Full wwPDB X-ray Structure Validation Report (i)

Feb 2, 2016 – 12:01 AM GMT

PDB ID	:	8AAT
Title	:	X-RAY STRUCTURE REFINEMENT AND COMPARISON OF THREE FORMS OF MITOCHONDRIAL ASPARTATE AMINOTRANSFERASE
Authors	:	Mcphalen, C.A.; Vincent, M.G.; Jansonius, J.N.
Deposited on	:	1991-12-02
Resolution	:	2.30 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 (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

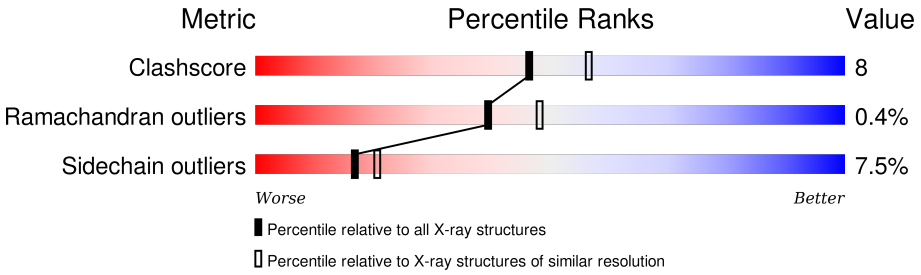
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)

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 <=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	401	80% 18%
1	B	401	79% 18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6952 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.

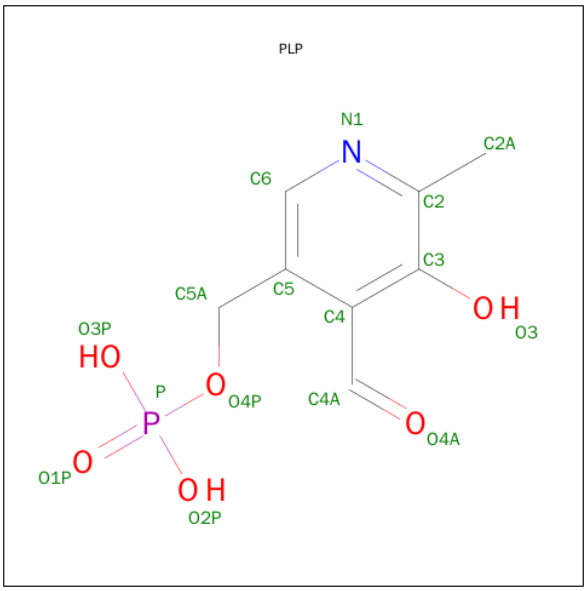
- Molecule 1 is a protein called ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	0	0
			3161	2004	558	581	18			
1	B	401	Total	C	N	O	S	0	0	0
			3161	2004	558	581	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	PRO	SER	CONFLICT	UNP P00508
B	47	PRO	SER	CONFLICT	UNP P00508

- 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		

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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 water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	297	Total	O	0	0
			297	297		
3	B	303	Total	O	0	0
			303	303		

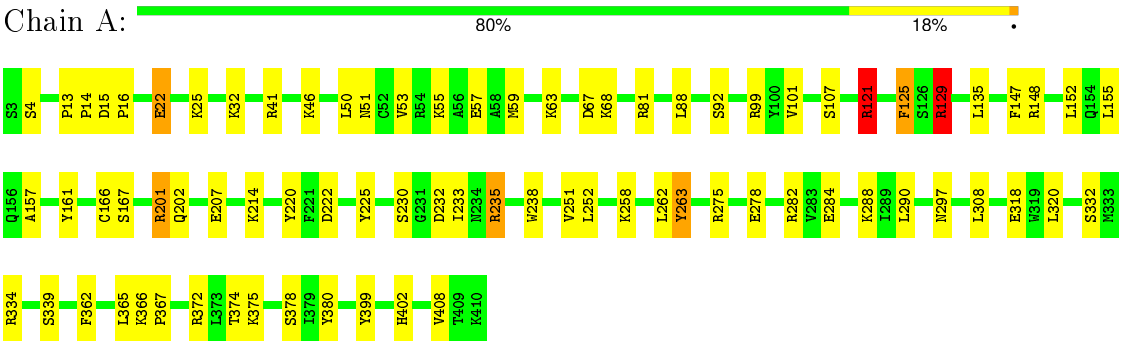
3 Residue-property plots (i)

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.

Note EDS was not executed.

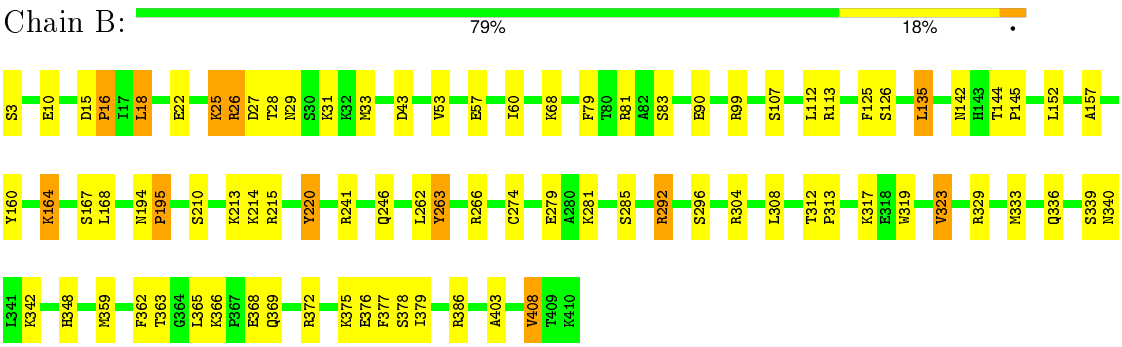
- Molecule 1: ASPARTATE AMINOTRANSFERASE

Chain A:



- Molecule 1: ASPARTATE AMINOTRANSFERASE

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	55.53Å 58.68Å 75.77Å 85.10° 109.20° 115.90°	Depositor
Resolution (Å)	10.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.130 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6952	wwPDB-VP
Average B, all atoms (Å^2)	20.0	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: 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	0.65	0/3231	1.24	13/4360 (0.3%)
1	B	0.65	0/3231	1.22	10/4360 (0.2%)
All	All	0.65	0/6462	1.23	23/8720 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	282	ARG	NE-CZ-NH1	10.53	125.56	120.30
1	A	121	ARG	NE-CZ-NH1	10.13	125.37	120.30
1	B	81	ARG	NE-CZ-NH1	9.01	124.80	120.30
1	B	215	ARG	NE-CZ-NH1	-8.39	116.11	120.30
1	A	121	ARG	NE-CZ-NH2	-8.36	116.12	120.30
1	B	81	ARG	CD-NE-CZ	8.30	135.22	123.60
1	A	129	ARG	NE-CZ-NH1	-8.11	116.24	120.30
1	A	334	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	A	41	ARG	NE-CZ-NH2	-7.65	116.47	120.30
1	B	99	ARG	NE-CZ-NH1	7.36	123.98	120.30
1	B	113	ARG	NE-CZ-NH2	-7.36	116.62	120.30
1	B	215	ARG	CD-NE-CZ	-6.49	114.52	123.60
1	B	215	ARG	NE-CZ-NH2	6.13	123.36	120.30
1	A	99	ARG	CD-NE-CZ	5.78	131.69	123.60
1	B	241	ARG	NE-CZ-NH1	-5.77	117.41	120.30
1	B	126	SER	N-CA-CB	-5.51	102.24	110.50
1	B	266	ARG	NE-CZ-NH2	-5.37	117.61	120.30
1	A	252	LEU	N-CA-CB	-5.34	99.71	110.40
1	A	297	ASN	CB-CA-C	5.27	120.93	110.40
1	A	41	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	99	ARG	NE-CZ-NH1	5.19	122.89	120.30
1	A	53	VAL	CB-CA-C	5.12	121.13	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	251	VAL	CB-CA-C	5.06	121.01	111.40

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	3161	0	3152	51	0
1	B	3161	0	3152	45	0
2	A	15	0	6	2	0
2	B	15	0	6	1	0
3	A	297	0	0	8	0
3	B	303	0	0	13	0
All	All	6952	0	6316	96	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 8.

All (96) 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:201:ARG:HH11	1:A:201:ARG:HG3	1.01	1.14
1:A:129:ARG:HA	1:A:129:ARG:HE	1.16	1.05
1:B:26:ARG:HH11	1:B:26:ARG:HG3	1.20	1.03
1:A:22:GLU:HG3	3:A:690:HOH:O	1.67	0.94
1:A:4:SER:HA	3:B:639:HOH:O	1.66	0.93
1:A:201:ARG:HH11	1:A:201:ARG:CG	1.82	0.92
1:A:201:ARG:HG3	1:A:201:ARG:NH1	1.81	0.87
1:B:33:MET:HG2	1:B:379:ILE:HG12	1.55	0.86
1:A:167:SER:HB2	3:A:491:HOH:O	1.77	0.83
1:B:26:ARG:HH11	1:B:26:ARG:CG	1.91	0.83
1:A:129:ARG:CA	1:A:129:ARG:HE	1.90	0.80
1:A:22:GLU:OE1	1:A:22:GLU:HA	1.84	0.78
1:A:81:ARG:HD3	3:A:614:HOH:O	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ARG:HA	1:A:129:ARG:NE	1.99	0.76
1:B:15:ASP:OD2	1:B:18:LEU:HB2	1.85	0.76
1:B:376:GLU:OE1	3:B:687:HOH:O	2.06	0.73
1:B:292:ARG:NH1	3:B:679:HOH:O	2.23	0.72
1:A:125:PHE:CE1	3:B:634:HOH:O	2.44	0.70
1:B:3:SER:O	3:B:691:HOH:O	2.09	0.70
1:B:363:THR:OG1	1:B:365:LEU:HD12	1.93	0.68
1:B:33:MET:CG	1:B:379:ILE:HG12	2.24	0.67
1:A:57:GLU:HG2	3:A:508:HOH:O	1.93	0.67
1:B:26:ARG:HG3	1:B:26:ARG:NH1	1.99	0.67
1:A:232:ASP:HB3	1:A:235:ARG:HH11	1.60	0.67
1:A:202:GLN:HG3	1:A:238:TRP:CH2	2.30	0.66
1:B:29:ASN:OD1	1:B:31:LYS:HB2	1.95	0.66
1:A:232:ASP:HB3	1:A:235:ARG:NH1	2.11	0.66
1:B:213:LYS:HE3	1:B:246:GLN:O	1.97	0.65
1:B:368:GLU:OE2	3:B:705:HOH:O	2.15	0.64
1:B:377:PHE:O	1:B:378:SER:HB2	1.95	0.64
1:A:125:PHE:HE1	3:B:634:HOH:O	1.83	0.61
1:B:22:GLU:HG3	3:B:441:HOH:O	2.00	0.61
1:A:101:VAL:HG21	1:A:284:GLU:HB2	1.81	0.61
1:A:22:GLU:CG	3:A:690:HOH:O	2.37	0.60
1:A:225:TYR:CE1	1:A:258:LYS:HD3	2.38	0.58
1:B:27:ASP:OD1	1:B:28:THR:N	2.37	0.57
1:B:79:PHE:O	1:B:83:SER:HB2	2.05	0.57
1:B:386:ARG:NH1	3:B:465:HOH:O	2.36	0.57
1:A:201:ARG:CG	1:A:201:ARG:NH1	2.48	0.56
1:B:329:ARG:O	1:B:333:MET:HG2	2.05	0.56
1:A:68:LYS:O	1:B:263:TYR:HB2	2.04	0.56
1:A:129:ARG:CA	1:A:129:ARG:NE	2.64	0.55
1:B:160:TYR:O	1:B:168:LEU:HD12	2.07	0.54
1:B:142:ASN:O	1:B:145:PRO:HD2	2.08	0.53
1:B:26:ARG:NH1	1:B:26:ARG:CG	2.58	0.51
1:A:88:LEU:O	1:A:92:SER:HB2	2.12	0.50
1:B:167:SER:HB2	3:B:500:HOH:O	2.12	0.50
1:A:372:ARG:HG2	1:A:408:VAL:HG12	1.93	0.50
2:B:411:PLP:O4P	2:B:411:PLP:C4A	2.59	0.50
1:B:10:GLU:OE1	3:B:414:HOH:O	2.19	0.49
1:B:53:VAL:O	1:B:57:GLU:HG3	2.13	0.48
1:B:274:CYS:HB3	1:B:279:GLU:HG2	1.95	0.48
1:A:135:LEU:O	1:A:157:ALA:HA	2.15	0.47
1:A:13:PRO:HA	1:A:14:PRO:HD3	1.76	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:SER:OG	1:A:235:ARG:HD2	2.14	0.47
1:A:263:TYR:HB2	1:B:68:LYS:O	2.15	0.47
1:A:202:GLN:HG3	1:A:238:TRP:CZ3	2.50	0.47
1:B:308:LEU:HD23	1:B:308:LEU:HA	1.73	0.46
1:A:233:ILE:HD13	1:A:320:LEU:HD21	1.98	0.46
1:A:262:LEU:O	1:A:263:TYR:C	2.54	0.46
1:B:144:THR:HB	1:B:145:PRO:HD3	1.98	0.46
1:A:32:LYS:HA	1:A:378:SER:O	2.16	0.46
1:A:59:MET:O	1:A:63:LYS:HG3	2.15	0.46
1:B:112:LEU:HA	1:B:220:TYR:OH	2.16	0.45
1:B:135:LEU:O	1:B:157:ALA:HA	2.16	0.45
1:A:50:LEU:N	1:A:50:LEU:HD12	2.32	0.45
1:A:15:ASP:HA	1:A:16:PRO:HD3	1.77	0.45
1:B:164:LYS:HB2	1:B:164:LYS:HE2	1.60	0.45
1:B:319:TRP:O	1:B:323:VAL:HG22	2.16	0.45
1:A:101:VAL:CG2	1:A:284:GLU:HB2	2.47	0.45
1:A:284:GLU:O	1:A:288:LYS:HG3	2.16	0.45
1:B:262:LEU:O	1:B:263:TYR:C	2.55	0.45
1:A:148:ARG:HD2	3:A:486:HOH:O	2.15	0.45
1:A:399:TYR:O	1:A:402:HIS:HB3	2.17	0.44
1:B:25:LYS:C	1:B:27:ASP:H	2.19	0.43
1:B:336:GLN:O	1:B:340:ASN:ND2	2.51	0.43
1:A:374:THR:HG23	1:A:380:TYR:CE1	2.54	0.43
1:B:366:LYS:HE2	3:B:705:HOH:O	2.18	0.43
1:B:348:HIS:CG	3:B:636:HOH:O	2.71	0.43
1:B:15:ASP:HA	1:B:16:PRO:HD3	1.73	0.43
1:B:60:ILE:HD13	1:B:304:ARG:HB3	2.01	0.43
1:B:363:THR:CB	1:B:365:LEU:HD12	2.49	0.42
1:A:67:ASP:OD1	1:A:67:ASP:C	2.58	0.42
1:A:366:LYS:HB3	1:A:367:PRO:HD2	2.02	0.42
1:A:46:LYS:NZ	3:A:538:HOH:O	2.39	0.42
1:A:372:ARG:HG2	1:A:408:VAL:CG1	2.50	0.41
2:A:411:PLP:O4P	2:A:411:PLP:C4A	2.68	0.41
1:A:147:PHE:HD1	1:A:152:LEU:HD12	1.84	0.41
1:B:403:ALA:O	1:B:408:VAL:HG13	2.21	0.41
1:A:55:LYS:NZ	1:A:318:GLU:OE1	2.44	0.41
1:A:121:ARG:HD3	3:A:469:HOH:O	2.21	0.41
1:A:161:TYR:OH	1:A:166:CYS:HA	2.22	0.40
1:A:101:VAL:HG21	1:A:284:GLU:CB	2.51	0.40
1:A:222:ASP:OD2	2:A:411:PLP:N1	2.54	0.40
1:B:312:THR:HA	1:B:313:PRO:HD3	1.94	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:194:ASN:HA	1:B:195:PRO:HA	1.78	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	399/401 (100%)	391 (98%)	7 (2%)	1 (0%)	46 57
1	B	399/401 (100%)	384 (96%)	13 (3%)	2 (0%)	34 41
All	All	798/802 (100%)	775 (97%)	20 (2%)	3 (0%)	39 48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	TYR
1	B	263	TYR
1	B	16	PRO

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	335/335 (100%)	313 (93%)	22 (7%)	21 27
1	B	335/335 (100%)	307 (92%)	28 (8%)	14 16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	670/670 (100%)	620 (92%)	50 (8%)	17 21

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	25	LYS
1	A	51	ASN
1	A	107	SER
1	A	121	ARG
1	A	125	PHE
1	A	129	ARG
1	A	155	LEU
1	A	201	ARG
1	A	207	GLU
1	A	214	LYS
1	A	220	TYR
1	A	235	ARG
1	A	275	ARG
1	A	278	GLU
1	A	290	LEU
1	A	308	LEU
1	A	332	SER
1	A	339	SER
1	A	362	PHE
1	A	365	LEU
1	A	375	LYS
1	B	18	LEU
1	B	25	LYS
1	B	26	ARG
1	B	43	ASP
1	B	90	GLU
1	B	107	SER
1	B	125	PHE
1	B	135	LEU
1	B	152	LEU
1	B	164	LYS
1	B	195	PRO
1	B	210	SER
1	B	214	LYS
1	B	220	TYR
1	B	281	LYS

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Mol	Chain	Res	Type
1	B	285	SER
1	B	292	ARG
1	B	296	SER
1	B	317	LYS
1	B	323	VAL
1	B	339	SER
1	B	342	LYS
1	B	359	MET
1	B	362	PHE
1	B	369	GLN
1	B	372	ARG
1	B	375	LYS
1	B	408	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	154	GLN
1	A	216	ASN
1	A	226	GLN
1	A	336	GLN
1	A	351	GLN
1	B	216	ASN
1	B	226	GLN
1	B	336	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)

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	411	1	15,15,16	1.17	1 (6%)	21,22,23	2.20	8 (38%)
2	PLP	B	411	1	15,15,16	1.32	1 (6%)	21,22,23	2.13	8 (38%)

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	411	1	-	0/6/6/8	0/1/1/1
2	PLP	B	411	1	-	0/6/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	411	PLP	C4A-C4	-3.48	1.44	1.51
2	A	411	PLP	C4A-C4	-3.28	1.45	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	411	PLP	C5-C6-N1	-2.93	118.78	123.86
2	B	411	PLP	C5-C6-N1	-2.88	118.86	123.86
2	A	411	PLP	C3-C2-N1	-2.80	116.75	120.61
2	B	411	PLP	C3-C2-N1	-2.39	117.31	120.61
2	A	411	PLP	C3-C4-C5	-2.37	116.19	118.78
2	B	411	PLP	C3-C4-C5	-2.03	116.56	118.78
2	A	411	PLP	O2P-P-O1P	2.01	117.06	110.58
2	B	411	PLP	C6-N1-C2	2.29	123.96	119.28
2	B	411	PLP	O2P-P-O1P	2.31	118.02	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	411	PLP	C6-N1-C2	2.52	124.42	119.28
2	B	411	PLP	C2A-C2-C3	3.71	125.51	121.04
2	A	411	PLP	C2A-C2-C3	4.01	125.88	121.04
2	B	411	PLP	C6-C5-C4	4.25	121.75	118.15
2	A	411	PLP	C6-C5-C4	4.33	121.82	118.15
2	A	411	PLP	O4P-C5A-C5	4.43	116.32	108.99
2	B	411	PLP	O4P-C5A-C5	4.50	116.43	108.99

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	411	PLP	2	0
2	B	411	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

6.1 Protein, DNA and RNA chains

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

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

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

6.3 Carbohydrates

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

6.4 Ligands

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

6.5 Other polymers

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