



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:44 AM UTC

PDB ID : 1S07 / pdb_00001s07
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

<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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

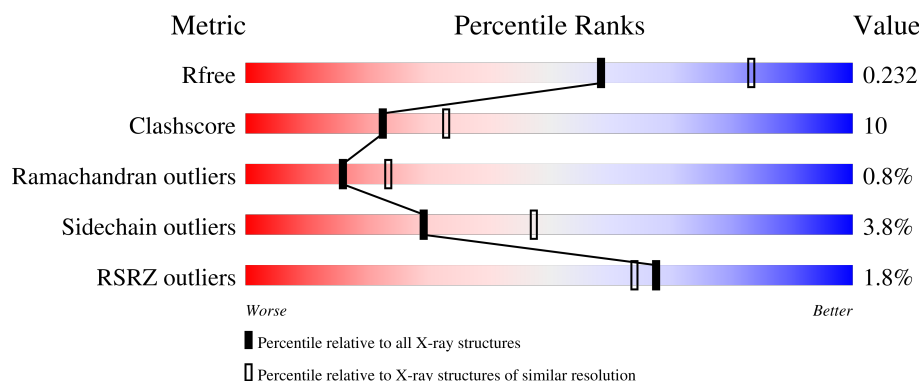
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (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 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	429	% 78% 20% .
1	B	429	2% 76% 21% .

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
3	PLP	A	430	-	-	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 mutation	UNP P12995
B	14	LEU	TRP	SEE REMARK 999	UNP P12995
B	253	ALA	ARG	engineered mutation	UNP P12995

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

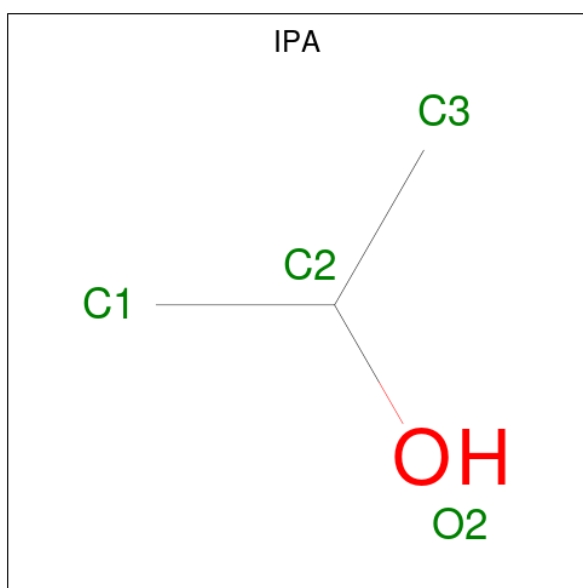
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		
2	B	1	Total	Na	0	0
			1	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C₈H₁₀NO₆P).



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 (CCD ID: IPA) (formula: C_3H_8O).



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

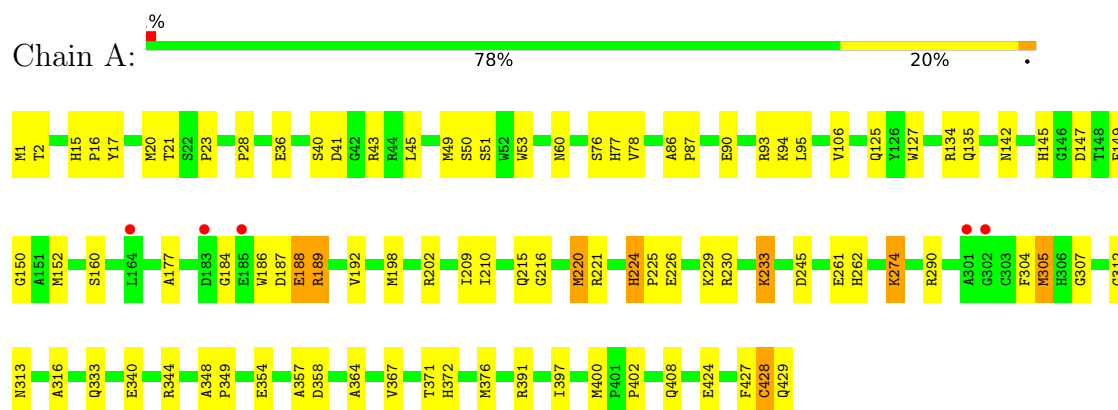
- Molecule 5 is water.

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

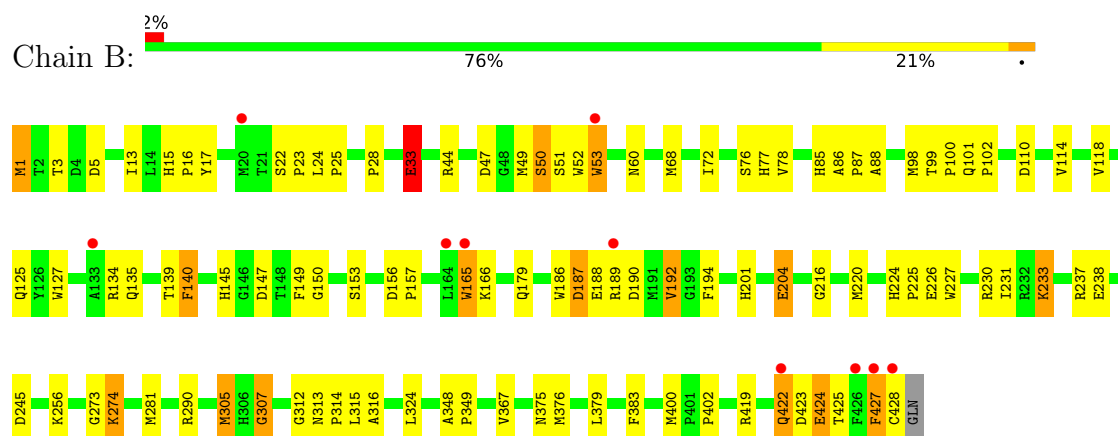
3 Residue-property plots

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: Adenosylmethionine-8-amino-7-oxononanoate aminotransferase



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4 Data and refinement statistics

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) 89.1 (20.00-2.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.188 , 0.224 0.195 , 0.232	Depositor DCC
R_{free} test set	1413 reflections (5.27%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	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.*

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: NA, PLP, IPA

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	0.98	5/3403 (0.1%)	1.13	8/4620 (0.2%)
1	B	0.98	3/3391 (0.1%)	1.11	10/4604 (0.2%)
All	All	0.98	8/6794 (0.1%)	1.12	18/9224 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	220	MET	SD-CE	-8.54	1.58	1.79
1	B	305	MET	SD-CE	6.60	1.96	1.79
1	A	20	MET	SD-CE	-6.25	1.64	1.79
1	B	220	MET	SD-CE	-6.24	1.64	1.79
1	A	152	MET	SD-CE	-5.99	1.64	1.79
1	B	187	ASP	CA-C	5.45	1.58	1.52
1	A	397	ILE	CA-CB	-5.30	1.47	1.53
1	A	305	MET	SD-CE	5.13	1.92	1.79

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	184	GLY	N-CA-C	10.19	125.79	112.13
1	A	333	GLN	CB-CG-CD	9.01	127.92	112.60
1	B	427	PHE	N-CA-C	6.72	120.48	109.72
1	B	140[A]	PHE	CB-CA-C	-6.60	96.32	109.33
1	B	140[B]	PHE	CB-CA-C	-6.60	96.32	109.33
1	B	33	GLU	CB-CG-CD	6.42	123.52	112.60
1	B	53	TRP	N-CA-C	-6.14	105.64	112.57
1	A	106	VAL	N-CA-C	6.05	116.64	108.17
1	B	13	ILE	N-CA-C	5.76	115.90	107.37
1	B	110	ASP	N-CA-C	5.61	120.25	112.90
1	A	149	PHE	N-CA-C	5.57	117.03	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	ASP	N-CA-C	-5.55	105.13	111.07
1	A	408	GLN	CA-CB-CG	5.54	125.17	114.10
1	A	305	MET	N-CA-C	5.53	118.14	108.52
1	B	367	VAL	N-CA-C	5.27	115.49	108.11
1	A	224	HIS	N-CA-C	5.16	117.09	109.50
1	A	2	THR	N-CA-C	5.05	117.92	110.59
1	B	192	VAL	N-CA-CB	5.02	115.89	110.62

There are no chirality outliers.

There are no planarity outliers.

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	3314	0	3265	65	0
1	B	3303	0	3260	72	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	137	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 10.

All (137) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:TRP:HE1	1:A:188:GLU:HG2	1.40	0.85
1:B:139:THR:C	1:B:140[B]:PHE:HD2	1.84	0.85
1:B:51:SER:HB2	1:B:400:MET:HG2	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: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:HH11	1:A:93:ARG:HG2	1.51	0.74
1:A:49:MET:HB3	1:A:53:TRP:HZ3	1.52	0.74
1:A:274:LYS:NZ	3:A:430:PLP:C4	2.50	0.74
1:B:139:THR:C	1:B:140[B]:PHE:CD2	2.66	0.73
1:A:49:MET:HB3	1:A:53:TRP:CZ3	2.24	0.72
1:A:340:GLU:CD	1:A:344:ARG:HH12	1.98	0.71
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:A:340:GLU:CD	1:A:344:ARG:NH1	2.50	0.70
1:A:230:ARG:HD2	5:A:542:HOH:O	1.92	0.68
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:HE2	1:B:194:PHE:CE1	2.16	0.62
1:B:140[B]:PHE:CZ	1:B:231:ILE:HD11	2.34	0.62
1:B:281:MET:HE1	1:B:315:LEU:HG	1.82	0.61
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:423:ASP:CG	1:B:425:THR:HG1	2.12	0.57
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:1:MET:HE3	1:B:28:PRO:HB2	1.87	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HD2	1:A:233:LYS:N	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:423:ASP:C	1:B:423:ASP:OD2	2.51	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:1:MET:HE3	1:A:40:SER:OG	2.11	0.50
1:B:375:ASN:O	1:B:376:MET:C	2.53	0.50
1:A:86:ALA:HB3	1:A:87:PRO:HD3	1.93	0.50
1:A:150:GLY:HA3	1:B:149:PHE:HD1	1.76	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:B:312:GLY:O	1:B:313:ASN:C	2.55	0.49
1:A:15:HIS:HB3	1:A:16:PRO:HD2	1.94	0.48
1:A:17[A]:TYR:OH	1:A:147:ASP:OD2	2.28	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: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:261:GLU:O	1:A:262:HIS:C	2.57	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:A:189:ARG:O	1:A:192:VAL:HG22	2.15	0.47
1:A:427:PHE:O	1:A:428:CYS:HB2	2.15	0.47
1:B:140[B]:PHE:HZ	1:B:231:ILE:HD11	1.78	0.47
1:A:245:ASP:C	1:A:245:ASP:OD1	2.54	0.46
1:A:376:MET:CE	1:A:391:ARG:NH1	2.78	0.46
1:B:99:THR:O	1:B:100:PRO:C	2.58	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HE2	1:B:1:MET:HB2	1.72	0.46
1:A:274:LYS:CE	3:A:430:PLP:O4A	2.63	0.46
1:A:312:GLY:O	1:A:313:ASN:C	2.58	0.46
1:A:36:GLU:HA	1:A:45:LEU:O	2.16	0.45
1:A:41:ASP:C	1:A:41:ASP:OD1	2.58	0.45
1:B:50:SER:HB2	1:B:402:PRO:HG3	1.98	0.45
1:A:51:SER:HB3	1:A:400:MET:HE2	1.98	0.45
1:A:209:ILE:C	1:A:210:ILE:HG23	2.41	0.45
1:A:93:ARG:HG2	1:A:93:ARG:NH1	2.23	0.45
1:B:127:TRP:CD2	1:B:134:ARG:HD2	2.51	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:376:MET:HE1	1:A:391:ARG:NH1	2.33	0.44
1:A:90:GLU:OE2	1:A:94:LYS:NZ	2.50	0.44
1:B:17:TYR:OH	1:B:147:ASP:OD2	2.27	0.44
1:B:150:GLY:O	1:B:153:SER:OG	2.26	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: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
1:B:24:LEU:O	1:B:25:PRO:C	2.62	0.42
1:B:190:ASP:CG	1:B:227:TRP:HE1	2.27	0.42
1:A:224:HIS:HA	1:A:225:PRO:HD3	1.94	0.42
1:B:165:TRP:O	1:B:166:LYS:C	2.63	0.42
1:A:21:THR:C	1:A:23:PRO:HD3	2.45	0.42
1:B:273:GLY:O	1:B:274:LYS:C	2.62	0.42
1:B:423:ASP:OD1	1:B:425:THR:OG1	2.36	0.42
1:A:364:ALA:O	1:A:400:MET:HA	2.20	0.42
1:B:424:GLU:O	1:B:425:THR:C	2.63	0.42
1:A:348:ALA:N	1:A:349:PRO:CD	2.83	0.41
1:B:15:HIS:HB3	1:B:16:PRO:CD	2.50	0.41
1:A:1:MET:HE3	1:A:28:PRO:HB2	2.03	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: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:98:MET:HE1	1:B:324:LEU:HD22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:ASP:HA	1:B:157:PRO:HD3	1.88	0.41
1:B:101:GLN:HB3	1:B:102:PRO:HD3	2.02	0.41
1:B:233:LYS:HE3	1:B:233:LYS:HB2	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:A:95:LEU:HD23	1:A:95:LEU:HA	1.91	0.40
1:B:186:TRP:HE1	1:B:188:GLU:HG2	1.85	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%)	14	20
1	B	428/429 (100%)	411 (96%)	14 (3%)	3 (1%)	18	27
All	All	857/858 (100%)	826 (96%)	24 (3%)	7 (1%)	16	23

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	274	LYS
1	A	274	LYS

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Mol	Chain	Res	Type
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%)	333 (96%)	12 (4%)	32	51
1	B	344/344 (100%)	330 (96%)	14 (4%)	27	44
All	All	689/688 (100%)	663 (96%)	26 (4%)	29	47

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	76	SER
1	A	135	GLN
1	A	160	SER
1	A	187	ASP
1	A	188	GLU
1	A	189	ARG
1	A	229	LYS
1	A	233	LYS
1	A	290	ARG
1	A	354	GLU
1	A	424	GLU
1	B	1	MET
1	B	3	THR
1	B	33	GLU
1	B	50	SER
1	B	60	ASN
1	B	76	SER

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Mol	Chain	Res	Type
1	B	165	TRP
1	B	189	ARG
1	B	192	VAL
1	B	204	GLU
1	B	233	LYS
1	B	290	ARG
1	B	422	GLN
1	B	424	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	71	GLN
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	334	GLN
1	A	335	GLN
1	A	342	GLN
1	A	346	GLN
1	A	372	HIS
1	A	408	GLN
1	A	411	GLN
1	A	418	ASN
1	B	125	GLN
1	B	135	GLN
1	B	142	ASN
1	B	145	HIS
1	B	201	HIS
1	B	313	ASN
1	B	334	GLN
1	B	335	GLN
1	B	342	GLN
1	B	346	GLN
1	B	418	ASN
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 oligosaccharides 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	2.84	5 (31%)	20,23,23	2.06	6 (30%)
4	IPA	B	600	-	3,3,3	0.61	0	3,3,3	0.54	0
3	PLP	B	431	-	16,16,16	2.72	4 (25%)	20,23,23	2.27	6 (30%)

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	-	-	1/8/8/8	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	430	PLP	C3-C2	7.10	1.48	1.41
3	B	431	PLP	C3-C2	7.09	1.48	1.41
3	A	430	PLP	C4-C3	5.52	1.50	1.41
3	A	430	PLP	O3-C3	-5.15	1.25	1.36
3	B	431	PLP	O3-C3	-4.82	1.25	1.36
3	B	431	PLP	C4-C3	4.40	1.48	1.41
3	B	431	PLP	C4-C5	4.15	1.47	1.42
3	A	430	PLP	C4-C5	3.29	1.46	1.42
3	A	430	PLP	P-O3P	-2.14	1.46	1.54

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	431	PLP	C4-C3-C2	-6.45	116.51	120.14
3	A	430	PLP	C3-C4-C5	-4.14	114.95	118.28
3	A	430	PLP	C3-C4-C4A	3.96	125.29	119.84
3	A	430	PLP	C4-C3-C2	-3.88	117.96	120.14
3	B	431	PLP	O3P-P-O2P	3.77	121.95	107.80
3	B	431	PLP	O4A-C4A-C4	-3.03	117.62	124.80
3	B	431	PLP	C6-N1-C2	2.98	124.61	119.20
3	A	430	PLP	O3P-P-O2P	2.80	118.30	107.80
3	B	431	PLP	O3-C3-C2	2.56	122.89	117.58
3	A	430	PLP	C6-N1-C2	2.28	123.33	119.20
3	A	430	PLP	O4A-C4A-C4	-2.26	119.45	124.80
3	B	431	PLP	C3-C4-C5	-2.11	116.58	118.28

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	431	PLP	C4-C5-C5A-O4P

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
4	B	600	IPA	3	0
3	B	431	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	429/429 (100%)	-0.18	5 (1%) 76 74	11, 27, 52, 64	9 (2%)
1	B	427/429 (99%)	-0.14	10 (2%) 61 57	15, 27, 52, 67	5 (1%)
All	All	856/858 (99%)	-0.16	15 (1%) 67 64	11, 27, 53, 67	14 (1%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	164	LEU	3.6
1	B	428	CYS	3.5
1	A	183	ASP	3.4
1	A	301	ALA	3.3
1	A	302	GLY	2.9
1	A	164	LEU	2.8
1	B	165	TRP	2.6
1	B	20	MET	2.5
1	B	53	TRP	2.5
1	B	427	PHE	2.3
1	B	133	ALA	2.3
1	B	189	ARG	2.3
1	B	426	PHE	2.1
1	B	422	GLN	2.1
1	A	185	GLU	2.0

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 oligosaccharides 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. 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
4	IPA	B	600	4/4	0.77	0.15	28,28,28,29	0
2	NA	B	502	1/1	0.93	0.20	14,14,14,14	0
2	NA	A	501	1/1	0.94	0.16	10,10,10,10	0
3	PLP	A	430	16/16	0.96	0.07	11,18,23,31	0
3	PLP	B	431	16/16	0.97	0.07	13,22,25,29	0

6.5 Other polymers

There are no such residues in this entry.