



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:45 AM UTC

PDB ID : 5EAA / pdb_00005eaa
Title : ASPARTATE AMINOTRANSFERASE FROM E. COLI, C191S MUTATION
Authors : Jeffery, C.J.; Gloss, L.M.; Petsko, G.A.; Ringe, D.
Deposited on : 1998-12-29
Resolution : 2.40 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

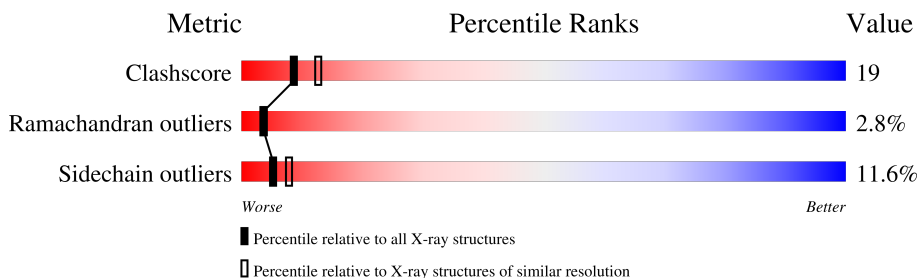
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	396	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3224 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
			Total	C	N	O	S			
1	A	396	3069	1936	536	585	12	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	191	SER	CYS	engineered mutation	UNP P00509

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



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

- Molecule 3 is water.

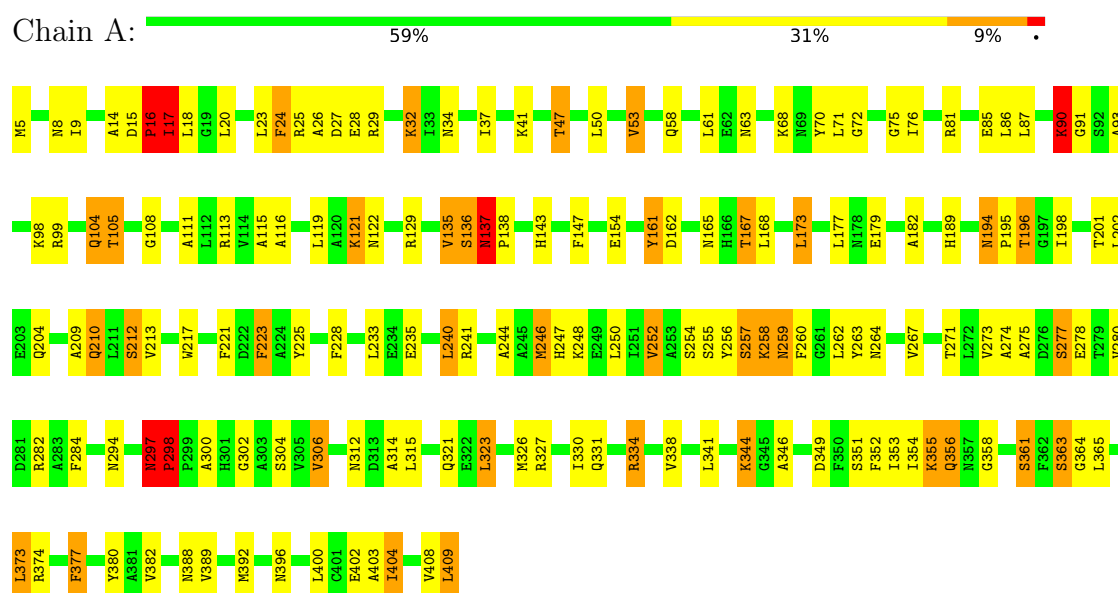
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	140	Total 140	O 140	0	0

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. 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. 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



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	156.27 Å 86.94 Å 79.77 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40	Depositor
% Data completeness (in resolution range)	92.0 (10.00-2.40)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.219 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3224	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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	1.05	1/3130 (0.0%)	1.36	40/4240 (0.9%)

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	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	258	LYS	C-N	28.70	1.76	1.33

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	SER	O-C-N	-12.44	107.12	122.17
1	A	257	SER	CA-C-N	12.10	137.47	120.29
1	A	257	SER	C-N-CA	12.10	137.47	120.29
1	A	356	GLN	N-CA-C	-11.99	87.46	108.69
1	A	47	THR	CA-C-N	10.26	130.19	120.03
1	A	47	THR	C-N-CA	10.26	130.19	120.03
1	A	137	ASN	CA-C-N	-10.23	108.79	119.83
1	A	137	ASN	C-N-CA	-10.23	108.79	119.83
1	A	17	ILE	N-CA-C	-9.05	90.52	109.34
1	A	87	LEU	N-CA-C	7.75	119.51	111.14
1	A	196	THR	N-CA-C	7.60	122.65	113.23
1	A	297	ASN	CA-C-N	7.49	128.09	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	297	ASN	C-N-CA	7.49	128.09	120.38
1	A	194	ASN	CA-C-N	-7.42	110.57	119.84
1	A	194	ASN	C-N-CA	-7.42	110.57	119.84
1	A	161	TYR	N-CA-C	7.41	121.76	109.46
1	A	24	PHE	N-CA-C	-7.37	102.41	111.40
1	A	298	PRO	N-CA-C	7.34	119.65	110.70
1	A	248	LYS	N-CA-C	-7.03	103.69	112.90
1	A	9	ILE	N-CA-C	6.97	118.43	109.30
1	A	70	TYR	N-CA-C	6.94	120.02	110.24
1	A	377	PHE	N-CA-C	6.77	121.14	113.01
1	A	72	GLY	N-CA-C	-6.51	102.12	112.66
1	A	228	PHE	N-CA-C	6.47	121.25	113.23
1	A	121	LYS	N-CA-C	6.17	121.09	113.50
1	A	344	LYS	N-CA-C	6.13	120.37	113.02
1	A	334	ARG	N-CA-C	-6.12	104.53	111.14
1	A	104	GLN	N-CA-C	-6.04	101.35	110.28
1	A	63	ASN	N-CA-C	5.95	120.54	113.16
1	A	58	GLN	N-CA-C	-5.80	104.55	111.69
1	A	355	LYS	CA-C-N	5.78	132.05	122.33
1	A	355	LYS	C-N-CA	5.78	132.05	122.33
1	A	71	LEU	N-CA-C	-5.67	102.53	110.35
1	A	264	ASN	N-CA-C	5.65	119.88	113.15
1	A	194	ASN	N-CA-C	5.64	120.58	113.25
1	A	300	ALA	N-CA-C	5.57	118.28	111.82
1	A	306	VAL	N-CA-C	5.52	117.99	111.09
1	A	364	GLY	N-CA-C	-5.33	107.67	115.72
1	A	115	ALA	N-CA-C	-5.33	105.55	111.36
1	A	93	ALA	N-CA-C	-5.28	105.42	111.07

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	225	TYR	Sidechain
1	A	257	SER	Mainchain

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	3069	0	3015	115	0
2	A	15	0	6	0	0
3	A	140	0	0	6	0
All	All	3224	0	3021	115	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 19.

All (115) 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:258:LYS:C	1:A:259:ASN:N	1.76	1.44
1:A:129:ARG:HD3	1:A:154:GLU:HB3	1.51	0.91
1:A:25:ARG:NE	1:A:25:ARG:HA	1.89	0.87
1:A:408:VAL:O	1:A:409:LEU:HB2	1.79	0.82
1:A:37:ILE:HG13	1:A:388:ASN:HD22	1.47	0.79
1:A:27:ASP:HB3	1:A:32:LYS:HZ3	1.49	0.77
1:A:50:LEU:HB2	1:A:53:VAL:HG13	1.65	0.77
1:A:90:LYS:HD3	1:A:90:LYS:H	1.50	0.75
1:A:338:VAL:HG21	1:A:354:ILE:HG13	1.69	0.73
1:A:17:ILE:HG23	1:A:18:LEU:H	1.56	0.70
1:A:323:LEU:HA	1:A:326:MET:HE3	1.75	0.69
1:A:392:MET:HE2	1:A:400:LEU:HD21	1.74	0.68
1:A:252:VAL:HG13	1:A:271:THR:HB	1.77	0.67
1:A:244:ALA:HA	1:A:250:LEU:HD21	1.78	0.66
1:A:76:ILE:HG13	1:A:104:GLN:HE22	1.59	0.66
1:A:24:PHE:CE1	1:A:32:LYS:HG2	2.31	0.64
1:A:201:THR:H	1:A:204:GLN:HE21	1.47	0.63
1:A:344:LYS:HD3	1:A:402:GLU:HG3	1.81	0.62
1:A:99:ARG:HD2	1:A:274:ALA:O	1.99	0.62
1:A:321:GLN:HG3	3:A:578:HOH:O	1.97	0.62
1:A:330:ILE:O	1:A:334:ARG:HG3	2.00	0.61
1:A:14:ALA:C	1:A:16:PRO:HD2	2.25	0.61
1:A:137:ASN:HB3	1:A:138:PRO:CD	2.31	0.60
1:A:177:LEU:O	1:A:217:TRP:HZ2	1.82	0.60
1:A:404:ILE:O	1:A:408:VAL:HG22	2.01	0.60
1:A:15:ASP:O	1:A:17:ILE:HG22	2.02	0.60
1:A:99:ARG:HD3	1:A:275:ALA:O	2.02	0.59
1:A:37:ILE:HG13	1:A:388:ASN:ND2	2.17	0.59
1:A:8:ASN:HD22	1:A:8:ASN:H	1.51	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:THR:HG21	1:A:111:ALA:N	2.17	0.58
1:A:338:VAL:CG2	1:A:353:ILE:HB	2.34	0.58
1:A:201:THR:H	1:A:204:GLN:NE2	2.03	0.56
1:A:27:ASP:HB3	1:A:32:LYS:NZ	2.18	0.56
1:A:15:ASP:HB3	1:A:18:LEU:HB2	1.87	0.56
1:A:25:ARG:HA	1:A:25:ARG:HE	1.70	0.55
1:A:99:ARG:HD3	1:A:275:ALA:C	2.33	0.53
1:A:17:ILE:HG23	1:A:18:LEU:N	2.24	0.52
1:A:209:ALA:O	1:A:213:VAL:HG23	2.10	0.52
1:A:162:ASP:OD2	1:A:165:ASN:HB2	2.10	0.52
1:A:382:VAL:HG11	3:A:583:HOH:O	2.10	0.52
1:A:143:HIS:O	1:A:147:PHE:CD2	2.63	0.51
1:A:338:VAL:HG11	1:A:351:SER:HA	1.92	0.51
1:A:212:SER:HB3	1:A:247:HIS:CE1	2.45	0.51
1:A:47:THR:HG22	3:A:540:HOH:O	2.10	0.51
1:A:8:ASN:H	1:A:8:ASN:ND2	2.07	0.51
1:A:34:ASN:OD1	1:A:37:ILE:HG23	2.10	0.51
1:A:380:TYR:CD1	1:A:380:TYR:N	2.78	0.51
1:A:334:ARG:HH22	1:A:361:SER:HB3	1.75	0.51
1:A:338:VAL:HG22	1:A:353:ILE:HB	1.94	0.50
1:A:196:THR:HB	1:A:198:ILE:HG13	1.94	0.49
1:A:129:ARG:CD	1:A:154:GLU:HB3	2.35	0.48
1:A:177:LEU:O	1:A:217:TRP:CZ2	2.65	0.48
1:A:297:ASN:HB2	1:A:298:PRO:HD2	1.95	0.48
1:A:352:PHE:O	1:A:356:GLN:HG3	2.13	0.48
1:A:223:PHE:O	1:A:254:SER:HA	2.14	0.47
1:A:294:ASN:C	1:A:294:ASN:HD22	2.21	0.47
1:A:161:TYR:CE1	1:A:198:ILE:HD11	2.49	0.47
1:A:327:ARG:O	1:A:331:GLN:HG3	2.14	0.47
1:A:256:TYR:HD2	1:A:267:VAL:HG12	1.81	0.46
1:A:105:THR:HG21	1:A:111:ALA:CA	2.45	0.46
1:A:356:GLN:OE1	1:A:361:SER:HB2	2.16	0.46
1:A:355:LYS:HA	1:A:355:LYS:HD3	1.56	0.46
1:A:221:PHE:CD1	1:A:240:LEU:HD23	2.50	0.46
1:A:373:LEU:HD11	1:A:403:ALA:O	2.16	0.46
1:A:338:VAL:HG12	1:A:338:VAL:O	2.16	0.46
1:A:68:LYS:HG2	3:A:600:HOH:O	2.16	0.45
1:A:135:VAL:HG23	1:A:136:SER:O	2.16	0.45
1:A:24:PHE:CE1	1:A:34:ASN:HB2	2.52	0.45
1:A:37:ILE:CD1	1:A:41:LYS:HE2	2.47	0.45
1:A:377:PHE:CG	1:A:403:ALA:HB1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLY:HA3	1:A:255:SER:HB2	1.97	0.45
1:A:91:GLY:HA3	3:A:548:HOH:O	2.17	0.45
1:A:37:ILE:HD13	1:A:41:LYS:HE2	1.99	0.45
1:A:29:ARG:O	1:A:32:LYS:HB2	2.17	0.44
1:A:168:LEU:HD11	1:A:173:LEU:HD12	1.98	0.44
1:A:105:THR:HG21	1:A:111:ALA:HB2	2.00	0.44
1:A:334:ARG:HG2	1:A:389:VAL:HG11	1.98	0.44
1:A:98:LYS:HB3	1:A:98:LYS:HE2	1.76	0.43
1:A:121:LYS:HG2	1:A:122:ASN:ND2	2.32	0.43
1:A:15:ASP:N	1:A:16:PRO:HD2	2.32	0.43
1:A:75:GLY:HA3	1:A:104:GLN:HB3	2.00	0.43
1:A:161:TYR:HD1	1:A:198:ILE:HD12	1.83	0.43
1:A:167:THR:HG23	1:A:168:LEU:N	2.33	0.43
1:A:210:GLN:HG2	1:A:246:MET:HE1	2.01	0.43
1:A:258:LYS:C	1:A:259:ASN:CA	2.80	0.43
1:A:137:ASN:O	1:A:138:PRO:C	2.59	0.42
1:A:177:LEU:HB3	1:A:217:TRP:CH2	2.54	0.42
1:A:306:VAL:O	1:A:306:VAL:HG12	2.19	0.42
1:A:28:GLU:O	1:A:29:ARG:C	2.61	0.42
1:A:86:LEU:HD11	1:A:233:LEU:HD22	2.02	0.42
1:A:363:SER:HB2	1:A:365:LEU:HG	2.00	0.42
1:A:137:ASN:HB3	1:A:138:PRO:HD2	2.02	0.42
1:A:113:ARG:O	1:A:116:ALA:HB3	2.18	0.42
1:A:312:ASN:OD1	1:A:314:ALA:HB3	2.20	0.42
1:A:330:ILE:HG21	1:A:358:GLY:O	2.20	0.41
1:A:25:ARG:NE	1:A:25:ARG:CA	2.71	0.41
1:A:402:GLU:HG2	3:A:535:HOH:O	2.19	0.41
1:A:76:ILE:H	1:A:104:GLN:HE21	1.69	0.41
1:A:81:ARG:O	1:A:85:GLU:HG3	2.20	0.41
1:A:262:LEU:O	1:A:263:TYR:C	2.63	0.41
1:A:302:GLY:O	1:A:306:VAL:HG23	2.21	0.41
1:A:294:ASN:C	1:A:294:ASN:ND2	2.79	0.41
1:A:27:ASP:C	1:A:32:LYS:HZ1	2.29	0.41
1:A:247:HIS:HB2	1:A:250:LEU:HD23	2.03	0.41
1:A:396:ASN:OD1	1:A:396:ASN:C	2.64	0.41
1:A:260:PHE:HB3	1:A:262:LEU:HD12	2.03	0.41
1:A:277:SER:O	1:A:280:VAL:HG12	2.21	0.41
1:A:373:LEU:HD13	1:A:408:VAL:HG11	2.02	0.40
1:A:15:ASP:N	1:A:16:PRO:CD	2.84	0.40
1:A:278:GLU:CD	1:A:282:ARG:HH21	2.29	0.40
1:A:280:VAL:O	1:A:284:PHE:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:TYR:N	1:A:380:TYR:HD1	2.18	0.40
1:A:373:LEU:HD12	1:A:373:LEU:HA	1.82	0.40
1:A:76:ILE:HG13	1:A:104:GLN:NE2	2.33	0.40
1:A:189:HIS:CD2	1:A:194:ASN:H	2.40	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	394/396 (100%)	355 (90%)	28 (7%)	11 (3%)	4 3

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	PRO
1	A	17	ILE
1	A	137	ASN
1	A	346	ALA
1	A	26	ALA
1	A	32	LYS
1	A	90	LYS
1	A	182	ALA
1	A	361	SER
1	A	136	SER
1	A	195	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	320/320 (100%)	283 (88%)	37 (12%)	5 8

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	MET
1	A	16	PRO
1	A	20	LEU
1	A	23	LEU
1	A	53	VAL
1	A	61	LEU
1	A	90	LYS
1	A	105	THR
1	A	119	LEU
1	A	135	VAL
1	A	167	THR
1	A	173	LEU
1	A	179	GLU
1	A	202	LEU
1	A	210	GLN
1	A	212	SER
1	A	223	PHE
1	A	235	GLU
1	A	240	LEU
1	A	241	ARG
1	A	246	MET
1	A	252	VAL
1	A	259	ASN
1	A	273	VAL
1	A	277	SER
1	A	297	ASN
1	A	298	PRO
1	A	304	SER
1	A	315	LEU
1	A	323	LEU
1	A	341	LEU
1	A	349	ASP
1	A	363	SER
1	A	373	LEU

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Mol	Chain	Res	Type
1	A	374	ARG
1	A	404	ILE
1	A	409	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	84	GLN
1	A	96	ASN
1	A	104	GLN
1	A	122	ASN
1	A	137	ASN
1	A	148	ASN
1	A	175	ASN
1	A	178	ASN
1	A	189	HIS
1	A	204	GLN
1	A	259	ASN
1	A	294	ASN
1	A	297	ASN
1	A	328	GLN
1	A	339	ASN
1	A	347	ASN
1	A	357	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

1 ligand is 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	458	1	15,15,16	2.73	5 (33%)	21,22,23	2.40	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	458	1	-	0/6/6/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	458	PLP	O4P-C5A	-7.68	1.16	1.44
2	A	458	PLP	C4A-C4	-3.72	1.44	1.51
2	A	458	PLP	P-O4P	-3.71	1.48	1.60
2	A	458	PLP	P-O3P	-3.33	1.42	1.54
2	A	458	PLP	P-O1P	-2.13	1.43	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	458	PLP	O4P-P-O1P	-5.52	91.52	106.44
2	A	458	PLP	C6-C5-C4	4.50	121.78	118.10
2	A	458	PLP	O2P-P-O4P	3.64	116.15	106.67
2	A	458	PLP	C3-C2-N1	-2.90	117.30	120.96
2	A	458	PLP	C5-C6-N1	-2.89	119.13	123.83
2	A	458	PLP	C4A-C4-C5	2.39	123.40	120.94
2	A	458	PLP	C6-N1-C2	2.35	123.46	119.20
2	A	458	PLP	C2A-C2-C3	2.19	123.36	120.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	258:LYS	C	259:ASN	N	1.76

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.