



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 05:32 PM UTC

PDB ID : 1I1L / pdb_00001i1l
Title : CRYSTAL STRUCTURE OF ESCHERICHIA COLI BRANCHED-CHAIN AMINO ACID AMINOTRANSFERASE.
Authors : Okada, K.; Hirotsu, K.; Hayashi, H.; Kagamiyama, H.
Deposited on : 2001-02-02
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

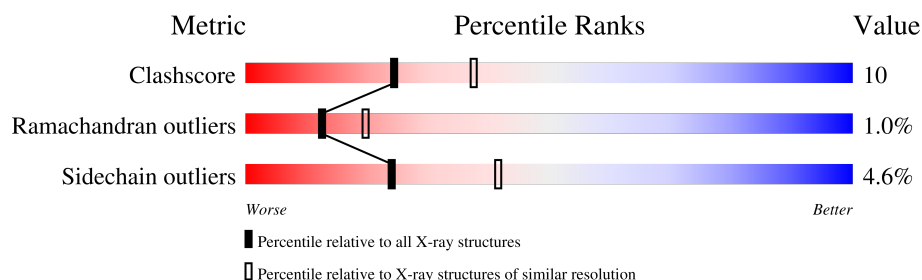
1 Overall quality at a glance

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	309	
1	B	309	
1	C	309	

2 Entry composition [i](#)

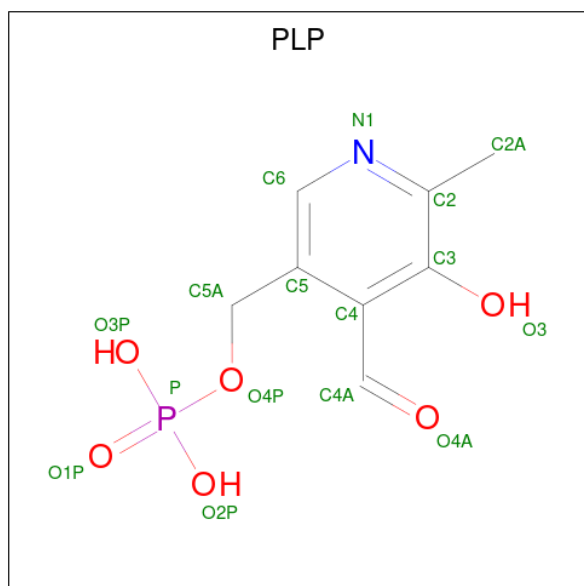
There are 4 unique types of molecules in this entry. The entry contains 7465 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BRANCHED-CHAIN AMINO ACID AMINOTRANSFERASE.

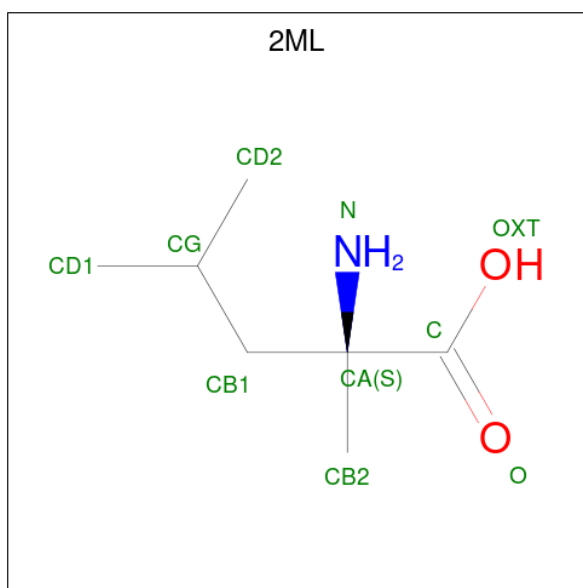
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2362	1499	410	443	10			
1	B	304	Total	C	N	O	S	0	0	0
			2362	1499	410	443	10			
1	C	304	Total	C	N	O	S	0	0	0
			2362	1499	410	443	10			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: 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		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is 2-METHYLLEUCINE (CCD ID: 2ML) (formula: $C_7H_{15}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	7	1	2		
3	B	1	Total	C	N	O	0	0
			10	7	1	2		
3	C	1	Total	C	N	O	0	0
			10	7	1	2		

- Molecule 4 is water.

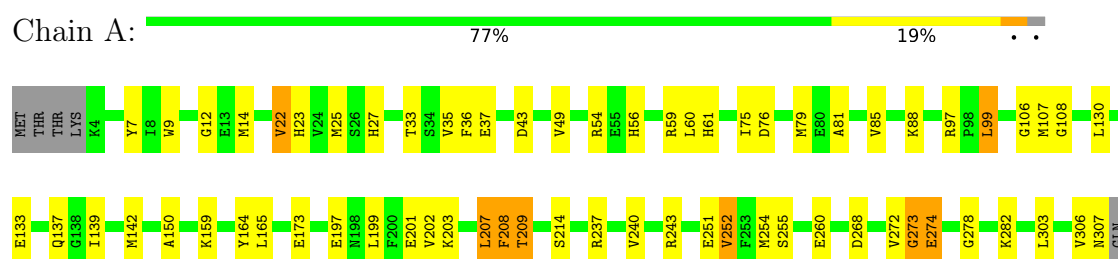
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	97	Total	O	0	0
			97	97		
4	B	102	Total	O	0	0
			102	102		
4	C	105	Total	O	0	0
			105	105		

3 Residue-property plots [i](#)

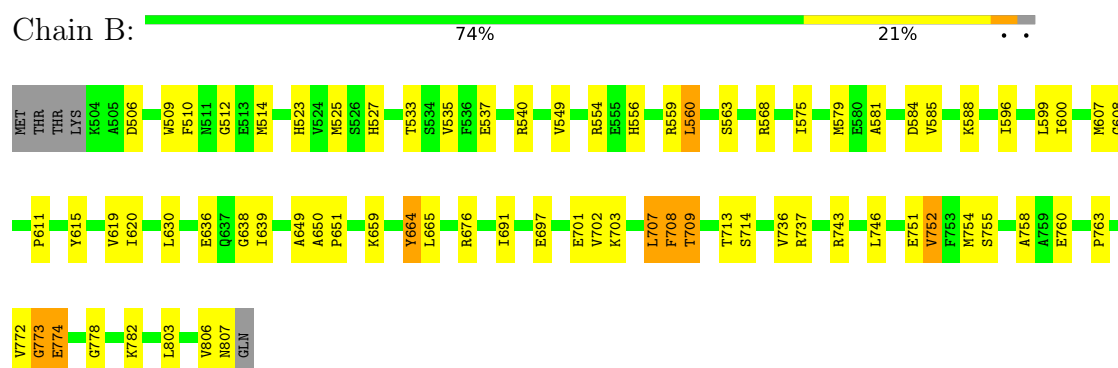
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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.

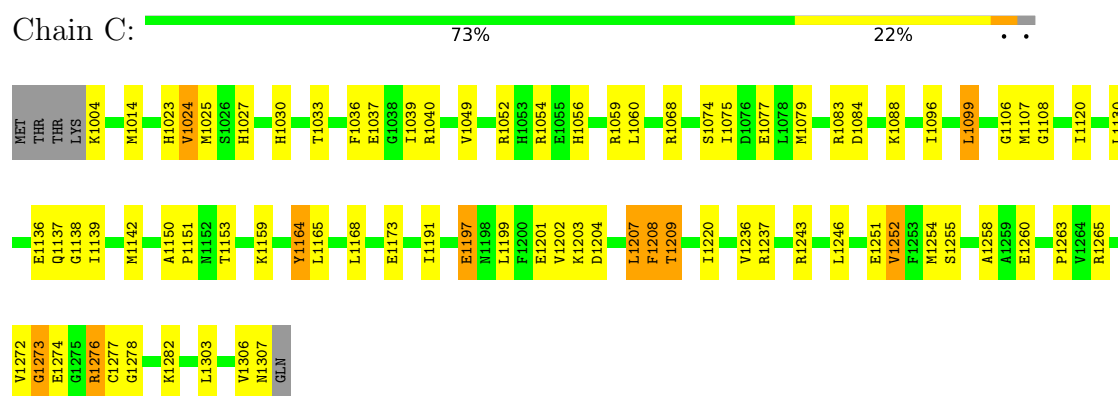
• Molecule 1: BRANCHED-CHAIN AMINO ACID AMINOTRANSFERASE



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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	156.09 Å 100.77 Å 141.56 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (40.00-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.182 , 0.224	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7465	wwPDB-VP
Average B, all atoms (Å ²)	21.0	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, 2ML

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	0.44	0/2418	0.91	8/3279 (0.2%)
1	B	0.44	0/2418	0.90	9/3279 (0.3%)
1	C	0.45	0/2418	0.91	6/3279 (0.2%)
All	All	0.44	0/7254	0.91	23/9837 (0.2%)

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	B	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	LEU	N-CA-C	-7.54	103.08	111.82
1	A	164	TYR	N-CA-C	7.37	121.95	113.19
1	B	664	TYR	N-CA-C	6.95	121.46	113.19
1	C	1164	TYR	N-CA-C	6.95	121.46	113.19
1	C	1278	GLY	CA-C-N	6.94	126.19	118.97
1	C	1278	GLY	C-N-CA	6.94	126.19	118.97
1	A	278	GLY	CA-C-N	6.89	126.40	119.24
1	A	278	GLY	C-N-CA	6.89	126.40	119.24
1	C	1165	LEU	N-CA-C	-6.84	103.88	111.82
1	B	665	LEU	N-CA-C	-6.53	104.25	111.82
1	A	150	ALA	CA-C-N	6.33	127.75	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	ALA	C-N-CA	6.33	127.75	119.84
1	B	650	ALA	CA-C-N	5.95	127.27	119.84
1	B	650	ALA	C-N-CA	5.95	127.27	119.84
1	C	1150	ALA	CA-C-N	5.78	127.06	119.84
1	C	1150	ALA	C-N-CA	5.78	127.06	119.84
1	B	506	ASP	N-CA-C	5.71	118.36	111.40
1	B	714	SER	N-CA-C	-5.53	106.42	112.72
1	B	778	GLY	CA-C-N	5.42	126.61	119.84
1	B	778	GLY	C-N-CA	5.42	126.61	119.84
1	B	649	ALA	N-CA-C	5.28	117.75	110.35
1	A	214	SER	N-CA-C	-5.19	106.81	112.72
1	A	43	ASP	N-CA-C	-5.02	101.56	109.25

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	664	TYR	Sidechain
1	C	1164	TYR	Sidechain

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	2362	0	2302	42	0
1	B	2362	0	2302	54	0
1	C	2362	0	2302	46	0
2	A	15	0	7	0	0
2	B	15	0	7	1	0
2	C	15	0	7	0	0
3	A	10	0	13	0	0
3	B	10	0	13	0	0
3	C	10	0	13	0	0
4	A	97	0	0	5	0
4	B	102	0	0	7	0
4	C	105	0	0	4	0
All	All	7465	0	6966	140	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 10.

All (140) 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:B:549:VAL:HG21	1:B:803:LEU:HB3	1.49	0.94
1:C:1049:VAL:HG21	1:C:1303:LEU:HB3	1.56	0.88
1:B:575:ILE:HG22	1:B:579:MET:HE2	1.64	0.80
1:A:202:VAL:HB	1:A:251:GLU:HB2	1.66	0.77
1:B:579:MET:SD	4:B:225:HOH:O	2.44	0.75
1:B:540:ARG:NH1	1:B:758:ALA:HA	2.03	0.74
1:A:23:HIS:HD2	1:A:25:MET:H	1.37	0.71
1:B:752:VAL:HG12	1:B:772:VAL:HG11	1.70	0.71
1:A:49:VAL:HG21	1:A:303:LEU:HB3	1.72	0.71
1:B:702:VAL:HB	1:B:751:GLU:HB2	1.72	0.71
1:C:1023:HIS:HD2	1:C:1025:MET:H	1.39	0.68
1:A:75:ILE:HG22	1:A:79:MET:HE2	1.77	0.67
1:C:1202:VAL:HB	1:C:1251:GLU:HB2	1.79	0.65
1:B:630:LEU:HD13	1:B:639:ILE:HD13	1.79	0.64
1:A:36:PHE:HB3	1:A:99:LEU:HD23	1.80	0.63
1:C:1142:MET:HE1	1:C:1173:GLU:HG2	1.79	0.63
1:C:1059:ARG:NH2	1:C:1159:LYS:HE3	2.14	0.61
1:C:1054:ARG:NH2	1:C:1079:MET:HE1	2.16	0.61
1:A:106:GLY:HA3	4:A:485:HOH:O	1.99	0.61
1:A:252:VAL:HG12	1:A:272:VAL:HG11	1.83	0.60
1:B:554:ARG:NH2	1:B:579:MET:HE1	2.17	0.60
1:B:708:PHE:N	4:B:45:HOH:O	2.35	0.60
1:B:523:HIS:HD2	1:B:525:MET:HB2	1.66	0.59
1:C:1207:LEU:O	1:C:1208:PHE:HB2	2.03	0.59
1:B:559:ARG:NH2	1:B:659:LYS:HE3	2.17	0.59
1:B:540:ARG:HH12	1:B:758:ALA:HA	1.67	0.59
1:C:1040:ARG:NH1	1:C:1258:ALA:HA	2.17	0.59
1:C:1027:HIS:HD2	1:C:1033:THR:OG1	1.86	0.58
1:A:9:TRP:CZ2	1:A:12:GLY:HA2	2.38	0.58
1:C:1273:GLY:O	1:C:1274:GLU:HB2	2.04	0.58
1:B:659:LYS:HZ1	2:B:913:PLP:C4A	2.17	0.58
1:A:27:HIS:HD2	1:A:33:THR:OG1	1.87	0.57
1:B:701:GLU:O	1:B:707:LEU:O	2.23	0.57
1:B:773:GLY:O	1:B:774:GLU:HB2	2.03	0.56
1:C:1277:CYS:SG	1:C:1282:LYS:HG2	2.45	0.56
1:A:23:HIS:CD2	1:A:25:MET:H	2.22	0.56
1:C:1106:GLY:HA3	4:C:282:HOH:O	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:707:LEU:O	1:B:708:PHE:HB2	2.05	0.55
1:A:54:ARG:NH2	1:A:79:MET:HE1	2.21	0.55
1:A:207:LEU:O	1:A:208:PHE:HB2	2.05	0.55
1:C:1023:HIS:CD2	1:C:1025:MET:H	2.23	0.55
1:A:59:ARG:NH2	1:A:159:LYS:HE3	2.22	0.54
1:C:1130:LEU:HD13	1:C:1139:ILE:HD13	1.88	0.54
1:C:1252:VAL:HG12	1:C:1272:VAL:HG11	1.90	0.54
1:C:1201:GLU:O	1:C:1207:LEU:O	2.24	0.54
1:A:306:VAL:O	1:A:307:ASN:HB2	2.07	0.54
1:A:23:HIS:HD2	1:A:25:MET:N	2.05	0.54
1:B:527:HIS:HD2	1:B:533:THR:OG1	1.90	0.54
1:B:527:HIS:HE1	1:B:608:GLY:O	1.91	0.54
1:B:773:GLY:HA2	4:B:297:HOH:O	2.08	0.54
1:B:584:ASP:O	1:B:588:LYS:HB2	2.08	0.53
1:A:142:MET:HE1	1:A:173:GLU:HG2	1.90	0.53
1:A:197:GLU:HG3	1:A:254:MET:HB3	1.91	0.53
1:B:535:VAL:HG12	1:B:560:LEU:HD23	1.90	0.53
1:A:273:GLY:O	1:A:274:GLU:HB2	2.09	0.53
1:C:1068:ARG:HD2	4:C:244:HOH:O	2.09	0.52
1:A:56:HIS:HE1	4:A:428:HOH:O	1.92	0.52
1:C:1014:MET:HE2	1:C:1088:LYS:HG2	1.91	0.51
1:A:208:PHE:HA	1:A:237:ARG:O	2.10	0.51
1:A:59:ARG:HG3	4:A:466:HOH:O	2.10	0.51
1:B:510:PHE:HA	1:B:620:ILE:HD13	1.92	0.51
1:A:22:VAL:HG22	4:A:441:HOH:O	2.11	0.50
1:C:1207:LEU:HD12	1:C:1236:VAL:HG22	1.94	0.50
1:C:1208:PHE:HA	1:C:1237:ARG:O	2.11	0.50
1:A:201:GLU:O	1:A:207:LEU:O	2.28	0.50
1:B:611:PRO:HB2	1:B:615:TYR:CE2	2.46	0.50
1:B:554:ARG:CZ	1:B:579:MET:HE1	2.41	0.50
1:B:510:PHE:CG	1:B:620:ILE:HD11	2.47	0.50
1:C:1136:GLU:HG2	1:C:1137:GLN:H	1.77	0.50
1:B:581:ALA:O	1:B:585:VAL:HG23	2.11	0.50
1:A:107:MET:HA	1:A:107:MET:HE2	1.94	0.50
1:C:1037:GLU:OE2	1:C:1056:HIS:HD2	1.95	0.50
1:C:1306:VAL:O	1:C:1307:ASN:HB2	2.12	0.49
1:A:133:GLU:O	1:A:137:GLN:HG3	2.13	0.49
1:C:1024:VAL:O	1:C:1030:HIS:HE1	1.96	0.49
1:A:27:HIS:HE1	1:A:108:GLY:O	1.94	0.49
1:B:707:LEU:HD12	1:B:736:VAL:HG22	1.93	0.49
1:B:697:GLU:HG3	1:B:754:MET:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:ARG:CZ	1:A:79:MET:HE1	2.44	0.48
1:B:600:ILE:HG12	1:B:619:VAL:HG22	1.94	0.48
1:C:1138:GLY:O	1:C:1265:ARG:NH1	2.46	0.48
1:C:1274:GLU:HB3	1:C:1276:ARG:HE	1.79	0.47
1:B:523:HIS:CD2	1:B:525:MET:H	2.32	0.47
1:C:1255:SER:HA	1:C:1260:GLU:O	2.14	0.47
1:C:1027:HIS:HE1	1:C:1108:GLY:O	1.98	0.47
1:C:1207:LEU:O	1:C:1208:PHE:CB	2.62	0.47
1:A:130:LEU:HD13	1:A:139:ILE:HD13	1.97	0.47
1:A:207:LEU:O	1:A:208:PHE:CB	2.63	0.47
1:A:81:ALA:O	1:A:85:VAL:HG23	2.14	0.46
1:B:707:LEU:O	1:B:708:PHE:CB	2.63	0.46
1:C:1059:ARG:HH21	1:C:1159:LYS:HE3	1.80	0.46
1:B:509:TRP:CZ2	1:B:512:GLY:HA2	2.51	0.46
1:A:37:GLU:OE2	1:A:56:HIS:HD2	1.98	0.46
1:B:709:THR:HG22	4:B:6:HOH:O	2.15	0.46
1:B:638:GLY:HA2	1:B:763:PRO:HG2	1.98	0.45
1:B:514:MET:HE2	1:B:588:LYS:HG2	1.97	0.45
1:B:676:ARG:NH2	4:B:184:HOH:O	2.50	0.45
1:C:1153:THR:HG21	4:C:1:HOH:O	2.16	0.45
1:C:1197:GLU:HG3	1:C:1254:MET:HB3	1.98	0.45
1:C:1207:LEU:HD23	1:C:1207:LEU:HA	1.80	0.45
1:A:268:ASP:O	1:B:568:ARG:NH2	2.49	0.45
1:B:535:VAL:HG12	1:B:560:LEU:CD2	2.46	0.45
1:C:1083:ARG:HD2	4:C:165:HOH:O	2.17	0.45
1:B:708:PHE:HA	1:B:737:ARG:O	2.17	0.45
1:C:1191:ILE:HD11	1:C:1246:LEU:HG	1.99	0.45
1:A:14:MET:HE2	1:A:88:LYS:HG2	1.98	0.45
1:B:556:HIS:HE1	4:B:52:HOH:O	1.99	0.45
1:B:537:GLU:OE2	1:B:556:HIS:HD2	2.00	0.44
1:C:1023:HIS:HD2	1:C:1025:MET:N	2.11	0.44
1:B:596:ILE:HD12	1:B:596:ILE:N	2.33	0.44
1:B:607:MET:HA	1:B:607:MET:HE2	1.99	0.44
1:C:1138:GLY:HA2	1:C:1263:PRO:HG2	1.99	0.43
1:B:575:ILE:O	1:B:579:MET:HG3	2.18	0.43
1:C:1197:GLU:HB2	1:C:1255:SER:O	2.18	0.43
1:A:56:HIS:CE1	4:A:428:HOH:O	2.70	0.43
1:B:691:ILE:HD11	1:B:746:LEU:HG	2.00	0.43
1:B:535:VAL:HG11	1:B:563:SER:HB2	2.00	0.43
1:A:203:LYS:HB2	1:A:208:PHE:HE1	1.83	0.43
1:A:255:SER:HA	1:A:260:GLU:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:755:SER:HA	1:B:760:GLU:O	2.19	0.42
1:A:199:LEU:O	1:A:209:THR:HG23	2.19	0.42
1:B:611:PRO:HB2	1:B:615:TYR:CD2	2.54	0.42
1:A:59:ARG:NH2	1:A:159:LYS:CE	2.82	0.42
1:B:703:LYS:HB2	1:B:708:PHE:CE1	2.54	0.42
1:A:203:LYS:HB2	1:A:208:PHE:CE1	2.54	0.41
1:C:1203:LYS:HB3	1:C:1204:ASP:H	1.72	0.41
1:A:7:TYR:CB	1:A:14:MET:HE3	2.50	0.41
1:A:240:VAL:HG12	1:B:713:THR:HG21	2.01	0.41
1:B:559:ARG:NH2	1:B:659:LYS:CE	2.83	0.41
1:C:1039:ILE:HB	1:C:1096:ILE:HB	2.03	0.41
1:A:207:LEU:HD23	1:A:207:LEU:HA	1.94	0.41
1:C:1052:ARG:HD3	1:C:1220:ILE:CG2	2.51	0.41
1:C:1074:SER:OG	1:C:1077:GLU:HG3	2.20	0.41
1:C:1199:LEU:O	1:C:1209:THR:HG22	2.21	0.41
1:B:523:HIS:CD2	1:B:525:MET:HB2	2.51	0.41
1:B:737:ARG:N	4:B:45:HOH:O	2.48	0.40
1:C:1107:MET:HA	1:C:1107:MET:HE2	2.03	0.40
1:B:806:VAL:O	1:B:807:ASN:HB2	2.22	0.40
1:C:1036:PHE:HB3	1:C:1099:LEU:HD23	2.04	0.40
1:C:1075:ILE:O	1:C:1079:MET:HG3	2.21	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	302/309 (98%)	285 (94%)	14 (5%)	3 (1%)	12	20
1	B	302/309 (98%)	282 (93%)	17 (6%)	3 (1%)	12	20
1	C	302/309 (98%)	284 (94%)	15 (5%)	3 (1%)	12	20
All	All	906/927 (98%)	851 (94%)	46 (5%)	9 (1%)	12	20

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	207	LEU
1	B	707	LEU
1	C	1207	LEU
1	A	273	GLY
1	B	708	PHE
1	B	773	GLY
1	C	1273	GLY
1	A	208	PHE
1	C	1208	PHE

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	246/251 (98%)	234 (95%)	12 (5%)	22	39
1	B	246/251 (98%)	237 (96%)	9 (4%)	30	51
1	C	246/251 (98%)	233 (95%)	13 (5%)	20	36
All	All	738/753 (98%)	704 (95%)	34 (5%)	24	41

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	35	VAL
1	A	60	LEU
1	A	61	HIS
1	A	76	ASP
1	A	97	ARG
1	A	99	LEU
1	A	209	THR
1	A	243	ARG
1	A	252	VAL
1	A	274	GLU
1	A	282	LYS

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Mol	Chain	Res	Type
1	B	560	LEU
1	B	599	LEU
1	B	636	GLU
1	B	651	PRO
1	B	709	THR
1	B	743	ARG
1	B	752	VAL
1	B	774	GLU
1	B	782	LYS
1	C	1004	LYS
1	C	1024	VAL
1	C	1060	LEU
1	C	1084	ASP
1	C	1099	LEU
1	C	1120	ILE
1	C	1151	PRO
1	C	1168	LEU
1	C	1197	GLU
1	C	1209	THR
1	C	1243	ARG
1	C	1252	VAL
1	C	1276	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	HIS
1	A	27	HIS
1	A	30	HIS
1	A	45	HIS
1	A	56	HIS
1	A	58	GLN
1	A	61	HIS
1	B	523	HIS
1	B	527	HIS
1	B	545	HIS
1	B	556	HIS
1	B	647	ASN
1	B	771	GLN
1	C	1023	HIS
1	C	1027	HIS
1	C	1030	HIS

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Mol	Chain	Res	Type
1	C	1045	HIS
1	C	1056	HIS
1	C	1058	GLN
1	C	1147	ASN
1	C	1177	HIS
1	C	1239	GLN

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 ⓘ

6 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	$\# Z > 2$
3	2ML	A	414	2	6,9,9	0.72	0	8,13,13	1.25	1 (12%)
2	PLP	B	913	3	15,15,16	1.89	4 (26%)	21,22,23	1.41	5 (23%)
3	2ML	C	1414	2	6,9,9	0.77	0	8,13,13	1.28	1 (12%)
3	2ML	B	914	2	6,9,9	0.71	0	8,13,13	1.24	1 (12%)
2	PLP	A	413	3	15,15,16	1.43	2 (13%)	21,22,23	1.15	1 (4%)
2	PLP	C	1413	3	15,15,16	1.62	3 (20%)	21,22,23	1.16	3 (14%)

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	2ML	A	414	2	-	2/11/11/11	-
2	PLP	B	913	3	-	1/6/6/8	0/1/1/1
3	2ML	C	1414	2	-	3/11/11/11	-
3	2ML	B	914	2	-	4/11/11/11	-
2	PLP	A	413	3	-	0/6/6/8	0/1/1/1
2	PLP	C	1413	3	-	1/6/6/8	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	913	PLP	C3-C2	-4.16	1.36	1.41
2	B	913	PLP	C5-C4	-4.12	1.35	1.40
2	C	1413	PLP	C2A-C2	3.46	1.55	1.50
2	A	413	PLP	C3-C2	-2.91	1.38	1.41
2	A	413	PLP	C5-C4	-2.43	1.37	1.40
2	B	913	PLP	C2A-C2	2.32	1.54	1.50
2	C	1413	PLP	P-O3P	-2.25	1.46	1.54
2	B	913	PLP	P-O3P	-2.25	1.46	1.54
2	C	1413	PLP	C4A-C4	2.11	1.55	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	913	PLP	O2P-P-O4P	-3.38	97.86	106.67
3	B	914	2ML	O-C-CA	-2.94	117.51	122.78
3	C	1414	2ML	O-C-CA	-2.80	117.75	122.78
3	A	414	2ML	O-C-CA	-2.68	117.97	122.78
2	C	1413	PLP	C6-C5-C4	2.54	120.18	118.10
2	A	413	PLP	O3P-P-O1P	2.54	120.73	110.83
2	B	913	PLP	O3P-P-O1P	2.44	120.35	110.83
2	B	913	PLP	C6-C5-C4	2.24	119.93	118.10
2	C	1413	PLP	O3P-P-O1P	2.23	119.53	110.83
2	B	913	PLP	C4A-C4-C5	-2.12	118.75	120.94
2	C	1413	PLP	O2P-P-O4P	-2.12	101.14	106.67
2	B	913	PLP	C5-C6-N1	-2.00	120.58	123.83

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	414	2ML	O-C-CA-N
3	C	1414	2ML	O-C-CA-N
2	B	913	PLP	C4-C5-C5A-O4P
3	A	414	2ML	C-CA-CB1-CG
2	C	1413	PLP	C5A-O4P-P-O1P
3	B	914	2ML	OXT-C-CA-CB1
3	B	914	2ML	O-C-CA-N
3	B	914	2ML	OXT-C-CA-CB2
3	C	1414	2ML	C-CA-CB1-CG
3	B	914	2ML	O-C-CA-CB1
3	C	1414	2ML	OXT-C-CA-CB1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	913	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 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.