



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2016 – 07:49 PM EDT

PDB ID : 3VSC
Title : Crystal Structure of the K127A Mutant of O-Phosphoserine Sulfhydrylase Complexed with External Schiff Base of Pyridoxal 5'-Phosphate with O-Phospho-L-Serine
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

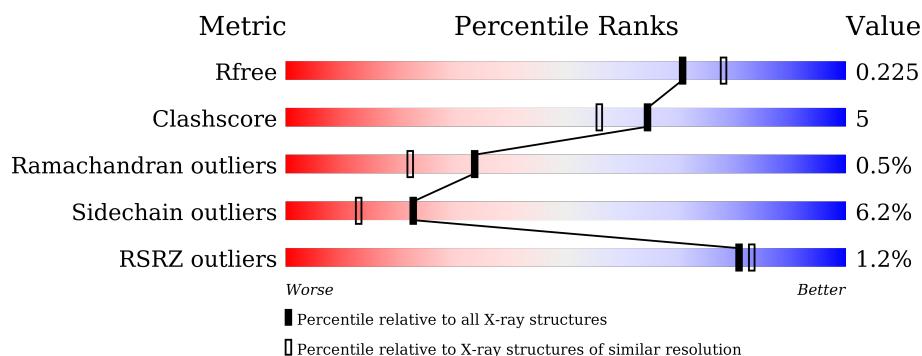
1 Overall quality at a glance

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}	91344	1799 (2.08-2.04)
Clashscore	102246	1910 (2.08-2.04)
Ramachandran outliers	100387	1893 (2.08-2.04)
Sidechain outliers	100360	1893 (2.08-2.04)
RSRZ outliers	91569	1802 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	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;"> % 82% 13% ... </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;"> % 80% 16% ... </div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6159 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.

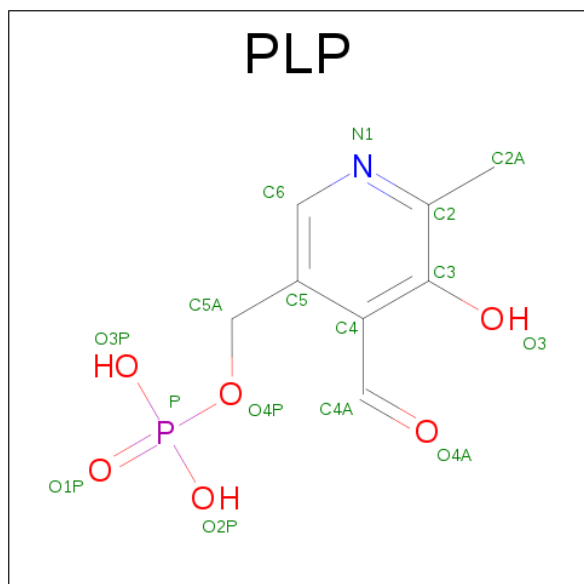
- Molecule 1 is a protein called Protein CysO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	4	0
			2949	1869	518	553	9			
1	B	383	Total	C	N	O	S	0	5	0
			2961	1875	521	556	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	ALA	LYS	ENGINEERED MUTATION	UNP Q9YBL2
B	127	ALA	LYS	ENGINEERED MUTATION	UNP Q9YBL2

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



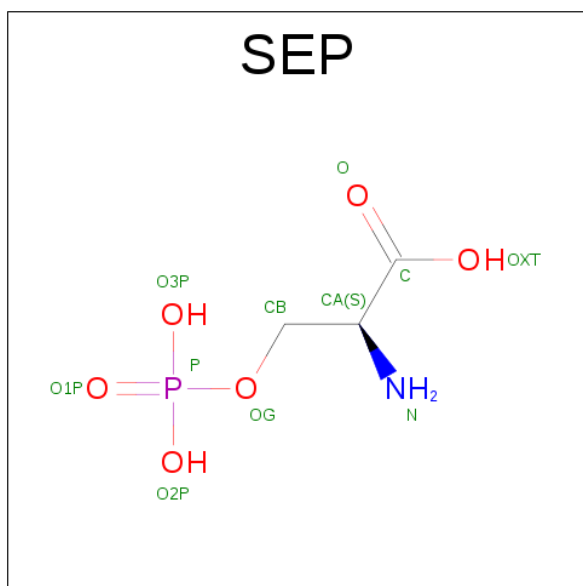
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

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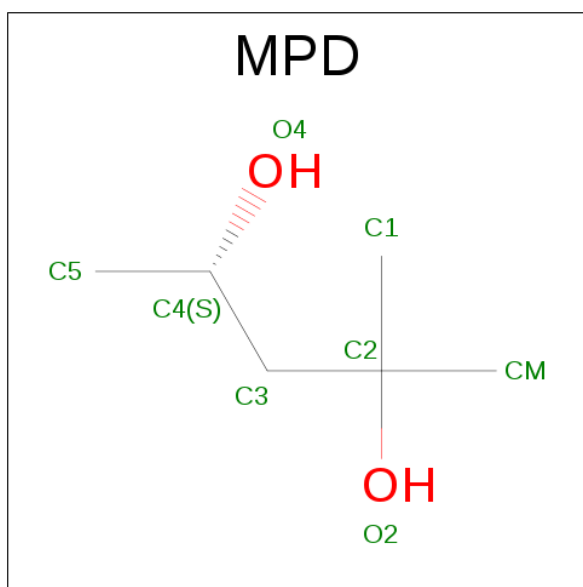
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is PHOSPHOSERINE (three-letter code: SEP) (formula: $C_3H_8NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			11	3	1	6	1		
3	B	1	Total	C	N	O	P	0	0
			11	3	1	6	1		

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



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

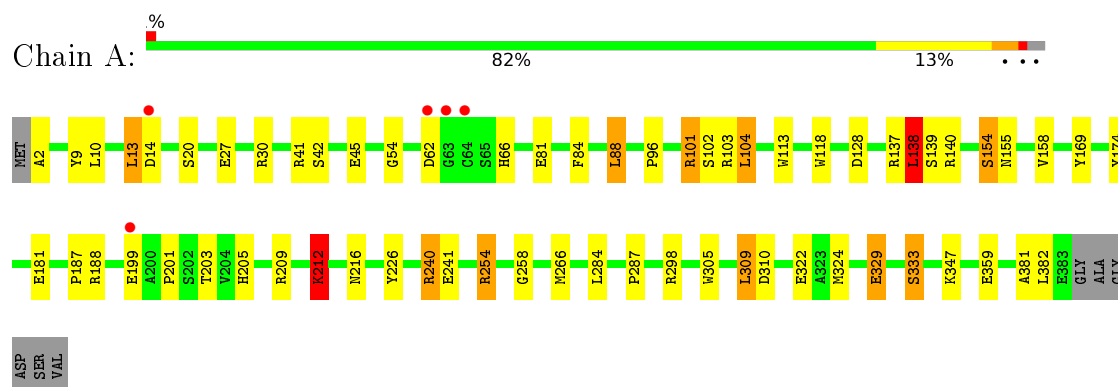
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	102	Total	O	0	0
			102	102		
5	B	87	Total	O	0	0
			87	87		

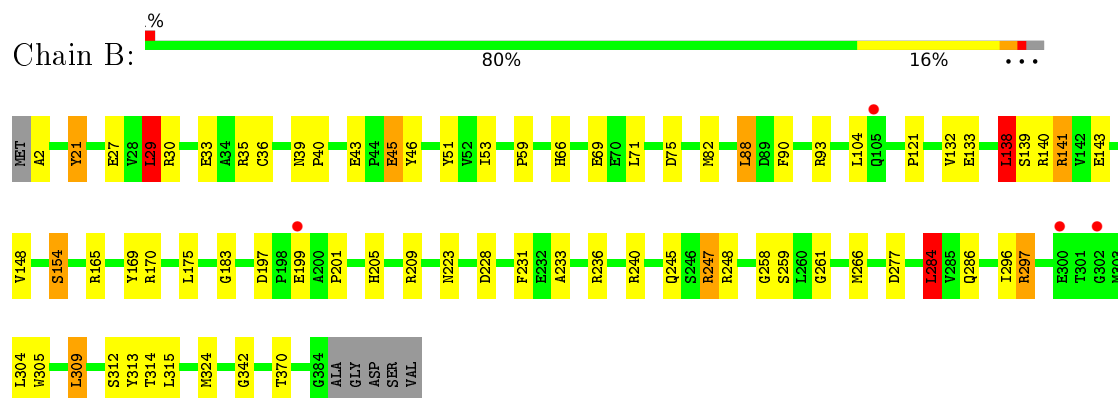
3 Residue-property plots [i](#)

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: 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	74.68 Å 74.68 Å 276.12 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.07 34.88 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.00-2.07) 99.3 (34.88-2.07)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.08 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.169 , 0.227 0.169 , 0.225	Depositor DCC
R_{free} test set	2439 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48293 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6159	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, SEP

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	1.24	9/3009 (0.3%)	1.08	12/4085 (0.3%)
1	B	1.26	8/3021 (0.3%)	1.13	21/4101 (0.5%)
All	All	1.25	17/6030 (0.3%)	1.10	33/8186 (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

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	181	GLU	CG-CD	6.32	1.61	1.51
1	B	148	VAL	CB-CG1	6.26	1.66	1.52
1	B	27	GLU	CG-CD	6.10	1.61	1.51
1	B	33	GLU	CG-CD	6.09	1.61	1.51
1	B	342	GLY	N-CA	6.08	1.55	1.46
1	A	118	TRP	CE3-CZ3	5.67	1.48	1.38
1	A	241	GLU	CB-CG	5.52	1.62	1.52
1	A	226	TYR	CG-CD2	5.46	1.46	1.39
1	A	181	GLU	CB-CG	5.45	1.62	1.52
1	A	329	GLU	CG-CD	5.44	1.60	1.51
1	B	51	TYR	CD1-CE1	5.38	1.47	1.39
1	B	21	TYR	CD1-CE1	5.23	1.47	1.39
1	B	132	VAL	CB-CG1	5.23	1.63	1.52
1	A	174	TYR	CD2-CE2	5.20	1.47	1.39
1	B	231	PHE	CE1-CZ	5.07	1.47	1.37
1	A	84	PHE	CE2-CZ	5.07	1.47	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	84	PHE	CD1-CE1	5.05	1.49	1.39

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ARG	NE-CZ-NH1	10.28	125.44	120.30
1	B	297	ARG	NE-CZ-NH1	10.16	125.38	120.30
1	B	93	ARG	NE-CZ-NH2	7.71	124.15	120.30
1	A	212	LYS	CD-CE-NZ	7.54	129.03	111.70
1	B	297	ARG	NE-CZ-NH2	-7.46	116.57	120.30
1	A	137	ARG	NE-CZ-NH1	-7.04	116.78	120.30
1	B	29	LEU	CB-CG-CD2	6.99	122.88	111.00
1	B	104	LEU	CB-CG-CD1	-6.76	99.51	111.00
1	A	103	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	B	88	LEU	CB-CG-CD1	6.58	122.18	111.00
1	B	170	ARG	NE-CZ-NH2	6.49	123.55	120.30
1	A	240	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	B	247	ARG	NE-CZ-NH2	6.40	123.50	120.30
1	A	88	LEU	CA-CB-CG	6.26	129.71	115.30
1	B	138	LEU	CA-CB-CG	6.08	129.29	115.30
1	B	228	ASP	CB-CG-OD1	5.98	123.69	118.30
1	A	41	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	128	ASP	CB-CG-OD2	-5.73	113.15	118.30
1	A	188	ARG	NE-CZ-NH1	-5.62	117.49	120.30
1	B	165	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	B	93	ARG	NE-CZ-NH1	-5.54	117.53	120.30
1	B	165	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	B	324	MET	CG-SD-CE	5.48	108.96	100.20
1	B	248	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	A	138	LEU	CA-CB-CG	5.28	127.45	115.30
1	B	277	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	B	141	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	B	284	LEU	CB-CG-CD1	5.16	119.76	111.00
1	A	254	ARG	NE-CZ-NH2	-5.15	117.73	120.30
1	A	254	ARG	NE-CZ-NH1	5.05	122.82	120.30
1	B	247	ARG	CG-CD-NE	5.05	122.40	111.80
1	B	277	ASP	CB-CG-OD1	5.04	122.83	118.30
1	B	30	ARG	NE-CZ-NH2	-5.00	117.80	120.30

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	2949	0	2948	34	0
1	B	2961	0	2956	29	0
2	A	15	0	7	1	0
2	B	15	0	7	0	0
3	A	11	0	3	0	0
3	B	11	0	3	2	0
4	A	8	0	14	0	0
5	A	102	0	0	2	0
5	B	87	0	0	3	0
All	All	6159	0	5938	61	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 5.

All (61) 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:212:LYS:HE3	1:A:216:ASN:ND2	1.76	1.01
1:A:254:ARG:HD2	5:A:572:HOH:O	1.69	0.92
1:B:286:GLN:HG2	1:B:315:LEU:HD22	1.68	0.75
1:B:35[B]:ARG:HD3	1:B:35[B]:ARG:O	1.87	0.74
1:A:138:LEU:HD23	1:A:138:LEU:C	2.07	0.74
1:A:212:LYS:HE3	1:A:216:ASN:HD22	1.50	0.74
1:A:139:SER:HA	1:A:169:TYR:OH	1.88	0.73
1:A:212:LYS:CE	1:A:216:ASN:HD22	2.02	0.72
1:A:212:LYS:CE	1:A:216:ASN:ND2	2.51	0.71
1:A:45:GLU:HG2	5:A:504:HOH:O	1.92	0.69
1:B:21:TYR:CB	1:B:45:GLU:HG3	2.29	0.63
1:B:21:TYR:HB3	1:B:45:GLU:HG3	1.83	0.61
1:B:261:GLY:CA	3:B:402:SEP:O3P	2.49	0.61
1:B:284:LEU:HD22	1:B:313:TYR:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:SER:OG	1:A:104:LEU:HB2	2.05	0.57
1:B:35[B]:ARG:HD2	1:B:46:TYR:CZ	2.40	0.57
1:A:212:LYS:NZ	1:A:216:ASN:HD22	2.03	0.56
1:A:27[A]:GLU:OE1	1:A:30:ARG:NH2	2.37	0.56
1:B:29:LEU:HD13	1:B:53:ILE:HD12	1.90	0.54
1:A:305:TRP:HB3	1:A:309:LEU:HD22	1.91	0.53
1:A:329:GLU:O	1:A:333:SER:HB3	2.09	0.53
1:B:197:ASP:OD1	1:B:199:GLU:HB2	2.10	0.52
1:A:201:PRO:HD2	1:A:205[B]:HIS:CG	2.45	0.51
1:A:212:LYS:HE3	1:A:216:ASN:HD21	1.70	0.51
1:B:261:GLY:HA2	3:B:402:SEP:O3P	2.11	0.49
1:B:258:GLY:HA3	1:B:266:MET:SD	2.52	0.49
1:B:201:PRO:HD2	1:B:205[B]:HIS:CG	2.47	0.48
1:B:138:LEU:HD23	1:B:138:LEU:C	2.34	0.48
1:A:254:ARG:NE	1:A:254:ARG:HA	2.29	0.47
1:A:101:ARG:HG2	1:A:113:TRP:CE2	2.49	0.47
1:B:35[B]:ARG:HD3	1:B:35[B]:ARG:C	2.34	0.47
1:A:10:LEU:O	1:A:13:LEU:HB2	2.15	0.46
1:A:2:ALA:N	1:A:66:HIS:HE2	2.13	0.46
1:B:154:SER:OG	1:B:183:GLY:HA3	2.16	0.46
1:A:54:GLY:HA3	1:A:240:ARG:HD3	1.97	0.45
1:B:139:SER:HA	1:B:169:TYR:OH	2.15	0.45
1:A:104:LEU:HA	1:A:104:LEU:HD22	1.78	0.45
1:A:101:ARG:HG2	1:A:113:TRP:CD2	2.52	0.45
1:A:322:GLU:HB3	1:A:347:LYS:HG2	1.99	0.44
1:A:324:MET:HB3	1:A:381:ALA:HB2	1.99	0.44
1:B:143:GLU:HG2	5:B:520:HOH:O	2.17	0.44
1:A:81:GLU:O	1:B:245:GLN:HG2	2.17	0.44
1:B:2:ALA:N	1:B:66:HIS:HE2	2.16	0.44
1:A:155:ASN:ND2	2:A:401:PLP:H2A1	2.34	0.43
1:B:305:TRP:HB3	1:B:309:LEU:HD22	2.00	0.43
1:B:36:CYS:HA	1:B:59:PRO:HB2	2.01	0.43
1:B:69:GLU:HB2	5:B:509:HOH:O	2.19	0.42
1:A:158:VAL:HG22	1:A:187:PRO:HD3	2.02	0.42
1:A:96:PRO:O	1:B:82:MET:HG3	2.20	0.42
1:A:212:LYS:HE2	1:A:212:LYS:C	2.41	0.41
1:B:45:GLU:HB2	5:B:504:HOH:O	2.20	0.41
1:A:254:ARG:HD3	1:A:359:GLU:OE1	2.19	0.41
1:A:9:TYR:CD1	1:A:9:TYR:N	2.89	0.41
1:A:138:LEU:CD2	1:A:138:LEU:C	2.85	0.41
1:A:287:PRO:O	1:A:298:ARG:NH1	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:GLY:HA3	1:A:266:MET:SD	2.62	0.40
1:B:304:LEU:HD23	1:B:305:TRP:CE2	2.56	0.40
1:B:39:ASN:HA	1:B:40:PRO:HD3	1.55	0.40
1:B:90:PHE:CE2	1:B:121:PRO:HB3	2.56	0.40
1:B:133:GLU:HB2	1:B:233:ALA:HB2	2.03	0.40
1:B:259:SER:HB2	1:B:296:ILE:HG21	2.04	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	384/389 (99%)	369 (96%)	12 (3%)	3 (1%)	24	12
1	B	386/389 (99%)	376 (97%)	8 (2%)	2 (0%)	34	22
All	All	770/778 (99%)	745 (97%)	20 (3%)	5 (1%)	34	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	SER
1	B	154	SER
1	B	223	ASN
1	A	14[A]	ASP
1	A	14[B]	ASP

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	312/312 (100%)	294 (94%)	18 (6%)	25	14
1	B	313/312 (100%)	293 (94%)	20 (6%)	22	12
All	All	625/624 (100%)	587 (94%)	38 (6%)	23	13

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	20	SER
1	A	42	SER
1	A	62	ASP
1	A	88	LEU
1	A	101	ARG
1	A	104	LEU
1	A	138	LEU
1	A	140	ARG
1	A	154	SER
1	A	199	GLU
1	A	203	THR
1	A	209	ARG
1	A	212	LYS
1	A	284	LEU
1	A	309	LEU
1	A	310	ASP
1	A	333	SER
1	B	29	LEU
1	B	43	GLU
1	B	45	GLU
1	B	71	LEU
1	B	75	ASP
1	B	88	LEU
1	B	138	LEU
1	B	140	ARG
1	B	141	ARG
1	B	175	LEU
1	B	209	ARG
1	B	236	ARG
1	B	240	ARG
1	B	247	ARG
1	B	284	LEU

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Mol	Chain	Res	Type
1	B	297	ARG
1	B	309	LEU
1	B	312	SER
1	B	314	THR
1	B	370	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	ASN

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](#)

5 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	3	15,15,16	1.43	2 (13%)	21,22,23	3.49	11 (52%)
3	SEP	A	402	2	7,10,10	2.10	3 (42%)	7,14,14	4.85	4 (57%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPD	A	403	-	6,7,7	0.70	0	6,10,10	0.68	0
2	PLP	B	401	3	15,15,16	2.23	5 (33%)	21,22,23	3.31	11 (52%)
3	SEP	B	402	2	7,10,10	2.18	2 (28%)	7,14,14	2.92	2 (28%)

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	3	-	0/6/6/8	0/1/1/1
3	SEP	A	402	2	-	0/6/10/10	0/0/0/0
4	MPD	A	403	-	-	0/5/5/5	0/0/0/0
2	PLP	B	401	3	-	0/6/6/8	0/1/1/1
3	SEP	B	402	2	-	0/6/10/10	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	PLP	P-O3P	-5.55	1.35	1.54
2	B	401	PLP	C3-C2	-3.00	1.38	1.40
3	A	402	SEP	P-O3P	2.04	1.61	1.54
2	B	401	PLP	O3-C3	2.32	1.42	1.37
2	A	401	PLP	P-O2P	2.50	1.63	1.54
2	A	401	PLP	O4P-C5A	2.61	1.54	1.45
3	A	402	SEP	P-O1P	2.97	1.60	1.50
2	B	401	PLP	O4P-C5A	3.12	1.56	1.45
2	B	401	PLP	C2A-C2	3.27	1.56	1.50
3	A	402	SEP	P-OG	3.31	1.69	1.59
3	B	402	SEP	P-OG	3.34	1.69	1.59
3	B	402	SEP	P-O1P	4.18	1.64	1.50

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PLP	O4P-P-O1P	-7.05	89.35	107.08
2	B	401	PLP	O2P-P-O1P	-6.69	88.79	110.63
2	A	401	PLP	O2P-P-O1P	-6.15	90.54	110.63
2	A	401	PLP	O3P-P-O1P	-5.95	91.22	110.63
2	B	401	PLP	P-O4P-C5A	-5.77	107.42	120.79
2	B	401	PLP	O4P-P-O1P	-5.14	94.17	107.08
2	A	401	PLP	P-O4P-C5A	-4.78	109.72	120.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PLP	C4A-C4-C5	-4.29	116.48	120.90
2	B	401	PLP	C4A-C4-C5	-4.18	116.60	120.90
3	A	402	SEP	OG-P-O1P	-3.38	98.59	107.08
2	B	401	PLP	O3P-P-O1P	-2.82	101.43	110.63
2	B	401	PLP	C5-C6-N1	-2.67	119.19	123.86
2	A	401	PLP	C4-C3-C2	-2.34	116.18	120.12
2	A	401	PLP	C2A-C2-C3	-2.06	118.81	120.90
2	B	401	PLP	O2P-P-O4P	2.08	112.78	106.72
2	B	401	PLP	C6-N1-C2	2.29	123.85	119.26
2	A	401	PLP	O3-C3-C4	2.31	124.40	118.11
3	A	402	SEP	O2P-P-OG	2.56	114.20	106.72
2	A	401	PLP	C5A-C5-C4	2.68	125.25	121.63
2	B	401	PLP	O3P-P-O2P	2.94	118.23	107.44
3	B	402	SEP	O2P-P-OG	2.95	115.35	106.72
3	A	402	SEP	O3P-P-OG	3.35	116.50	106.72
2	B	401	PLP	C3-C4-C5	3.78	123.12	118.68
2	A	401	PLP	O3P-P-O4P	5.33	122.28	106.72
2	A	401	PLP	O3P-P-O2P	5.36	127.11	107.44
3	B	402	SEP	OG-CB-CA	6.60	113.07	108.16
2	B	401	PLP	O3P-P-O4P	6.81	126.60	106.72
3	A	402	SEP	OG-CB-CA	11.55	116.75	108.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PLP	1	0
3	B	402	SEP	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%)	-0.40	5 (1%) 79 81	22, 33, 51, 64	0
1	B	383/389 (98%)	-0.35	4 (1%) 84 86	23, 33, 51, 64	0
All	All	765/778 (98%)	-0.37	9 (1%) 81 83	22, 33, 51, 64	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	199	GLU	3.7
1	B	300	GLU	3.1
1	A	199	GLU	2.8
1	A	62	ASP	2.5
1	A	64	CYS	2.5
1	A	14[A]	ASP	2.4
1	B	105	GLN	2.4
1	A	63	GLY	2.4
1	B	302	GLY	2.3

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, 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	LLDF	B-factors(Å ²)	Q<0.9
3	SEP	A	402	11/11	0.94	0.11	0.94	32,35,54,57	0
4	MPD	A	403	8/8	0.96	0.12	0.70	32,38,40,42	0
2	PLP	A	401	15/16	0.97	0.12	0.60	23,29,32,33	0
2	PLP	B	401	15/16	0.99	0.13	0.28	22,25,29,30	0
3	SEP	B	402	11/11	0.93	0.11	0.27	30,32,55,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.