



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2025 – 11:30 AM EDT

PDB ID : 1XL0  
Title : Kinetic and crystallographic studies on 2-(beta-D-glucopyranosyl)-5-methyl-1,3,4-oxadiazole,-benzothiazole, and-benzimidazole, inhibitors of muscle glycogen phosphorylase b. Evidence for a new binding site.  
Authors : Chrysina, E.D.; Kosmopoulou, M.N.; Tiraidis, C.; Kardakaris, R.; Bischler, N.; Leonidas, D.D.; Hadady, Z.; Somsak, L.; Docsa, T.; Gergely, P.; Oikonomakos, N.G.  
Deposited on : 2004-09-30  
Resolution : 1.92 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

<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>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

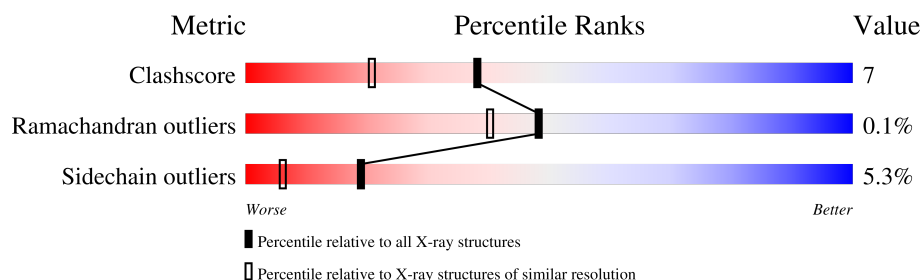
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.92 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	180529	1100 (1.92-1.92)
Ramachandran outliers	177936	1087 (1.92-1.92)
Sidechain outliers	177891	1087 (1.92-1.92)

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  $\geq 3$ , 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	842	 80% 14% • •

## 2 Entry composition [i](#)

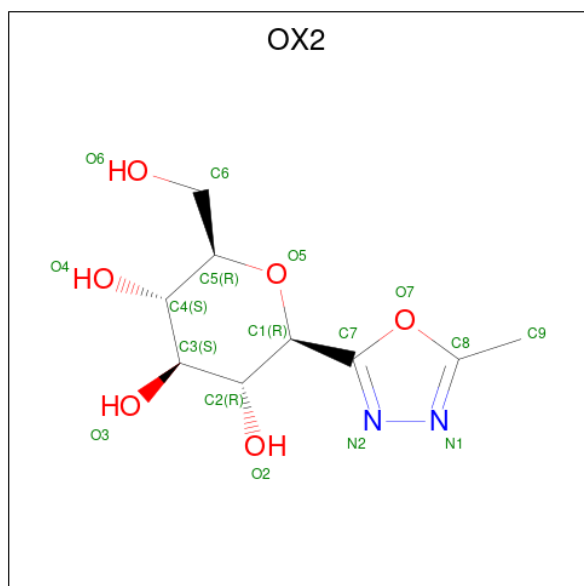
There are 4 unique types of molecules in this entry. The entry contains 6989 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycogen phosphorylase, muscle form.

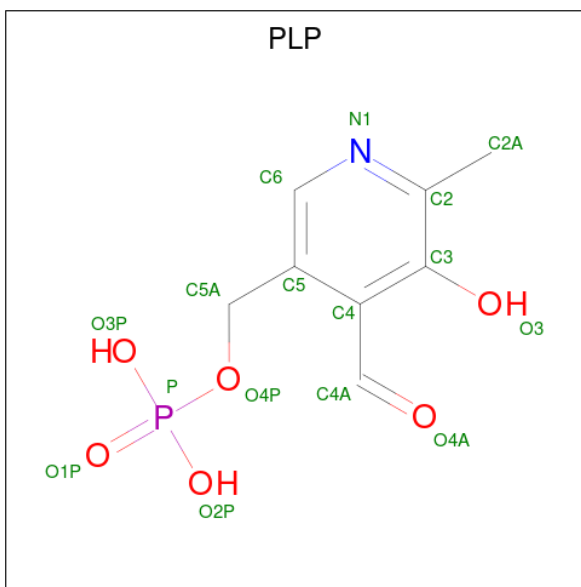
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	810	6589	4201	1160	1199	29	0	0	0

- Molecule 2 is (1R)-1,5-anhydro-1-(5-methyl-1,3,4-oxadiazol-2-yl)-D-glucitol (three-letter code: OX2) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>2</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	17	9	2	6	0	0

- 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		

- Molecule 4 is water.

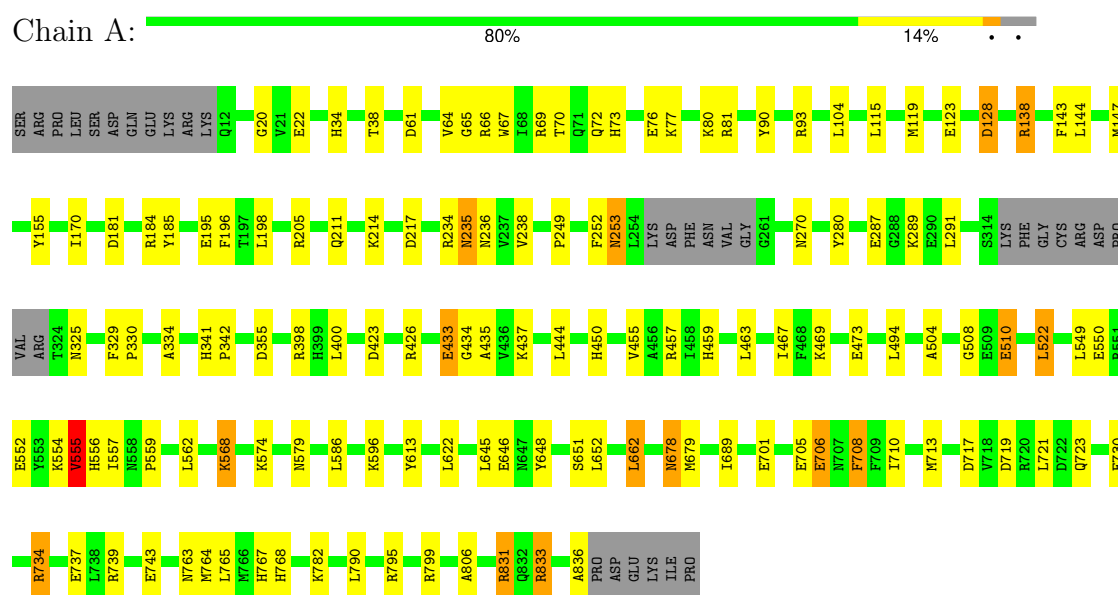
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	368	Total	O	0	0
			368	368		

### 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. 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. 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: Glycogen phosphorylase, muscle form



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	128.13Å 128.13Å 115.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.32 – 1.92	Depositor
% Data completeness (in resolution range)	98.8 (29.32-1.92)	Depositor
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.182 , 0.200	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6989	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

## 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, OX2

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	0.31	0/6736	0.59	0/9116

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 [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	6589	0	6533	89	0
2	A	17	0	14	0	0
3	A	15	0	7	0	0
4	A	368	0	0	14	0
All	All	6989	0	6554	89	0

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 7.

All (89) 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:270:ASN:HB2	4:A:1346:HOH:O	1.68	0.94
1:A:455:VAL:H	1:A:459:HIS:HD2	1.27	0.83
1:A:138:ARG:O	1:A:138:ARG:HD3	1.80	0.82
1:A:708:PHE:HB3	1:A:710:ILE:CD1	2.12	0.77
1:A:708:PHE:HB3	1:A:710:ILE:HD11	1.65	0.77
1:A:678:ASN:HD22	1:A:679:MET:H	1.34	0.73
1:A:433:GLU:OE1	1:A:437:LYS:HD2	1.91	0.70
1:A:61:ASP:O	1:A:64:VAL:HG22	1.97	0.65
1:A:355:ASP:OD1	1:A:398:ARG:HD3	1.97	0.64
1:A:65:GLY:O	1:A:69:ARG:HG2	1.97	0.64
1:A:329:PHE:HB3	1:A:330:PRO:HD3	1.81	0.62
1:A:235:ASN:H	1:A:235:ASN:HD22	1.46	0.62
1:A:34:HIS:HD2	1:A:38:THR:OG1	1.84	0.61
1:A:423:ASP:O	1:A:426:ARG:HG3	2.02	0.59
1:A:66:ARG:HG3	4:A:1333:HOH:O	2.02	0.59
1:A:730:GLU:O	1:A:734:ARG:HG3	2.03	0.58
1:A:739:ARG:O	1:A:743:GLU:HG3	2.04	0.58
1:A:236:ASN:CG	1:A:836:ALA:HB2	2.23	0.58
1:A:713:MET:HB3	1:A:717:ASP:HB2	1.86	0.57
1:A:678:ASN:HD22	1:A:679:MET:N	2.02	0.56
1:A:678:ASN:ND2	1:A:679:MET:H	2.02	0.56
1:A:550:GLU:O	1:A:554:LYS:HD3	2.05	0.56
1:A:549:LEU:HD12	1:A:557:ILE:HD13	1.87	0.56
1:A:795:ARG:O	1:A:799:ARG:HG3	2.06	0.56
1:A:34:HIS:HE1	1:A:61:ASP:OD1	1.90	0.55
1:A:69:ARG:HG3	4:A:1336:HOH:O	2.06	0.55
1:A:423:ASP:HB2	4:A:1229:HOH:O	2.06	0.55
1:A:708:PHE:HB3	1:A:710:ILE:HD12	1.88	0.55
1:A:235:ASN:HA	1:A:833:ARG:HG3	1.90	0.54
1:A:510:GLU:HB2	4:A:1179:HOH:O	2.07	0.54
1:A:184:ARG:HE	1:A:185:TYR:HE2	1.56	0.53
1:A:556:HIS:CD2	1:A:556:HIS:N	2.76	0.52
1:A:705:GLU:HG3	4:A:1029:HOH:O	2.10	0.51
1:A:128:ASP:OD1	1:A:651:SER:HB3	2.10	0.51
1:A:144:LEU:HD23	1:A:147:MET:CE	2.40	0.51
1:A:170:ILE:HG12	1:A:646:GLU:HG3	1.93	0.51
1:A:568:LYS:HG3	1:A:574:LYS:HD3	1.92	0.51
1:A:20:GLY:HA3	4:A:1329:HOH:O	2.10	0.50
1:A:138:ARG:HD3	1:A:138:ARG:C	2.31	0.50
1:A:469:LYS:O	1:A:473:GLU:HG3	2.12	0.50
1:A:450:HIS:HE1	4:A:1154:HOH:O	1.95	0.50
1:A:678:ASN:HD22	1:A:678:ASN:N	2.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:ARG:HH12	1:A:701:GLU:CD	2.15	0.49
1:A:522:LEU:HD13	1:A:806:ALA:CB	2.43	0.48
1:A:450:HIS:HD2	4:A:1267:HOH:O	1.96	0.48
1:A:554:LYS:O	1:A:555:VAL:HG13	2.14	0.48
1:A:235:ASN:H	1:A:235:ASN:ND2	2.12	0.47
1:A:763:ASN:OD1	1:A:767:HIS:HD2	1.98	0.47
1:A:144:LEU:HD23	1:A:147:MET:HE3	1.97	0.46
1:A:455:VAL:H	1:A:459:HIS:CD2	2.18	0.46
1:A:710:ILE:HD12	1:A:710:ILE:N	2.31	0.46
1:A:494:LEU:C	1:A:494:LEU:HD23	2.36	0.45
1:A:730:GLU:O	1:A:734:ARG:CG	2.64	0.45
1:A:138:ARG:HB3	4:A:1135:HOH:O	2.15	0.45
1:A:831:ARG:O	1:A:833:ARG:NH1	2.49	0.45
1:A:253:ASN:ND2	4:A:1252:HOH:O	2.50	0.44
1:A:557:ILE:O	1:A:559:PRO:HD3	2.17	0.44
1:A:562:LEU:HD21	1:A:662:LEU:HB2	2.00	0.44
1:A:81:ARG:HG3	1:A:155:TYR:HE2	1.82	0.44
1:A:181:ASP:HB3	1:A:184:ARG:NH2	2.33	0.44
1:A:341:HIS:HB2	1:A:342:PRO:HD3	1.99	0.44
1:A:737:GLU:HG3	4:A:1173:HOH:O	2.18	0.44
1:A:706:GLU:H	1:A:706:GLU:CD	2.21	0.44
1:A:249:PRO:O	1:A:252:PHE:HB2	2.18	0.43
1:A:434:GLY:O	1:A:435:ALA:C	2.57	0.43
1:A:504:ALA:HA	1:A:508:GLY:O	2.19	0.43
1:A:64:VAL:HG23	1:A:65:GLY:N	2.34	0.42
1:A:67:TRP:HA	1:A:238:VAL:HB	2.01	0.42
1:A:93:ARG:HD2	4:A:1224:HOH:O	2.19	0.42
1:A:678:ASN:ND2	1:A:679:MET:N	2.65	0.42
1:A:214:LYS:NZ	1:A:214:LYS:HB2	2.35	0.42
1:A:184:ARG:NE	1:A:185:TYR:CE2	2.81	0.42
1:A:287:GLU:HG2	1:A:289:LYS:HG2	2.02	0.42
1:A:734:ARG:HD3	4:A:1364:HOH:O	2.19	0.42
1:A:719:ASP:O	1:A:723:GLN:HG3	2.19	0.42
1:A:80:LYS:HE2	1:A:334:ALA:HB2	2.01	0.42
1:A:556:HIS:CD2	1:A:556:HIS:H	2.38	0.42
1:A:764:MET:HA	1:A:768:HIS:CE1	2.54	0.41
1:A:648:TYR:HA	1:A:652:LEU:HD23	2.01	0.41
1:A:689:ILE:HG23	1:A:689:ILE:O	2.20	0.41
1:A:463:LEU:CD2	1:A:467:ILE:HD11	2.50	0.41
1:A:72:GLN:O	1:A:76:GLU:HG2	2.21	0.41
1:A:195:GLU:HG2	1:A:196:PHE:CD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:TYR:OH	1:A:291:LEU:HB3	2.20	0.41
1:A:119:MET:O	1:A:123:GLU:HG3	2.21	0.41
1:A:143:PHE:O	1:A:147:MET:HG3	2.20	0.41
1:A:70:THR:O	1:A:73:HIS:HB3	2.21	0.40
1:A:73:HIS:CE1	1:A:77:LYS:HD2	2.57	0.40
1:A:205:ARG:NE	1:A:217:ASP:OD2	2.55	0.40

There are no symmetry-related clashes.

## 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	804/842 (96%)	772 (96%)	31 (4%)	1 (0%)	48 40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	555	VAL

### 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	701/731 (96%)	664 (95%)	37 (5%)	19 6

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	90	TYR
1	A	104	LEU
1	A	115	LEU
1	A	128	ASP
1	A	138	ARG
1	A	198	LEU
1	A	211	GLN
1	A	234	ARG
1	A	235	ASN
1	A	253	ASN
1	A	325	ASN
1	A	400	LEU
1	A	433	GLU
1	A	444	LEU
1	A	510	GLU
1	A	522	LEU
1	A	552	GLU
1	A	555	VAL
1	A	568	LYS
1	A	579	ASN
1	A	586	LEU
1	A	596	LYS
1	A	613	TYR
1	A	622	LEU
1	A	645	LEU
1	A	662	LEU
1	A	678	ASN
1	A	706	GLU
1	A	708	PHE
1	A	721	LEU
1	A	734	ARG
1	A	765	LEU
1	A	782	LYS
1	A	790	LEU
1	A	831	ARG
1	A	833	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	73	HIS
1	A	211	GLN
1	A	235	ASN
1	A	264	GLN
1	A	270	ASN
1	A	325	ASN
1	A	412	ASN
1	A	450	HIS
1	A	459	HIS
1	A	481	ASN
1	A	484	ASN
1	A	539	GLN
1	A	556	HIS
1	A	560	ASN
1	A	566	GLN
1	A	579	ASN
1	A	678	ASN
1	A	767	HIS
1	A	832	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

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	OX2	A	998	-	14,18,18	1.24	1 (7%)	18,26,26	0.99	1 (5%)
3	PLP	A	999	1	15,15,16	1.49	4 (26%)	21,22,23	1.17	3 (14%)

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	OX2	A	998	-	-	0/2/26/26	0/2/2/2
3	PLP	A	999	1	-	2/6/6/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	998	OX2	O5-C1	2.68	1.47	1.43
3	A	999	PLP	C4A-C4	2.63	1.56	1.51
3	A	999	PLP	C3-C2	-2.44	1.38	1.41
3	A	999	PLP	C2A-C2	2.43	1.54	1.50
3	A	999	PLP	C5A-C5	2.03	1.56	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	998	OX2	C9-C8-N1	3.09	126.43	119.65
3	A	999	PLP	O3P-P-O1P	2.39	120.13	110.83
3	A	999	PLP	C5-C6-N1	-2.12	120.38	123.83
3	A	999	PLP	C6-C5-C4	2.11	119.83	118.10

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	999	PLP	C4-C5-C5A-O4P
3	A	999	PLP	C6-C5-C5A-O4P

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

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.