Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:04 PM GMT

PDB ID	:	1S07
Title	:	Crystal Structure of the R253A Mutant of 7,8-Diaminopelargonic Acid Synthase
Authors	:	Sandmark, J.; Eliot, A.C.; Famm, K.; Schneider, G.; Kirsch, J.F.
Deposited on	:	2003-12-30
Resolution	:	2.42 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 (1)) were used in the production of this report:

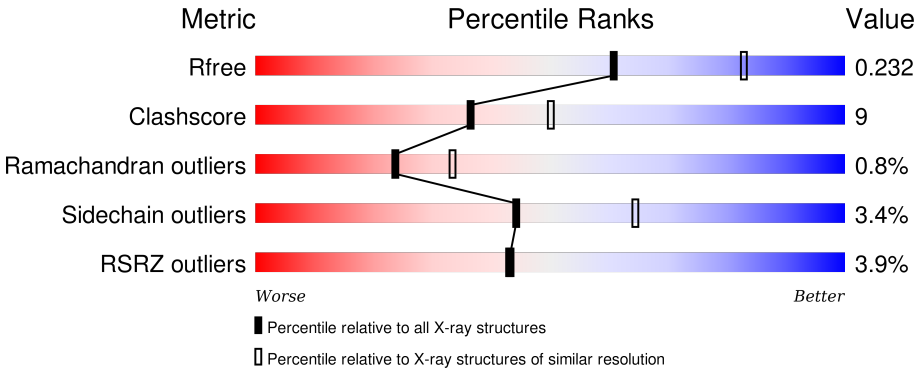
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.42 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	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

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.

Mol	Chain	Length	Quality of chain
1	A	429	4% 80% 19%
1	B	429	4% 79% 19%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NA	A	501	-	-	-	X
2	NA	B	502	-	-	-	X
3	PLP	A	430	-	-	X	-
4	IPA	B	600	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6898 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 Adenosylmethionine-8-amino-7-oxononanoate aminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	30	2	0
			3314	2106	573	603	32			
1	B	428	Total	C	N	O	S	19	2	0
			3303	2098	574	599	32			

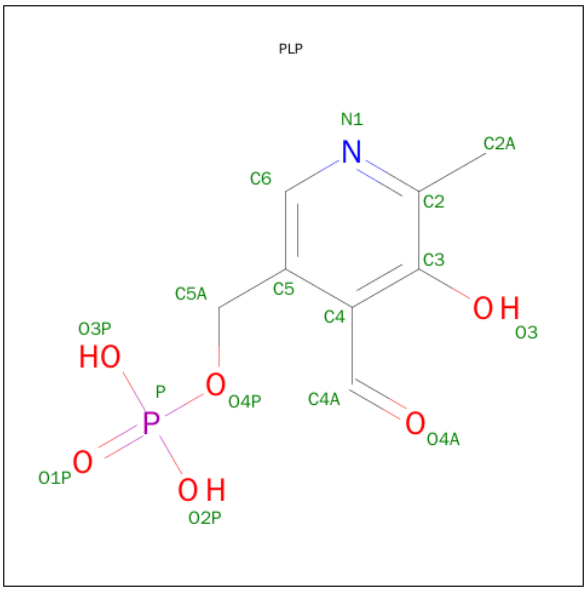
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	LEU	TRP	SEE REMARK 999	UNP P12995
A	253	ALA	ARG	ENGINEERED	UNP P12995
B	14	LEU	TRP	SEE REMARK 999	UNP P12995
B	253	ALA	ARG	ENGINEERED	UNP P12995

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

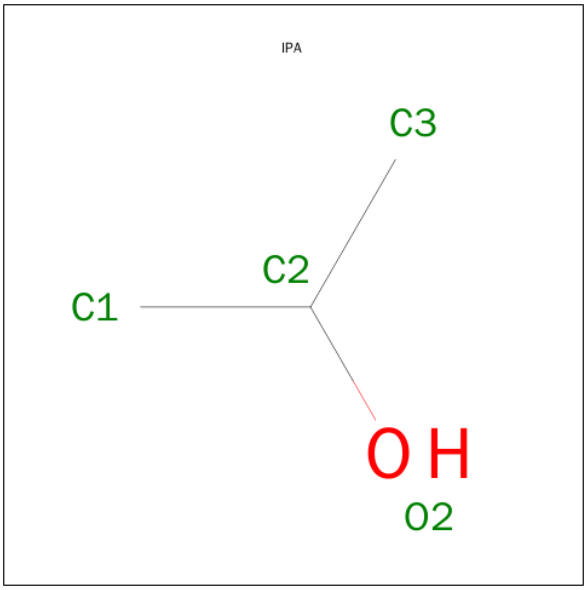
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	B	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

- Molecule 4 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C_3H_8O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	O		0	0
			4	3	1			

- Molecule 5 is water.

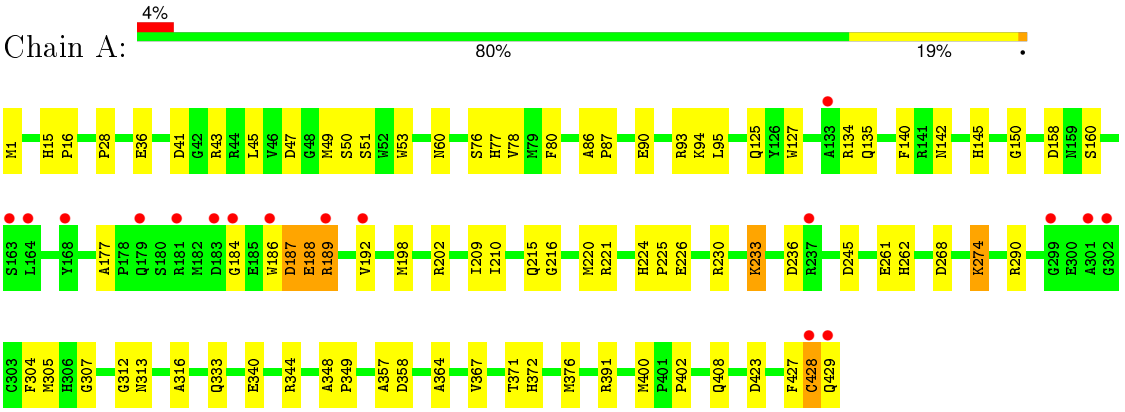
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	135	Total O 135 135	0	0
5	B	108	Total O 108 108	0	0

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.

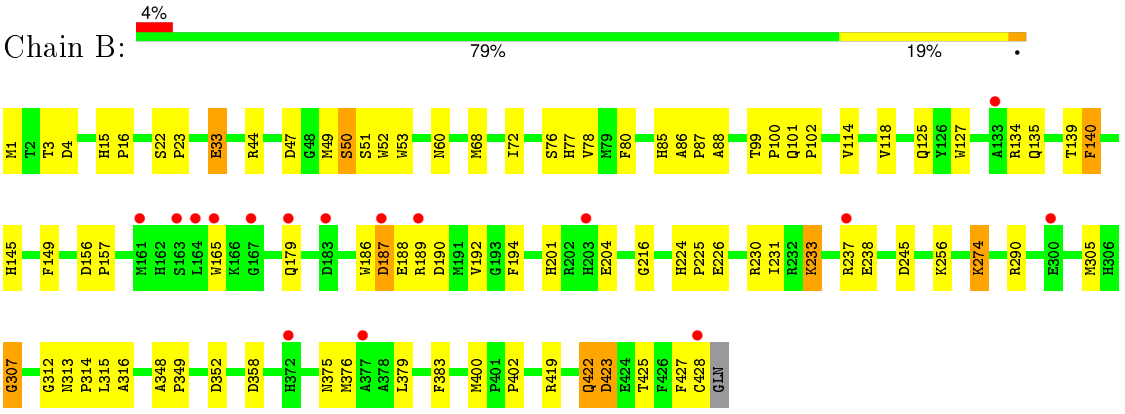
- Molecule 1: Adenosylmethionine-8-amino-7-oxononanoate aminotransferase

Chain A:



- Molecule 1: Adenosylmethionine-8-amino-7-oxononanoate aminotransferase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	58.08Å 56.53Å 120.99Å 90.00° 96.32° 90.00°	Depositor
Resolution (Å)	20.00 – 2.42 20.00 – 2.42	Depositor EDS
% Data completeness (in resolution range)	89.3 (20.00-2.42) 89.3 (20.00-2.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	1.81 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.188 , 0.224 0.195 , 0.232	Depositor DCC
R_{free} test set	1413 reflections (5.56%)	DCC
Wilson B-factor (Å^2)	30.5	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 36.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 26820 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6898	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 12.48% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NA, IPA, 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	0.72	0/3403	0.94	13/4620 (0.3%)
1	B	0.72	0/3391	0.88	7/4604 (0.2%)
All	All	0.72	0/6794	0.91	20/9224 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	140[A]	PHE	CB-CG-CD1	-10.19	113.67	120.80
1	A	140[B]	PHE	CB-CG-CD1	-10.19	113.67	120.80
1	A	140[A]	PHE	CB-CG-CD2	9.08	127.16	120.80
1	A	140[B]	PHE	CB-CG-CD2	9.08	127.16	120.80
1	B	140[A]	PHE	CB-CA-C	-7.04	96.32	110.40
1	B	140[B]	PHE	CB-CA-C	-7.04	96.32	110.40
1	A	423	ASP	CB-CG-OD2	6.41	124.07	118.30
1	A	333	GLN	CB-CG-CD	6.28	127.92	111.60
1	A	236	ASP	CB-CG-OD2	6.28	123.95	118.30
1	A	47	ASP	CB-CG-OD2	5.68	123.41	118.30
1	B	4	ASP	CB-CG-OD2	5.64	123.38	118.30
1	B	423	ASP	CB-CG-OD2	5.54	123.29	118.30
1	A	158	ASP	CB-CG-OD2	5.43	123.19	118.30
1	B	187	ASP	CB-CG-OD2	5.37	123.13	118.30
1	A	408	GLN	CA-CB-CG	5.35	125.17	113.40
1	A	187	ASP	CB-CG-OD2	5.33	123.09	118.30
1	A	268	ASP	CB-CG-OD2	5.29	123.06	118.30
1	B	358	ASP	CB-CG-OD2	5.26	123.04	118.30
1	B	352	ASP	CB-CG-OD2	5.14	122.92	118.30
1	A	184	GLY	N-CA-C	5.08	125.79	113.10

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	3314	0	3265	60	0
1	B	3303	0	3260	64	2
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	16	0	7	9	0
3	B	16	0	7	2	0
4	B	4	0	8	3	0
5	A	135	0	0	2	0
5	B	108	0	0	5	0
All	All	6898	0	6547	124	2

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 9.

All (124) 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:274:LYS:HZ1	3:A:430:PLP:C4A	1.60	1.15
1:A:274:LYS:NZ	3:A:430:PLP:C4A	2.29	0.94
3:A:430:PLP:O4A	4:B:600:IPA:H31	1.71	0.90
1:A:274:LYS:HZ1	3:A:430:PLP:C4	1.88	0.85
1:A:186:TRP:HE1	1:A:188:GLU:HG2	1.40	0.85
1:B:51:SER:HB2	1:B:400:MET:HG2	1.60	0.83
1:A:1:MET:HE3	1:A:28:PRO:HB2	1.60	0.83
1:A:77:HIS:HD2	5:B:601:HOH:O	1.61	0.83
1:B:33:GLU:OE2	1:B:44:ARG:NH2	2.12	0.81
1:B:49:MET:HB3	1:B:53:TRP:HZ3	1.45	0.81
1:B:139:THR:C	1:B:140[B]:PHE:HD2	1.84	0.80
1:A:51:SER:HB2	1:A:400:MET:HG2	1.66	0.77
1:B:226:GLU:OE1	1:B:230:ARG:NH1	2.18	0.77
1:A:226:GLU:OE1	1:A:230:ARG:NH1	2.20	0.74
1:A:93:ARG:HG2	1:A:93:ARG:HH11	1.51	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:LYS:NZ	3:A:430:PLP:C4	2.50	0.74
1:A:49:MET:HB3	1:A:53:TRP:HZ3	1.52	0.74
1:A:49:MET:HB3	1:A:53:TRP:CZ3	2.24	0.72
1:A:313:ASN:ND2	1:A:316:ALA:H	1.89	0.70
1:A:274:LYS:NZ	3:A:430:PLP:O4A	2.24	0.70
1:A:186:TRP:NE1	1:A:188:GLU:HG2	2.07	0.70
1:B:139:THR:C	1:B:140[B]:PHE:CD2	2.66	0.68
1:A:230:ARG:HD2	5:A:542:HOH:O	1.92	0.68
1:A:340:GLU:CD	1:A:344:ARG:HH12	1.98	0.66
1:A:340:GLU:CD	1:A:344:ARG:NH1	2.50	0.65
1:B:140[B]:PHE:HE2	1:B:194:PHE:HE1	1.46	0.64
1:B:49:MET:HB3	1:B:53:TRP:CZ3	2.29	0.63
1:B:135:GLN:CB	5:B:614:HOH:O	2.45	0.63
1:B:201:HIS:HD2	1:B:204:GLU:OE2	1.81	0.63
1:B:140[B]:PHE:CZ	1:B:231:ILE:HD11	2.34	0.62
1:B:140[B]:PHE:HE2	1:B:194:PHE:CE1	2.16	0.62
1:B:135:GLN:HB3	5:B:614:HOH:O	2.02	0.60
1:B:237[A]:ARG:NH1	1:B:238:GLU:OE1	2.35	0.59
1:A:150:GLY:HA3	1:B:149:PHE:CD1	2.36	0.59
1:B:313:ASN:HD22	1:B:316:ALA:H	1.51	0.59
1:A:50:SER:OG	1:A:402:PRO:HG3	2.03	0.58
1:A:125:GLN:HE22	1:A:305:MET:H	1.51	0.58
1:B:140[B]:PHE:CE2	1:B:194:PHE:CE1	2.91	0.58
1:B:348:ALA:N	1:B:349:PRO:CD	2.68	0.57
1:B:427:PHE:O	1:B:428:CYS:HB2	2.05	0.56
3:A:430:PLP:C4A	4:B:600:IPA:H31	2.35	0.56
1:B:274:LYS:HE2	3:B:431:PLP:O4A	2.05	0.56
1:B:15:HIS:HB3	1:B:16:PRO:HD2	1.88	0.56
1:A:220:MET:HB3	1:A:367:VAL:HG21	1.89	0.55
1:A:188:GLU:OE2	1:A:230:ARG:NH2	2.38	0.54
1:B:22:SER:O	1:B:22:SER:OG	2.21	0.54
1:A:233:LYS:N	1:A:233:LYS:HD2	2.22	0.54
1:B:419:ARG:O	1:B:422:GLN:OE1	2.26	0.54
1:B:135:GLN:HB2	5:B:614:HOH:O	2.05	0.53
1:B:125:GLN:HE22	1:B:305:MET:H	1.56	0.52
1:B:187:ASP:O	1:B:190:ASP:HB2	2.10	0.52
1:B:313:ASN:ND2	1:B:316:ALA:H	2.07	0.52
1:A:371:THR:OG1	1:A:372:HIS:CD2	2.63	0.51
1:A:142:ASN:HD22	1:A:177:ALA:HB2	1.76	0.51
1:B:85:HIS:CE1	1:B:88:ALA:HB2	2.45	0.51
1:A:150:GLY:HA3	1:B:149:PHE:HD1	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:ALA:HB3	1:A:87:PRO:HD3	1.93	0.50
1:B:348:ALA:HB3	1:B:349:PRO:HD3	1.93	0.49
1:B:47:ASP:OD2	1:B:50:SER:OG	2.29	0.49
1:B:274:LYS:CE	3:B:431:PLP:O4A	2.60	0.49
1:B:307:GLY:O	4:B:600:IPA:H12	2.11	0.49
1:A:15:HIS:HB3	1:A:16:PRO:HD2	1.94	0.48
1:A:313:ASN:HD22	1:A:316:ALA:CB	2.26	0.48
1:B:23:PRO:HD2	5:B:628:HOH:O	2.12	0.48
1:B:423:ASP:C	1:B:423:ASP:OD2	2.51	0.48
1:B:22:SER:N	1:B:23:PRO:HD3	2.28	0.48
1:A:428:CYS:O	1:A:429:GLN:HB2	2.13	0.48
1:A:43:ARG:HD2	5:A:526:HOH:O	2.14	0.48
1:A:348:ALA:HB3	1:A:349:PRO:HD3	1.96	0.47
1:A:371:THR:OG1	1:A:372:HIS:HD2	1.97	0.47
1:B:52:TRP:CE2	1:B:274:LYS:NZ	2.83	0.47
1:B:140[B]:PHE:HZ	1:B:231:ILE:HD11	1.78	0.47
1:A:427:PHE:O	1:A:428:CYS:HB2	2.15	0.47
1:A:189:ARG:O	1:A:192:VAL:HG22	2.15	0.47
1:A:376:MET:CE	1:A:391:ARG:NH1	2.78	0.46
1:B:77:HIS:HA	1:B:314:PRO:HD2	1.98	0.46
1:B:313:ASN:HD21	1:B:315:LEU:HB3	1.81	0.46
1:B:423:ASP:CG	1:B:425:THR:HG1	2.19	0.46
1:A:274:LYS:CE	3:A:430:PLP:O4A	2.63	0.46
1:B:50:SER:HB2	1:B:402:PRO:HG3	1.98	0.45
1:A:36:GLU:HA	1:A:45:LEU:O	2.16	0.45
1:A:245:ASP:OD1	1:A:245:ASP:C	2.54	0.45
1:B:375:ASN:O	1:B:376:MET:C	2.53	0.45
1:A:93:ARG:NH1	1:A:93:ARG:HG2	2.23	0.45
1:B:127:TRP:CD2	1:B:134:ARG:HD2	2.51	0.45
1:B:312:GLY:O	1:B:313:ASN:C	2.55	0.45
1:A:198:MET:O	1:A:202:ARG:N	2.51	0.44
1:A:77:HIS:CG	1:A:78:VAL:N	2.85	0.44
1:B:77:HIS:CG	1:B:78:VAL:N	2.86	0.44
1:A:90:GLU:OE2	1:A:94:LYS:NZ	2.50	0.44
1:A:127:TRP:CD2	1:A:134:ARG:HD2	2.53	0.43
1:B:379:LEU:HG	1:B:383:PHE:CE1	2.53	0.43
1:B:86:ALA:HB3	1:B:87:PRO:HD3	2.00	0.43
1:A:274:LYS:HZ1	3:A:430:PLP:C3	2.28	0.43
1:A:261:GLU:O	1:A:262:HIS:C	2.57	0.43
1:B:145:HIS:HD2	1:B:245:ASP:OD2	2.01	0.43
1:B:139:THR:O	1:B:140[B]:PHE:HD2	2.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:423:ASP:OD1	1:B:425:THR:OG1	2.36	0.43
1:A:224:HIS:HA	1:A:225:PRO:HD3	1.94	0.42
1:B:47:ASP:CG	1:B:50:SER:HG	2.23	0.42
1:A:364:ALA:O	1:A:400:MET:HA	2.20	0.42
1:B:99:THR:O	1:B:100:PRO:C	2.58	0.42
1:A:348:ALA:N	1:A:349:PRO:CD	2.83	0.41
1:A:312:GLY:O	1:A:313:ASN:C	2.58	0.41
1:B:15:HIS:HB3	1:B:16:PRO:CD	2.50	0.41
1:A:340:GLU:CG	1:A:344:ARG:NH1	2.83	0.41
1:B:52:TRP:CZ2	1:B:274:LYS:NZ	2.88	0.41
1:B:423:ASP:OD2	1:B:425:THR:OG1	2.38	0.41
1:B:224:HIS:HA	1:B:225:PRO:HD3	1.84	0.41
1:B:68:MET:O	1:B:72:ILE:HG13	2.21	0.41
1:A:215:GLN:OE1	1:A:221:ARG:HD3	2.20	0.41
1:B:156:ASP:HA	1:B:157:PRO:HD3	1.88	0.41
1:A:41:ASP:OD1	1:A:41:ASP:C	2.58	0.41
1:B:101:GLN:HB3	1:B:102:PRO:HD3	2.02	0.41
1:B:139:THR:O	1:B:140[B]:PHE:CD2	2.73	0.41
1:A:209:ILE:C	1:A:210:ILE:HG23	2.41	0.41
1:B:233:LYS:HB2	1:B:233:LYS:HE3	1.70	0.41
1:A:125:GLN:NE2	1:A:304:PHE:CD1	2.89	0.40
1:B:114:VAL:O	1:B:118:VAL:HG23	2.21	0.40
1:A:125:GLN:NE2	1:A:304:PHE:HD1	2.19	0.40
1:B:186:TRP:HE1	1:B:188:GLU:HG2	1.85	0.40
1:A:95:LEU:HD23	1:A:95:LEU:HA	1.91	0.40
1:A:145:HIS:HD2	1:A:245:ASP:OD2	2.05	0.40
1:A:357:ALA:O	1:A:358:ASP:HB2	2.21	0.40

All (2) 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:B:237[A]:ARG:NE	1:B:256:LYS:NZ[2_646]	2.07	0.13
1:B:179:GLN:NE2	1:B:422:GLN:NE2[2_546]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone

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	429/429 (100%)	415 (97%)	10 (2%)	4 (1%)	21	29
1	B	428/429 (100%)	411 (96%)	14 (3%)	3 (1%)	26	37
All	All	857/858 (100%)	826 (96%)	24 (3%)	7 (1%)	24	33

All (7) Ramachandran outliers are listed below:

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

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	345/344 (100%)	335 (97%)	10 (3%)	50	70
1	B	344/344 (100%)	331 (96%)	13 (4%)	40	59
All	All	689/688 (100%)	666 (97%)	23 (3%)	44	65

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	76	SER
1	A	80	PHE
1	A	135	GLN
1	A	160	SER
1	A	187	ASP
1	A	188	GLU
1	A	189	ARG
1	A	233	LYS
1	A	290	ARG
1	B	1	MET
1	B	3	THR
1	B	33	GLU
1	B	50	SER
1	B	60	ASN
1	B	76	SER
1	B	80	PHE
1	B	165	TRP
1	B	189	ARG
1	B	192	VAL
1	B	233	LYS
1	B	290	ARG
1	B	422	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	HIS
1	A	125	GLN
1	A	142	ASN
1	A	145	HIS
1	A	201	HIS
1	A	313	ASN
1	A	335	GLN
1	A	342	GLN
1	A	346	GLN
1	A	372	HIS
1	A	408	GLN
1	B	125	GLN
1	B	135	GLN
1	B	142	ASN
1	B	145	HIS
1	B	201	HIS

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Mol	Chain	Res	Type
1	B	262	HIS
1	B	313	ASN
1	B	335	GLN
1	B	342	GLN
1	B	346	GLN
1	B	411	GLN
1	B	422	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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	430	-	16,16,16	3.74	5 (31%)	21,23,23	1.82	5 (23%)
3	PLP	B	431	-	16,16,16	3.57	4 (25%)	21,23,23	1.77	5 (23%)
4	IPA	B	600	-	3,3,3	0.62	0	3,3,3	0.52	0

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	430	-	-	0/8/8/8	0/1/1/1
3	PLP	B	431	-	-	0/8/8/8	0/1/1/1
4	IPA	B	600	-	-	0/0/0/0	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	430	PLP	O3-C3	-5.08	1.25	1.37
3	B	431	PLP	O3-C3	-4.76	1.25	1.37
3	A	430	PLP	P-O3P	-2.19	1.46	1.54
3	A	430	PLP	C4-C5	3.39	1.46	1.42
3	B	431	PLP	C4-C5	4.27	1.47	1.42
3	B	431	PLP	C4-C3	6.03	1.48	1.40
3	A	430	PLP	C4-C3	7.50	1.50	1.40
3	B	431	PLP	C3-C2	10.92	1.48	1.40
3	A	430	PLP	C3-C2	10.94	1.48	1.40

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	430	PLP	C3-C4-C5	-4.21	114.95	118.11
3	B	431	PLP	O4A-C4A-C4	-3.71	117.62	125.11
3	A	430	PLP	O4A-C4A-C4	-2.80	119.45	125.11
3	B	431	PLP	C3-C4-C5	-2.04	116.58	118.11
3	A	430	PLP	O4P-C5A-C5	2.37	112.92	108.99
3	B	431	PLP	C6-N1-C2	2.61	124.61	119.28
3	A	430	PLP	O3P-P-O2P	2.87	118.30	107.38
3	B	431	PLP	O3-C3-C2	3.01	122.89	117.66
3	A	430	PLP	C3-C4-C4A	3.61	125.29	119.84
3	B	431	PLP	O3P-P-O2P	3.83	121.95	107.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	430	PLP	9	0
3	B	431	PLP	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	600	IPA	3	0

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

6.1 Protein, DNA and RNA chains

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, 95^{th} 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	429/429 (100%)	-0.15	17 (3%)	42	42	15, 27, 53, 64	7 (1%)
1	B	427/429 (99%)	-0.11	16 (3%)	45	45	15, 28, 52, 67	3 (0%)
All	All	856/858 (99%)	-0.13	33 (3%)	43	43	15, 27, 53, 67	10 (1%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	ALA	5.6
1	A	183	ASP	5.5
1	A	164	LEU	5.2
1	A	163	SER	4.7
1	B	164	LEU	4.2
1	A	302	GLY	3.9
1	B	428	CYS	3.6
1	B	189	ARG	3.5
1	B	165	TRP	3.5
1	A	299	GLY	3.3
1	A	181	ARG	3.3
1	B	133	ALA	3.0
1	B	183	ASP	2.9
1	B	163	SER	2.9
1	B	187	ASP	2.8
1	A	189	ARG	2.7
1	A	168	TYR	2.7
1	A	428	CYS	2.6
1	B	179	GLN	2.5
1	B	161	MET	2.5
1	A	184	GLY	2.5
1	A	192	VAL	2.5
1	B	372	HIS	2.3
1	B	203	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	300	GLU	2.3
1	A	429	GLN	2.2
1	A	186	TRP	2.2
1	A	237	ARG	2.2
1	B	237[A]	ARG	2.2
1	A	179	GLN	2.1
1	B	167	GLY	2.1
1	A	133	ALA	2.0
1	B	377	ALA	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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, 95^{th} 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
2	NA	A	501	1/1	0.94	0.21	4.94	10,10,10,10	0
2	NA	B	502	1/1	0.92	0.26	3.37	14,14,14,14	0
4	IPA	B	600	4/4	0.79	0.26	2.96	28,28,28,29	0
3	PLP	A	430	16/16	0.96	0.12	0.77	11,18,23,31	0
3	PLP	B	431	16/16	0.97	0.11	0.47	13,22,25,29	0

6.5 Other polymers

There are no such residues in this entry.