



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:28 PM UTC

PDB ID : 1G4V / pdb_00001g4v
Title : ASPARTATE AMINOTRANSFERASE ACTIVE SITE MUTANT
N194A/Y225F
Authors : Mizuguchi, H.; Hayashi, H.; Okada, K.; Miyahara, I.; Hirotsu, K.;
Kagamiyama, H.
Deposited on : 2000-10-28
Resolution : 2.00 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

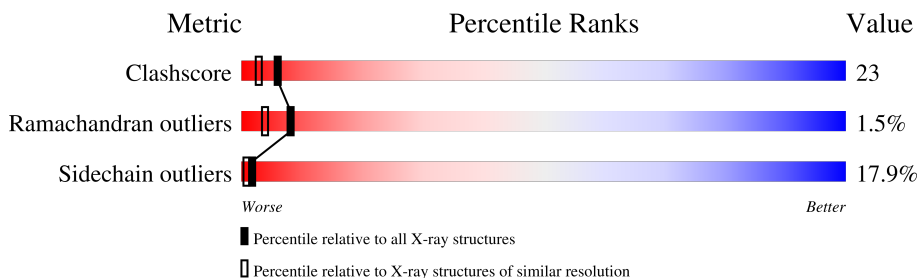
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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	50% 37% 12%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3295 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.

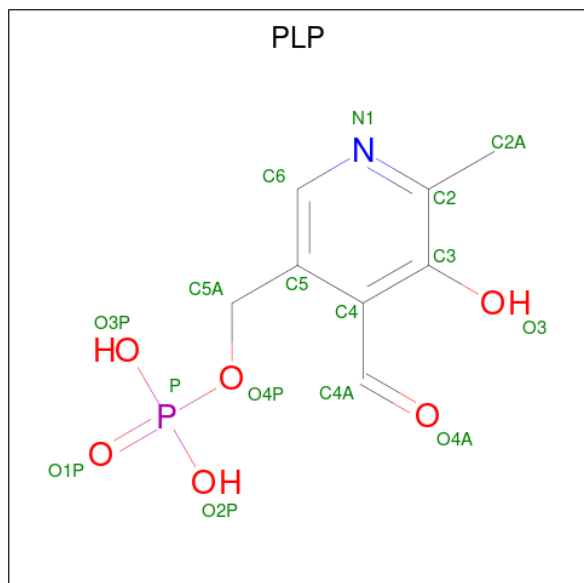
- Molecule 1 is a protein called ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	396	3065	1935	535	582	13	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	194	ALA	ASN	engineered mutation	UNP P00509
A	225	PHE	TYR	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.

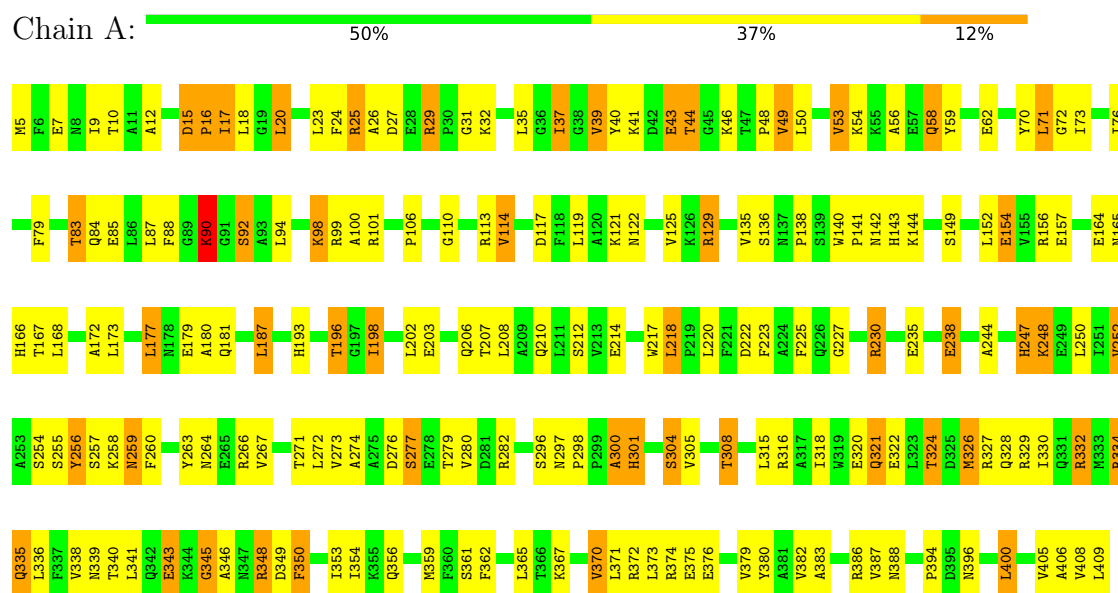
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	215	Total 215	O 215	0	0

3 Residue-property plots [i](#)

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	155.30Å 87.10Å 79.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.209 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3295	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	0.79	1/3126 (0.0%)	1.23	34/4234 (0.8%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	326	MET	SD-CE	-7.45	1.60	1.79

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	THR	N-CA-C	8.63	122.30	111.69
1	A	298	PRO	N-CA-C	8.35	120.89	110.70
1	A	300	ALA	N-CA-C	8.04	120.12	111.36
1	A	301	HIS	N-CA-C	7.90	120.66	111.02
1	A	297	ASN	N-CA-C	-7.80	101.20	108.22
1	A	140	TRP	CA-C-N	7.49	129.20	119.84
1	A	140	TRP	C-N-CA	7.49	129.20	119.84
1	A	15	ASP	CA-C-N	7.39	129.08	119.84
1	A	15	ASP	C-N-CA	7.39	129.08	119.84
1	A	37	ILE	N-CA-C	7.16	122.19	111.89
1	A	247	HIS	N-CA-C	7.14	121.68	110.32
1	A	334	ARG	N-CA-C	-6.67	103.93	111.07
1	A	238	GLU	N-CA-C	6.64	118.52	111.28
1	A	90	LYS	N-CA-C	6.35	119.02	111.33
1	A	31	GLY	N-CA-C	-6.30	104.65	115.61
1	A	58	GLN	N-CA-C	-6.29	104.50	111.36
1	A	350	PHE	N-CA-C	-5.92	104.45	112.26
1	A	345	GLY	N-CA-C	5.86	127.07	113.18
1	A	166	HIS	N-CA-C	-5.79	103.26	111.24
1	A	340	THR	N-CA-C	-5.77	105.73	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	12	ALA	CA-C-N	5.60	125.55	119.78
1	A	12	ALA	C-N-CA	5.60	125.55	119.78
1	A	298	PRO	CA-C-N	5.59	126.47	119.98
1	A	298	PRO	C-N-CA	5.59	126.47	119.98
1	A	70	TYR	N-CA-C	5.54	117.44	110.24
1	A	248	LYS	N-CA-C	-5.46	104.97	111.69
1	A	71	LEU	N-CA-C	-5.36	102.59	110.52
1	A	256	TYR	N-CA-C	-5.35	105.58	113.61
1	A	222	ASP	N-CA-C	-5.33	98.56	107.99
1	A	343	GLU	N-CA-C	-5.32	106.47	113.12
1	A	264	ASN	N-CA-C	5.31	119.62	113.20
1	A	122	ASN	N-CA-C	5.15	120.42	113.72
1	A	44	THR	N-CA-C	-5.12	107.03	113.28
1	A	218	LEU	N-CA-C	-5.02	99.51	109.10

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	3065	0	3015	139	0
2	A	15	0	7	0	0
3	A	215	0	0	27	0
All	All	3295	0	3022	139	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 23.

All (139) 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:179:GLU:HB3	3:A:501:HOH:O	1.32	1.28
1:A:43:GLU:HG2	1:A:329:ARG:HD2	1.48	0.94
1:A:152:LEU:HG	3:A:621:HOH:O	1.70	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:VAL:HG22	1:A:263:TYR:CE2	2.15	0.80
1:A:258:LYS:HE2	3:A:458:HOH:O	1.84	0.76
1:A:328:GLN:O	1:A:332:ARG:HD3	1.86	0.74
1:A:117:ASP:O	1:A:121:LYS:HG2	1.86	0.74
1:A:83:THR:HB	1:A:256:TYR:OH	1.87	0.74
1:A:17:ILE:HG13	1:A:18:LEU:N	2.02	0.73
1:A:282:ARG:NE	3:A:634:HOH:O	2.21	0.73
1:A:110:GLY:O	1:A:114:VAL:HG13	1.88	0.73
1:A:27:ASP:HB3	1:A:32:LYS:HD3	1.71	0.72
1:A:304:SER:O	1:A:308:THR:HG23	1.88	0.72
1:A:330:ILE:O	1:A:334:ARG:HG3	1.92	0.69
1:A:20:LEU:HD12	1:A:380:TYR:HD1	1.58	0.68
1:A:324:THR:HG22	3:A:438:HOH:O	1.93	0.68
1:A:373:LEU:HD13	1:A:408:VAL:HG21	1.75	0.67
1:A:50:LEU:HB2	1:A:53:VAL:HG13	1.75	0.67
1:A:332:ARG:HG2	1:A:332:ARG:HH21	1.59	0.66
1:A:320:GLU:O	1:A:324:THR:HG23	1.94	0.66
1:A:90:LYS:H	1:A:90:LYS:HD2	1.61	0.66
1:A:49:VAL:HG23	1:A:53:VAL:HG22	1.78	0.65
1:A:230:ARG:HD2	3:A:451:HOH:O	1.97	0.64
1:A:252:VAL:HG13	1:A:271:THR:HB	1.79	0.63
1:A:37:ILE:CD1	1:A:41:LYS:HD2	2.28	0.63
1:A:37:ILE:HD12	1:A:41:LYS:HD2	1.81	0.63
1:A:129:ARG:HH11	1:A:156:ARG:HD3	1.64	0.63
1:A:15:ASP:HB3	1:A:18:LEU:HB2	1.83	0.61
1:A:214:GLU:HG3	3:A:570:HOH:O	2.00	0.61
1:A:255:SER:OG	1:A:257:SER:HB2	2.01	0.60
1:A:129:ARG:NH1	1:A:154:GLU:HG2	2.15	0.60
1:A:203:GLU:O	1:A:207:THR:HG23	2.02	0.59
1:A:235:GLU:O	1:A:238:GLU:HG2	2.03	0.58
1:A:346:ALA:N	1:A:348:ARG:NH2	2.51	0.58
1:A:98:LYS:HB2	3:A:537:HOH:O	2.02	0.58
1:A:348:ARG:HH12	1:A:405:VAL:HG22	1.70	0.57
1:A:25:ARG:HH21	1:A:25:ARG:HB3	1.69	0.57
1:A:324:THR:HB	1:A:327:ARG:NH1	2.20	0.57
1:A:348:ARG:HA	3:A:597:HOH:O	2.04	0.57
1:A:172:ALA:HB1	3:A:569:HOH:O	2.06	0.55
1:A:20:LEU:CD1	1:A:380:TYR:HD1	2.18	0.55
1:A:356:GLN:NE2	1:A:361:SER:HB2	2.22	0.55
1:A:374:ARG:HG3	1:A:380:TYR:CE1	2.43	0.54
1:A:370:VAL:HG21	1:A:383:ALA:HA	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:PHE:O	1:A:83:THR:CG2	2.55	0.54
1:A:382:VAL:HG21	1:A:386:ARG:CZ	2.37	0.54
1:A:274:ALA:HB1	1:A:279:THR:HG23	1.89	0.54
1:A:318:ILE:O	1:A:322:GLU:HG3	2.07	0.54
1:A:165:ASN:HB2	1:A:167:THR:HG22	1.90	0.54
1:A:73:ILE:HD12	1:A:296:SER:O	2.08	0.53
1:A:223:PHE:HD1	3:A:423:HOH:O	1.90	0.53
1:A:101:ARG:HG2	1:A:280:VAL:HG22	1.90	0.53
1:A:46:LYS:O	1:A:48:PRO:HD3	2.07	0.53
1:A:135:VAL:O	1:A:157:GLU:HA	2.09	0.52
1:A:173:LEU:HD23	1:A:177:LEU:HD23	1.90	0.52
1:A:129:ARG:HA	1:A:154:GLU:O	2.10	0.52
1:A:338:VAL:HG21	1:A:354:ILE:HG12	1.90	0.52
1:A:5:MET:HE3	3:A:627:HOH:O	2.10	0.52
1:A:193:HIS:CD2	1:A:196:THR:OG1	2.63	0.52
1:A:24:PHE:HD1	1:A:380:TYR:CD2	2.28	0.51
1:A:88:PHE:O	1:A:92:SER:HB2	2.09	0.51
1:A:212:SER:OG	1:A:247:HIS:HE1	1.94	0.51
1:A:49:VAL:HG23	1:A:53:VAL:CG2	2.40	0.50
1:A:49:VAL:HG22	1:A:54:LYS:HG2	1.93	0.50
1:A:143:HIS:HA	3:A:544:HOH:O	2.10	0.50
1:A:394:PRO:HD2	3:A:589:HOH:O	2.11	0.50
1:A:18:LEU:HA	3:A:519:HOH:O	2.11	0.50
1:A:90:LYS:H	1:A:90:LYS:CD	2.25	0.49
1:A:328:GLN:NE2	3:A:606:HOH:O	2.46	0.49
1:A:56:ALA:O	1:A:59:TYR:HB3	2.13	0.49
1:A:73:ILE:N	1:A:73:ILE:HD13	2.27	0.49
1:A:180:ALA:HB3	1:A:217:TRP:CH2	2.48	0.49
1:A:71:LEU:HD11	1:A:300:ALA:HB2	1.94	0.49
1:A:79:PHE:O	1:A:83:THR:HG22	2.13	0.49
1:A:256:TYR:HA	1:A:259:ASN:HD21	1.78	0.48
1:A:202:LEU:O	1:A:206:GLN:HG3	2.13	0.48
1:A:121:LYS:HE3	3:A:624:HOH:O	2.13	0.48
1:A:196:THR:OG1	1:A:198:ILE:HG12	2.14	0.48
1:A:346:ALA:HB3	1:A:348:ARG:CZ	2.44	0.48
1:A:396:ASN:O	1:A:400:LEU:HB2	2.13	0.48
1:A:27:ASP:OD1	1:A:29:ARG:HB2	2.14	0.47
1:A:276:ASP:OD2	1:A:279:THR:HG22	2.14	0.47
1:A:277:SER:O	1:A:280:VAL:HG12	2.14	0.47
1:A:406:ALA:HB2	3:A:608:HOH:O	2.14	0.47
1:A:149:SER:HB3	3:A:629:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:HIS:O	1:A:305:VAL:HG23	2.15	0.47
1:A:44:THR:HG22	1:A:46:LYS:HE3	1.95	0.47
1:A:230:ARG:HD3	1:A:235:GLU:HB3	1.97	0.47
1:A:250:LEU:HD23	1:A:273:VAL:HB	1.97	0.47
1:A:35:LEU:HD11	1:A:400:LEU:HD21	1.97	0.46
1:A:94:LEU:HD11	1:A:244:ALA:HB1	1.98	0.46
1:A:274:ALA:CB	1:A:279:THR:HG23	2.45	0.46
1:A:15:ASP:OD1	1:A:18:LEU:HD12	2.14	0.46
1:A:193:HIS:HD2	1:A:196:THR:OG1	1.99	0.46
1:A:141:PRO:HD2	3:A:484:HOH:O	2.15	0.45
1:A:85:GLU:HG2	1:A:90:LYS:HA	1.97	0.45
1:A:98:LYS:HA	1:A:98:LYS:HZ2	1.82	0.45
1:A:321:GLN:HG2	3:A:469:HOH:O	2.15	0.45
1:A:15:ASP:HA	1:A:16:PRO:HD2	1.61	0.45
1:A:341:LEU:HD13	1:A:350:PHE:CD2	2.52	0.45
1:A:40:TYR:CD1	1:A:326:MET:HE3	2.53	0.44
1:A:322:GLU:O	1:A:326:MET:HG3	2.17	0.44
1:A:348:ARG:HB2	1:A:350:PHE:CE2	2.53	0.44
1:A:374:ARG:HA	1:A:379:VAL:O	2.17	0.44
1:A:343:GLU:HB2	3:A:456:HOH:O	2.18	0.44
1:A:144:LYS:HE3	1:A:144:LYS:HB2	1.83	0.44
1:A:227:GLY:O	1:A:327:ARG:HD3	2.18	0.44
1:A:282:ARG:CZ	3:A:634:HOH:O	2.62	0.44
1:A:73:ILE:HD11	3:A:491:HOH:O	2.17	0.44
1:A:53:VAL:HB	1:A:305:VAL:HG11	2.00	0.43
1:A:187:LEU:C	1:A:187:LEU:HD12	2.43	0.43
1:A:350:PHE:O	1:A:353:ILE:HB	2.19	0.43
1:A:27:ASP:HB2	3:A:506:HOH:O	2.18	0.43
1:A:277:SER:HA	1:A:280:VAL:HG12	2.01	0.43
1:A:79:PHE:O	1:A:83:THR:HG23	2.19	0.43
1:A:84:GLN:HE22	1:A:98:LYS:NZ	2.17	0.43
1:A:259:ASN:HD22	1:A:260:PHE:N	2.17	0.42
1:A:198:ILE:HG12	1:A:198:ILE:H	1.69	0.42
1:A:100:ALA:HA	1:A:272:LEU:O	2.19	0.42
1:A:72:GLY:HA2	3:A:497:HOH:O	2.20	0.42
1:A:113:ARG:NH1	3:A:548:HOH:O	2.52	0.42
1:A:29:ARG:O	1:A:32:LYS:HG2	2.20	0.41
1:A:20:LEU:HD12	1:A:380:TYR:CD1	2.46	0.41
1:A:37:ILE:HG13	1:A:388:ASN:ND2	2.35	0.41
1:A:49:VAL:CG2	1:A:54:LYS:HG2	2.50	0.41
1:A:39:VAL:HG22	1:A:263:TYR:CD2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:SER:HA	1:A:217:TRP:CD1	2.55	0.41
1:A:316:ARG:O	1:A:320:GLU:HG3	2.20	0.41
1:A:58:GLN:HE21	1:A:62:GLU:HG3	1.86	0.41
1:A:136:SER:OG	1:A:193:HIS:HE1	2.03	0.41
1:A:138:PRO:HG3	1:A:362:PHE:HE1	1.85	0.41
1:A:248:LYS:HD3	1:A:248:LYS:HA	1.64	0.41
1:A:225:PHE:HE2	1:A:359:MET:HE3	1.84	0.41
1:A:382:VAL:HG21	1:A:386:ARG:NH1	2.36	0.41
1:A:106:PRO:HA	1:A:266:ARG:O	2.20	0.41
1:A:338:VAL:CG2	1:A:353:ILE:HB	2.51	0.41
1:A:372:ARG:O	1:A:376:GLU:HB3	2.20	0.41
1:A:341:LEU:O	1:A:348:ARG:NH2	2.53	0.40
1:A:335:GLN:HE21	1:A:354:ILE:HD12	1.86	0.40

There are no symmetry-related clashes.

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	394/396 (100%)	365 (93%)	23 (6%)	6 (2%)	8 4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	ALA
1	A	345	GLY
1	A	349	ASP
1	A	16	PRO
1	A	142	ASN
1	A	92	SER

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	319/319 (100%)	262 (82%)	57 (18%)	2 1

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	9	ILE
1	A	10	THR
1	A	17	ILE
1	A	20	LEU
1	A	23	LEU
1	A	25	ARG
1	A	29	ARG
1	A	39	VAL
1	A	43	GLU
1	A	49	VAL
1	A	53	VAL
1	A	76	ILE
1	A	83	THR
1	A	87	LEU
1	A	90	LYS
1	A	98	LYS
1	A	99	ARG
1	A	114	VAL
1	A	119	LEU
1	A	125	VAL
1	A	129	ARG
1	A	154	GLU
1	A	164	GLU
1	A	168	LEU
1	A	177	LEU
1	A	181	GLN
1	A	187	LEU
1	A	198	ILE
1	A	208	LEU

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Mol	Chain	Res	Type
1	A	210	GLN
1	A	218	LEU
1	A	220	LEU
1	A	230	ARG
1	A	252	VAL
1	A	254	SER
1	A	259	ASN
1	A	267	VAL
1	A	277	SER
1	A	304	SER
1	A	308	THR
1	A	315	LEU
1	A	321	GLN
1	A	324	THR
1	A	332	ARG
1	A	335	GLN
1	A	336	LEU
1	A	339	ASN
1	A	348	ARG
1	A	365	LEU
1	A	367	LYS
1	A	370	VAL
1	A	371	LEU
1	A	375	GLU
1	A	387	VAL
1	A	400	LEU
1	A	409	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	84	GLN
1	A	96	ASN
1	A	122	ASN
1	A	142	ASN
1	A	148	ASN
1	A	166	HIS
1	A	193	HIS
1	A	206	GLN
1	A	226	GLN
1	A	247	HIS

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Mol	Chain	Res	Type
1	A	259	ASN
1	A	328	GLN
1	A	335	GLN
1	A	356	GLN
1	A	357	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	413	1	15,15,16	2.39	5 (33%)	21,22,23	2.09	7 (33%)

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	413	1	-	3/6/6/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	413	PLP	C3-C2	-5.35	1.35	1.41
2	A	413	PLP	C5-C4	-4.23	1.35	1.40
2	A	413	PLP	C2A-C2	3.93	1.56	1.50
2	A	413	PLP	O4P-C5A	-2.61	1.35	1.44
2	A	413	PLP	P-O1P	-2.53	1.42	1.50

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	413	PLP	O4P-C5A-C5	5.99	120.59	109.36
2	A	413	PLP	O2P-P-O4P	-4.42	95.13	106.67
2	A	413	PLP	O4P-P-O1P	2.72	113.79	106.44
2	A	413	PLP	O3P-P-O1P	2.55	120.79	110.83
2	A	413	PLP	C5-C6-N1	-2.23	120.21	123.83
2	A	413	PLP	O3P-P-O4P	-2.21	100.91	106.67
2	A	413	PLP	C6-C5-C4	2.03	119.76	118.10

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	413	PLP	C6-C5-C5A-O4P
2	A	413	PLP	C5A-O4P-P-O2P
2	A	413	PLP	C5A-O4P-P-O3P

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](#)

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 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.