### 3 Residue-property plots

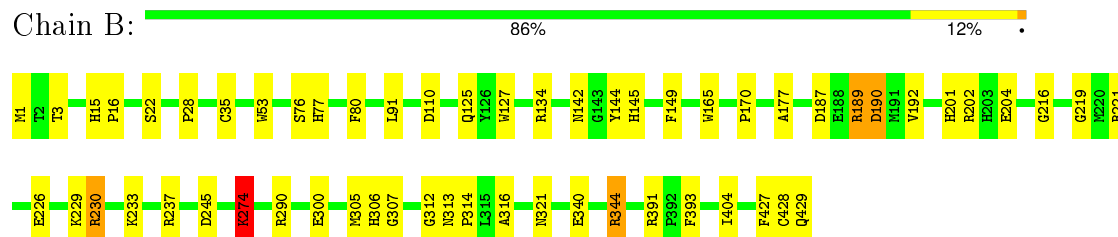
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

#### • Molecule 1: 7,8-DIAMINOPELARGONIC ACID SYNTHASE



#### • Molecule 1: 7,8-DIAMINOPELARGONIC ACID SYNTHASE



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.34 Å   55.56 Å   120.88 Å 90.00°   96.97°   90.00°	Depositor
Resolution (Å)	20.00 – 1.80	Depositor
% Data completeness (in resolution range)	94.4 (20.00-1.80)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.175 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7227	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K, 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	0.51	1/3431 (0.0%)	1.20	20/4653 (0.4%)
1	B	0.52	1/3406 (0.0%)	1.31	16/4623 (0.3%)
All	All	0.52	2/6837 (0.0%)	1.26	36/9276 (0.4%)

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	A	0	1
1	B	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	274	LYS	C-N	5.99	1.47	1.34
1	A	274	LYS	C-N	5.85	1.47	1.34

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	290	ARG	CD-NE-CZ	35.12	172.76	123.60
1	B	274	LYS	O-C-N	-17.45	94.79	122.70
1	A	274	LYS	O-C-N	-16.77	95.87	122.70
1	A	274	LYS	CA-C-O	16.00	153.70	120.10
1	B	274	LYS	CA-C-O	15.97	153.65	120.10
1	B	344	ARG	CD-NE-CZ	11.48	139.67	123.60
1	B	391	ARG	NE-CZ-NH2	-10.87	114.86	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	221	ARG	NE-CZ-NH2	-9.32	115.64	120.30
1	A	290	ARG	NE-CZ-NH2	-8.28	116.16	120.30
1	B	230	ARG	NE-CZ-NH2	7.98	124.29	120.30
1	A	141	ARG	NE-CZ-NH1	7.87	124.23	120.30
1	A	391	ARG	NE-CZ-NH2	-7.80	116.40	120.30
1	A	274	LYS	CD-CE-NZ	7.32	128.54	111.70
1	A	351	ARG	CA-C-N	7.28	133.21	117.20
1	B	221	ARG	NE-CZ-NH1	7.07	123.83	120.30
1	A	274	LYS	CA-CB-CG	6.63	127.99	113.40
1	B	144	TYR	CB-CG-CD1	-6.53	117.08	121.00
1	A	230	ARG	NE-CZ-NH2	6.23	123.41	120.30
1	B	427	PHE	C-N-CA	6.15	137.08	121.70
1	B	290	ARG	NE-CZ-NH1	-6.03	117.28	120.30
1	B	202	ARG	NE-CZ-NH1	5.91	123.25	120.30
1	A	141	ARG	CD-NE-CZ	5.86	131.81	123.60
1	A	274	LYS	CA-C-N	-5.86	104.31	117.20
1	A	27	TYR	CB-CG-CD1	-5.83	117.50	121.00
1	A	221	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	B	344	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	274	LYS	CB-CA-C	5.65	121.69	110.40
1	A	204	GLU	N-CA-CB	-5.53	100.64	110.60
1	A	93	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	A	358	ASP	CB-CG-OD2	5.24	123.02	118.30
1	A	398	TYR	CB-CG-CD2	-5.16	117.90	121.00
1	A	391	ARG	NH1-CZ-NH2	5.16	125.07	119.40
1	B	274	LYS	CA-C-N	-5.09	105.99	117.20
1	A	181	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	B	22	SER	N-CA-CB	-5.07	102.90	110.50
1	B	230	ARG	NE-CZ-NH1	-5.07	117.77	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	274	LYS	Mainchain
1	B	274	LYS	Mainchain

## 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	3324	0	3275	35	1
1	B	3315	0	3262	29	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	15	0	6	0	0
3	B	15	0	6	0	0
4	A	277	0	0	5	0
4	B	279	0	0	3	0
All	All	7227	0	6549	59	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 4.

All (59) 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:180:SER:OG	1:A:190[A]:ASP:OD2	1.88	0.90
1:B:125:GLN:HE22	1:B:305:MET:H	1.25	0.82
1:A:125:GLN:HE22	1:A:305:MET:H	1.36	0.73
1:A:121[B]:LYS:HG3	1:B:165:TRP:CH2	2.26	0.71
1:A:428:CYS:O	1:A:429:GLN:HB2	1.92	0.69
1:A:142:ASN:HD22	1:A:177:ALA:HB2	1.61	0.66
1:B:142:ASN:HD22	1:B:177:ALA:HB2	1.62	0.64
1:A:132[A]:GLU:CD	1:A:134:ARG:HE	2.02	0.63
1:B:226:GLU:OE2	1:B:230:ARG:NH1	2.32	0.62
1:A:226:GLU:OE2	1:A:230:ARG:NH1	2.34	0.61
1:B:1:MET:HG2	4:B:2013:HOH:O	2.03	0.58
1:B:77:HIS:HA	1:B:314:PRO:HD2	1.87	0.56
1:B:145:HIS:HD2	1:B:245:ASP:OD2	1.88	0.56
1:A:77:HIS:HA	1:A:314:PRO:HD2	1.86	0.56
1:B:149:PHE:CZ	1:B:170:PRO:HD3	2.43	0.53
1:A:156:ASP:O	1:A:160:SER:HB2	2.08	0.53
1:A:101:GLN:HB3	1:A:102:PRO:HD3	1.93	0.51
1:B:189:ARG:O	1:B:192:VAL:HG13	2.10	0.50
1:B:127:TRP:CD2	1:B:134:ARG:HD2	2.47	0.50
1:B:125:GLN:NE2	1:B:305:MET:H	2.02	0.49
1:B:145:HIS:HE1	4:B:2125:HOH:O	1.95	0.49
1:A:229:LYS:HG3	1:A:263:ALA:HB1	1.94	0.49
1:A:160:SER:HB3	4:A:2135:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132[A]:GLU:OE2	1:A:134:ARG:NE	2.30	0.49
1:A:232:ARG:HG2	1:A:265:ILE:HG13	1.95	0.48
1:A:313:ASN:ND2	1:A:316:ALA:H	2.12	0.48
1:A:300:GLU:HG3	4:A:2116:HOH:O	2.14	0.48
1:B:340:GLU:O	1:B:344:ARG:HG3	2.14	0.48
1:A:121[B]:LYS:CG	1:B:165:TRP:CH2	2.97	0.47
1:A:156:ASP:HB3	1:A:159:ASN:HB2	1.96	0.47
1:A:342:GLN:OE1	1:A:411:GLN:HG3	2.15	0.47
1:A:127:TRP:CD2	1:A:134:ARG:HD2	2.49	0.46
1:A:41:ASP:OD2	1:A:43:ARG:NE	2.46	0.46
1:B:15:HIS:HB3	1:B:16:PRO:HD2	1.98	0.46
1:A:204:GLU:HG2	4:A:2162:HOH:O	2.15	0.46
1:A:170:PRO:HD3	1:B:149:PHE:CZ	2.51	0.46
1:A:147:ASP:HB3	1:B:306:HIS:CE1	2.51	0.45
1:A:332:TRP:HA	1:A:335:GLN:HE21	1.80	0.45
1:B:91:LEU:HA	1:B:321[B]:ASN:OD1	2.17	0.45
1:B:53:TRP:CD1	1:B:274:LYS:HE2	2.52	0.45
1:A:53:TRP:CD1	1:A:274:LYS:HE2	2.52	0.44
1:A:224:HIS:HA	1:A:225:PRO:HD3	1.86	0.44
1:A:166:LYS:HE2	4:A:2138:HOH:O	2.17	0.44
1:B:1:MET:HE3	1:B:28:PRO:HB2	1.99	0.44
1:A:187:ASP:HB3	1:A:190[A]:ASP:OD1	2.18	0.43
1:A:165:TRP:CH2	1:B:125:GLN:HG3	2.53	0.43
1:B:229:LYS:HB2	1:B:229:LYS:HE2	1.88	0.42
1:A:346:GLN:NE2	1:A:411:GLN:HG2	2.34	0.42
1:A:346:GLN:HE22	1:A:411:GLN:CG	2.33	0.42
1:B:187:ASP:O	1:B:190:ASP:HB2	2.20	0.42
1:A:85:HIS:HE1	4:A:2086:HOH:O	2.02	0.41
1:B:312:GLY:O	1:B:313:ASN:C	2.58	0.41
1:B:201:HIS:HD2	1:B:204:GLU:OE2	2.04	0.41
1:B:313:ASN:ND2	1:B:316:ALA:H	2.19	0.41
1:B:300:GLU:HG3	4:B:2122:HOH:O	2.20	0.41
1:B:35[A]:CYS:SG	1:B:404:ILE:HG13	2.61	0.40
1:B:429:GLN:HG2	1:B:429:GLN:OXT	2.21	0.40
1:A:348:ALA:N	1:A:349:PRO:CD	2.84	0.40
1:A:198:MET:O	1:A:202:ARG:HB3	2.21	0.40

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)
1:A:351:ARG:O	1:A:408:GLN:NE2[2_645]	2.07	0.13

## 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	430/429 (100%)	417 (97%)	9 (2%)	4 (1%)	21	7
1	B	430/429 (100%)	416 (97%)	9 (2%)	5 (1%)	16	4
All	All	860/858 (100%)	833 (97%)	18 (2%)	9 (1%)	17	5

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	428	CYS
1	A	274	LYS
1	B	274	LYS
1	B	216	GLY
1	A	428	CYS
1	A	216	GLY
1	B	307	GLY
1	A	307	GLY
1	B	219	GLY

### 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	350/345 (101%)	336 (96%)	14 (4%)	38	20
1	B	347/345 (101%)	336 (97%)	11 (3%)	46	29
All	All	697/690 (101%)	672 (96%)	25 (4%)	46	24

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	80	PHE
1	A	110[A]	ASP
1	A	110[B]	ASP
1	A	132[A]	GLU
1	A	132[B]	GLU
1	A	160	SER
1	A	166	LYS
1	A	202	ARG
1	A	233	LYS
1	A	265	ILE
1	A	274	LYS
1	A	290	ARG
1	A	410	LEU
1	B	3	THR
1	B	76	SER
1	B	80	PHE
1	B	110[A]	ASP
1	B	110[B]	ASP
1	B	189	ARG
1	B	190	ASP
1	B	233	LYS
1	B	237	ARG
1	B	274	LYS
1	B	393	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	HIS
1	A	125	GLN
1	A	135	GLN
1	A	142	ASN
1	A	313	ASN

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Mol	Chain	Res	Type
1	A	335	GLN
1	B	63	GLN
1	B	125	GLN
1	B	135	GLN
1	B	142	ASN
1	B	145	HIS
1	B	201	HIS
1	B	262	HIS
1	B	313	ASN
1	B	335	GLN
1	B	342	GLN
1	B	346	GLN

### 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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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$
3	PLP	A	502	1	15,15,16	1.13	1 (6%)	21,22,23	2.15	7 (33%)
3	PLP	B	502	1	15,15,16	1.23	2 (13%)	21,22,23	1.85	7 (33%)

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
3	PLP	A	502	1	-	0/6/6/8	0/1/1/1
3	PLP	B	502	1	-	0/6/6/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	PLP	C5-C4	-2.79	1.37	1.40
3	A	502	PLP	O3-C3	-2.49	1.31	1.37
3	B	502	PLP	O3-C3	-2.45	1.31	1.37

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	PLP	C5-C6-N1	-3.50	117.78	123.86
3	A	502	PLP	C3-C2-N1	-3.24	116.14	120.61
3	B	502	PLP	C5-C6-N1	-2.65	119.26	123.86
3	A	502	PLP	C4A-C4-C5	-2.61	118.16	120.88
3	B	502	PLP	C3-C2-N1	-2.17	117.62	120.61
3	B	502	PLP	O3P-P-O1P	2.06	117.20	110.58
3	B	502	PLP	C5A-C5-C4	2.17	124.52	121.65
3	A	502	PLP	C3-C4-C5	2.33	121.32	118.78
3	B	502	PLP	O2P-P-O4P	2.49	113.73	106.56
3	B	502	PLP	O4P-C5A-C5	2.92	113.82	108.99
3	A	502	PLP	O4P-C5A-C5	3.68	115.07	108.99
3	B	502	PLP	C4A-C4-C5	3.71	124.75	120.88
3	A	502	PLP	C2A-C2-C3	3.71	125.52	121.04
3	A	502	PLP	C6-N1-C2	4.09	127.62	119.28

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.