



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2025 – 08:53 AM EDT

PDB ID : 3VSA
Title : Crystal Structure of O-phosphoserine sulphydrylase without acetate
Authors : Nakamura, T.; Kawai, Y.; Kataoka, M.; Ishikawa, K.
Deposited on : 2012-04-24
Resolution : 2.07 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

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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

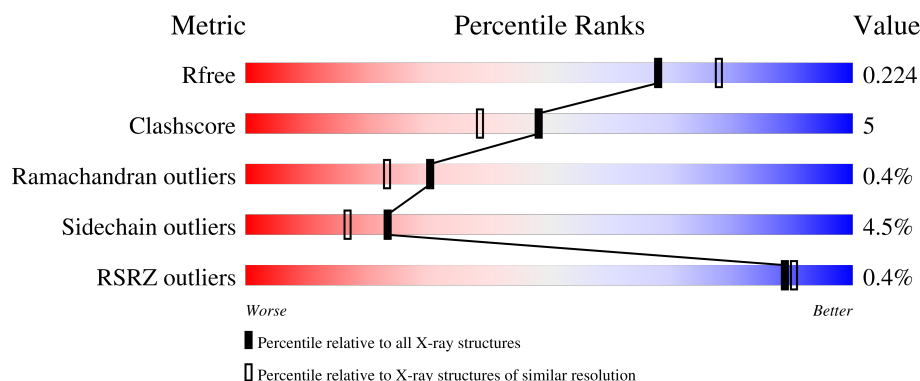
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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}	164625	3436 (2.08-2.04)
Clashscore	180529	3661 (2.08-2.04)
Ramachandran outliers	177936	3649 (2.08-2.04)
Sidechain outliers	177891	3649 (2.08-2.04)
RSRZ outliers	164620	3436 (2.08-2.04)

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	389	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 86% 11% ... </div> </div>
1	B	389	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 81% 15% ... </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6300 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 Protein CysO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	8	0
			2998	1896	537	556	9			
1	B	383	Total	C	N	O	S	0	7	0
			2988	1891	530	558	9			

- 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 (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	6	2		

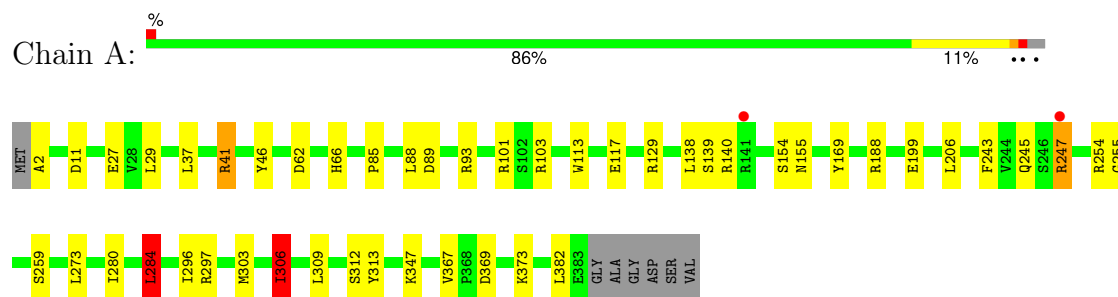
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	139	Total	O	0	0
			139	139		
4	B	137	Total	O	0	0
			137	137		

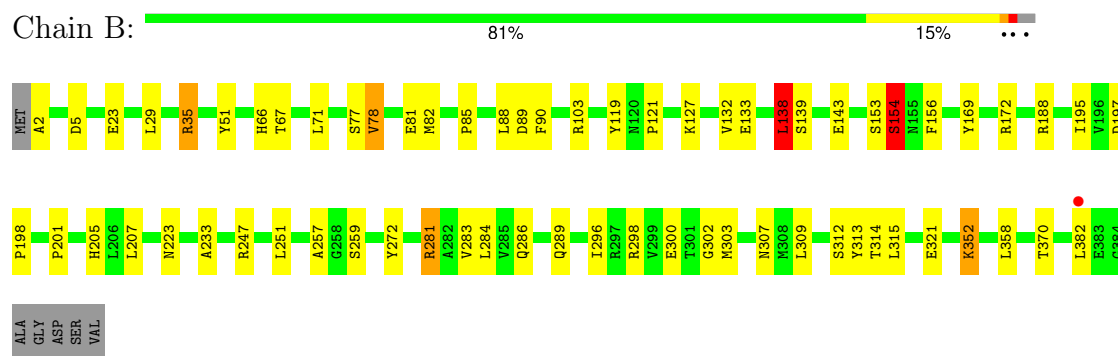
3 Residue-property plots [i](#)

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: Protein CysO



• Molecule 1: Protein CysO



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	74.51Å 74.51Å 275.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.14 – 2.07 39.14 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.3 (39.14-2.07) 99.3 (39.14-2.07)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.61 (at 2.06Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.167 , 0.224 0.168 , 0.224	Depositor DCC
R_{free} test set	2427 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6300	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% of the height of the origin peak. No significant pseudotranslation is detected.*

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: MPD, 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	1.05	2/3057 (0.1%)	0.98	10/4144 (0.2%)
1	B	1.09	4/3048 (0.1%)	0.96	2/4134 (0.0%)
All	All	1.07	6/6105 (0.1%)	0.97	12/8278 (0.1%)

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	272	TYR	CD2-CE2	6.84	1.49	1.39
1	B	51	TYR	CD2-CE2	6.22	1.48	1.39
1	B	132	VAL	CB-CG1	5.34	1.64	1.52
1	A	243	PHE	CE2-CZ	5.21	1.47	1.37
1	B	119	TYR	CE2-CZ	5.19	1.45	1.38
1	A	27	GLU	CG-CD	5.17	1.59	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ARG	NE-CZ-NH1	7.91	124.25	120.30
1	B	138	LEU	CA-CB-CG	6.84	131.04	115.30
1	A	306	ILE	CB-CA-C	-6.66	98.29	111.60
1	A	93	ARG	NE-CZ-NH1	6.39	123.50	120.30
1	A	188	ARG	NE-CZ-NH1	-6.08	117.26	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	A	206	LEU	CA-CB-CG	5.68	128.38	115.30
1	B	188	ARG	NE-CZ-NH1	-5.58	117.51	120.30
1	A	284	LEU	CB-CG-CD1	5.45	120.27	111.00
1	A	129	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	103	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	273	LEU	CB-CG-CD1	-5.01	102.48	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	382	LEU	Peptide

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	2998	0	3010	32	1
1	B	2988	0	2988	34	1
2	A	15	0	7	2	0
2	B	15	0	7	0	0
3	A	8	0	14	0	0
4	A	139	0	0	4	0
4	B	137	0	0	7	0
All	All	6300	0	6026	65	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 5.

All (65) 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:247[A]:ARG:HH11	1:A:247[A]:ARG:CG	1.60	1.12
1:A:247[A]:ARG:NH1	1:A:247[A]:ARG:HG2	1.53	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101[B]:ARG:HH21	1:A:101[B]:ARG:HG3	1.17	1.03
1:A:247[A]:ARG:HH11	1:A:247[A]:ARG:HG2	0.78	0.95
1:A:101[B]:ARG:HH21	1:A:101[B]:ARG:CG	1.88	0.87
1:B:284:LEU:CD2	1:B:313:TYR:HB2	2.04	0.86
1:B:284:LEU:HD23	1:B:313:TYR:HB2	1.62	0.80
1:B:5:ASP:HB2	4:B:608:HOH:O	1.83	0.77
1:A:303:MET:HB2	1:A:306:ILE:HD11	1.68	0.76
1:B:302:GLY:HA2	4:B:607:HOH:O	1.86	0.75
1:B:35[A]:ARG:HD3	4:B:634:HOH:O	1.89	0.71
1:A:247[B]:ARG:NH1	4:A:550:HOH:O	2.26	0.69
1:A:101[B]:ARG:HG3	1:A:101[B]:ARG:NH2	1.99	0.68
1:B:103:ARG:HG3	4:B:555:HOH:O	1.93	0.68
1:B:286:GLN:HG2	1:B:315:LEU:HD22	1.78	0.65
1:A:139:SER:HA	1:A:169:TYR:OH	1.98	0.64
1:A:101[B]:ARG:CG	1:A:101[B]:ARG:NH2	2.55	0.62
1:A:254:ARG:HD2	4:A:546:HOH:O	1.99	0.61
1:B:201:PRO:HD2	1:B:205[A]:HIS:CG	2.38	0.59
1:A:155:ASN:ND2	2:A:401:PLP:H2A1	2.20	0.57
1:A:2:ALA:N	1:A:66:HIS:HE2	2.03	0.56
1:B:172:ARG:HD3	1:B:195:ILE:HD12	1.87	0.56
1:A:259:SER:HB2	1:A:296:ILE:HG21	1.90	0.53
1:A:155:ASN:HD22	2:A:401:PLP:H2A1	1.74	0.52
1:B:153:SER:O	1:B:154:SER:CB	2.58	0.52
1:B:259:SER:HB2	1:B:296:ILE:HG21	1.92	0.52
1:A:11:ASP:OD1	1:A:41[A]:ARG:NH2	2.43	0.52
1:B:78:VAL:HG13	1:B:82:MET:HE1	1.91	0.51
1:A:101[A]:ARG:HD2	1:A:113:TRP:CE2	2.44	0.51
1:A:255:GLY:HA2	1:A:280:ILE:HG23	1.93	0.51
1:A:297:ARG:HB2	4:A:559:HOH:O	2.10	0.50
1:A:11:ASP:OD2	1:A:41[A]:ARG:NH1	2.44	0.50
1:B:300:GLU:HB3	4:B:541:HOH:O	2.12	0.49
1:B:247[B]:ARG:NH1	4:B:630:HOH:O	2.42	0.48
1:A:11:ASP:CG	1:A:41[A]:ARG:HH22	2.16	0.48
1:B:139:SER:HA	1:B:169:TYR:OH	2.13	0.48
1:B:257:ALA:HA	1:B:283:VAL:O	2.14	0.48
1:A:101[B]:ARG:NH2	4:A:567:HOH:O	2.09	0.47
1:B:197:ASP:HA	1:B:198:PRO:HD3	1.79	0.47
1:A:284:LEU:HD22	1:A:313:TYR:HB2	1.98	0.46
1:B:2:ALA:N	1:B:66:HIS:HE2	2.13	0.46
1:A:85:PRO:HD2	1:A:89:ASP:OD2	2.16	0.45
1:B:5:ASP:CB	4:B:608:HOH:O	2.55	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247[A]:ARG:CG	1:A:247[A]:ARG:NH1	2.34	0.45
1:A:37:LEU:HD21	1:A:46:TYR:CD2	2.52	0.44
1:B:382:LEU:HD23	1:B:382:LEU:HA	1.75	0.44
1:B:81:GLU:OE1	1:B:81:GLU:N	2.38	0.44
1:A:245:GLN:HG2	1:B:81:GLU:O	2.18	0.44
1:A:101[A]:ARG:HD2	1:A:113:TRP:CZ2	2.53	0.44
1:A:369[A]:ASP:OD1	1:A:373:LYS:NZ	2.38	0.43
1:B:303:MET:O	1:B:307:ASN:ND2	2.46	0.43
1:B:352:LYS:HB3	1:B:358:LEU:HD11	2.00	0.43
1:B:90:PHE:CE2	1:B:121:PRO:HB3	2.54	0.43
1:B:85:PRO:HD2	1:B:89:ASP:OD1	2.19	0.43
1:B:67:THR:O	1:B:71:LEU:HG	2.19	0.42
1:A:41[A]:ARG:HD2	1:A:41[A]:ARG:HA	1.78	0.41
1:B:133:GLU:HB2	1:B:233:ALA:HB2	2.02	0.41
1:B:138:LEU:C	1:B:138:LEU:HD23	2.41	0.41
1:A:117:GLU:HB2	1:A:367:VAL:O	2.21	0.41
1:B:207:LEU:HD23	1:B:207:LEU:HA	1.94	0.41
1:A:11:ASP:OD1	1:A:41[B]:ARG:NH1	2.52	0.41
1:B:127:LYS:HE2	1:B:156:PHE:HB2	2.03	0.40
1:B:281[A]:ARG:N	1:B:281[A]:ARG:HD3	2.36	0.40
1:B:251:LEU:HD12	1:B:251:LEU:N	2.36	0.40
1:B:289[A]:GLN:HG2	1:B:298:ARG:CZ	2.52	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:101[A]:ARG:NH2	1:B:23:GLU:OE2[5_654]	2.19	0.01

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	388/389 (100%)	376 (97%)	11 (3%)	1 (0%)	37	31
1	B	388/389 (100%)	376 (97%)	10 (3%)	2 (0%)	25	17
All	All	776/778 (100%)	752 (97%)	21 (3%)	3 (0%)	30	23

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	SER
1	B	154	SER
1	B	223	ASN

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	317/313 (101%)	302 (95%)	15 (5%)	22	15
1	B	316/313 (101%)	299 (95%)	17 (5%)	18	11
All	All	633/626 (101%)	601 (95%)	32 (5%)	23	13

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	41[A]	ARG
1	A	41[B]	ARG
1	A	62	ASP
1	A	88	LEU
1	A	138	LEU
1	A	140	ARG
1	A	199	GLU
1	A	247[A]	ARG
1	A	247[B]	ARG
1	A	284	LEU
1	A	306	ILE
1	A	309	LEU

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Mol	Chain	Res	Type
1	A	312	SER
1	A	347	LYS
1	B	29	LEU
1	B	35[A]	ARG
1	B	35[B]	ARG
1	B	77	SER
1	B	78	VAL
1	B	88	LEU
1	B	138	LEU
1	B	143	GLU
1	B	154	SER
1	B	281[A]	ARG
1	B	281[B]	ARG
1	B	309	LEU
1	B	312	SER
1	B	314	THR
1	B	321	GLU
1	B	352	LYS
1	B	370	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

3 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
3	MPD	A	402	-	7,7,7	0.64	0	9,10,10	0.69	0
2	PLP	B	400	1	15,15,16	2.11	2 (13%)	21,22,23	3.19	9 (42%)
2	PLP	A	401	1	15,15,16	1.53	3 (20%)	21,22,23	3.29	11 (52%)

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	MPD	A	402	-	-	1/5/5/5	-
2	PLP	B	400	1	-	1/6/6/8	0/1/1/1
2	PLP	A	401	1	-	1/6/6/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	400	PLP	P-O3P	-6.69	1.29	1.54
2	B	400	PLP	C2A-C2	3.25	1.55	1.50
2	A	401	PLP	O4P-C5A	2.85	1.55	1.44
2	A	401	PLP	C2-N1	2.56	1.38	1.33
2	A	401	PLP	C3-C2	-2.51	1.38	1.41

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	PLP	O3P-P-O4P	8.27	128.24	106.67
2	A	401	PLP	O4P-P-O1P	-7.59	85.92	106.44
2	B	400	PLP	O4P-P-O1P	-5.18	92.43	106.44
2	A	401	PLP	O3P-P-O1P	-5.11	90.92	110.83
2	B	400	PLP	O2P-P-O4P	4.89	119.42	106.67
2	B	400	PLP	O2P-P-O1P	-4.85	91.94	110.83
2	B	400	PLP	O3P-P-O1P	-4.76	92.29	110.83
2	A	401	PLP	O3P-P-O2P	4.68	125.35	107.80
2	A	401	PLP	O2P-P-O4P	4.63	118.75	106.67
2	A	401	PLP	O4P-C5A-C5	4.53	117.85	109.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PLP	O2P-P-O1P	-4.51	93.27	110.83
2	A	401	PLP	C4A-C4-C5	-4.35	116.46	120.94
2	B	400	PLP	C4A-C4-C5	-3.98	116.84	120.94
2	A	401	PLP	O3P-P-O4P	3.54	115.90	106.67
2	B	400	PLP	O4P-C5A-C5	3.33	115.61	109.36
2	A	401	PLP	C6-C5-C4	-2.74	115.84	118.10
2	A	401	PLP	C3-C4-C5	2.65	121.77	118.59
2	B	400	PLP	O3-C3-C2	2.46	122.68	117.58
2	B	400	PLP	C3-C4-C5	2.44	121.52	118.59
2	A	401	PLP	C2A-C2-C3	-2.07	118.38	120.80

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	PLP	C5A-O4P-P-O3P
2	B	400	PLP	C5A-O4P-P-O3P
3	A	402	MPD	O2-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PLP	2	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, 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	382/389 (98%)	-0.53	2 (0%) 87 89	11, 29, 46, 59	8 (2%)
1	B	383/389 (98%)	-0.48	1 (0%) 90 92	12, 30, 47, 58	7 (1%)
All	All	765/778 (98%)	-0.51	3 (0%) 89 90	11, 29, 46, 59	15 (1%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	141[A]	ARG	3.1
1	A	247[A]	ARG	2.7
1	B	382	LEU	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 monosaccharides 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
3	MPD	A	402	8/8	0.96	0.06	17,27,32,34	0
2	PLP	B	400	15/16	0.98	0.05	15,21,27,27	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PLP	A	401	15/16	0.98	0.05	17,23,26,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.