



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:32 PM UTC

PDB ID : 2VGZ / pdb_00002vgz
Title : CRYSTAL STRUCTURE OF HUMAN KYNURENINE AMINOTRANSFERASE II
Authors : Rossi, F.; Garavaglia, S.; Montalbano, V.; Walsh, M.A.; Rizzi, M.
Deposited on : 2007-11-16
Resolution : 2.30 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

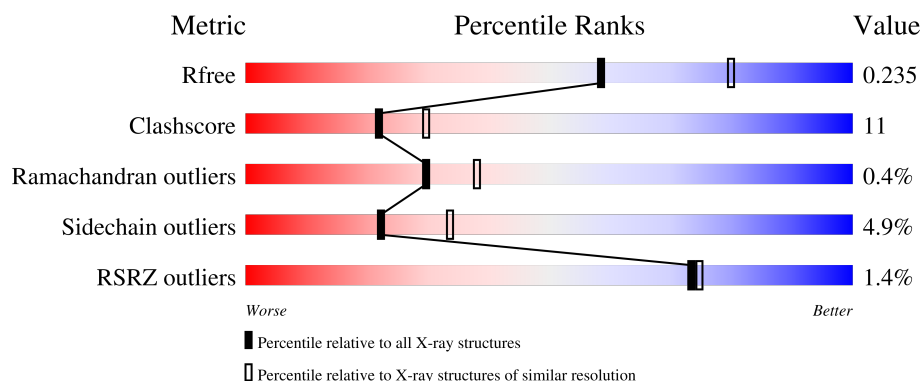
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

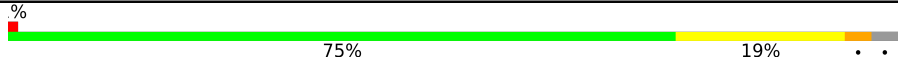
The reported resolution of this entry is 2.30 Å.

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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	427	 75% 19% . .
1	B	427	 77% 18% . .

2 Entry composition [i](#)

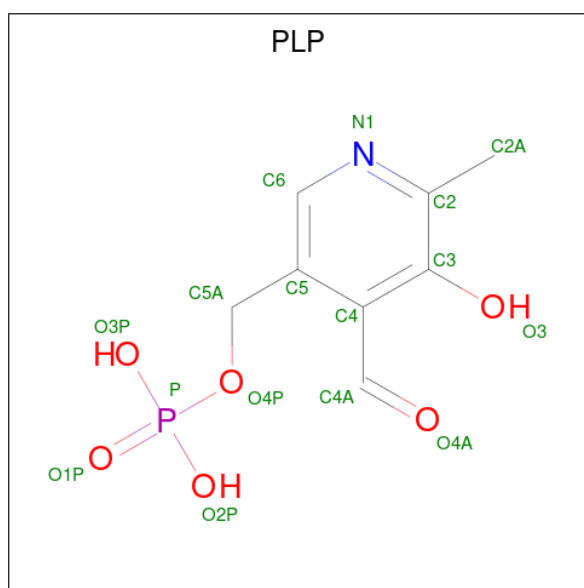
There are 4 unique types of molecules in this entry. The entry contains 6927 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 KYNURENINE/ α -AMINOADIPATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	0	0
			3246	2087	541	601	17			
1	B	414	Total	C	N	O	S	0	0	0
			3247	2087	541	602	17			

- 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		

- Molecule 3 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total I 1 1	0	0

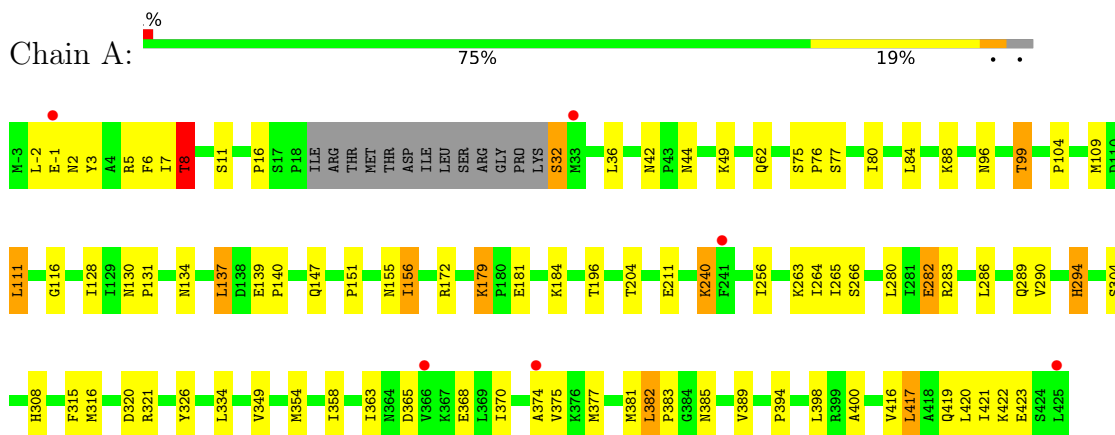
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	199	Total O 199 199	0	0
4	B	204	Total O 204 204	0	0

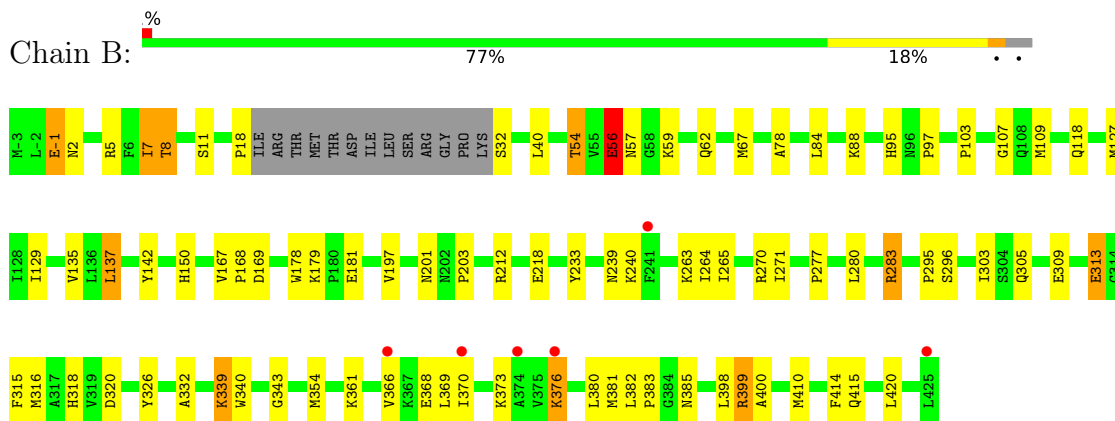
3 Residue-property plots [i](#)

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 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: KYNURENINE/ALPHA-AMINOADIPATE AMINOTRANSFERASE



• Molecule 1: KYNURENINE/ALPHA-AMINOADIPATE AMINOTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	99.63Å 154.96Å 61.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.59 – 2.30 29.59 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.59-2.30) 98.8 (29.59-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.21 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.182 , 0.243 (Not available) , 0.235	Depositor DCC
R_{free} test set	2780 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6927	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: IOD, 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	3/3325 (0.1%)	1.09	10/4511 (0.2%)
1	B	1.28	4/3327 (0.1%)	1.08	8/4516 (0.2%)
All	All	1.17	7/6652 (0.1%)	1.08	18/9027 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	-1	GLU	C-N	42.95	1.99	1.33
1	A	80	ILE	CA-C	7.36	1.59	1.52
1	B	271	ILE	CA-CB	6.32	1.62	1.54
1	A	96	ASN	C-O	-5.71	1.21	1.23
1	B	103	PRO	CA-C	5.25	1.57	1.52
1	A	290	VAL	CA-CB	5.16	1.60	1.54
1	B	167	VAL	CA-CB	5.00	1.58	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	-1	GLU	O-C-N	-10.61	108.48	122.59
1	A	-2	LEU	N-CA-C	8.57	122.23	111.69
1	A	179	LYS	CA-C-N	7.38	126.92	119.24
1	A	179	LYS	C-N-CA	7.38	126.92	119.24
1	A	130	ASN	N-CA-C	-6.41	101.23	110.40
1	B	-1	GLU	CA-C-N	-6.05	109.83	121.94
1	B	-1	GLU	C-N-CA	-6.05	109.83	121.94
1	B	40	LEU	CA-C-N	-6.01	114.44	120.21
1	B	40	LEU	C-N-CA	-6.01	114.44	120.21
1	A	128	ILE	N-CA-C	5.51	116.58	111.45
1	B	179	LYS	N-CA-C	-5.45	103.31	110.39
1	B	150	HIS	CA-C-N	-5.38	114.13	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	150	HIS	C-N-CA	-5.38	114.13	119.56
1	A	321	ARG	N-CA-C	5.27	117.03	111.28
1	A	116	GLY	N-CA-C	-5.09	106.32	112.48
1	A	294	HIS	CA-C-N	-5.04	115.37	120.31
1	A	294	HIS	C-N-CA	-5.04	115.37	120.31
1	A	8	THR	N-CA-CB	-5.01	103.47	110.29

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	3246	0	3252	74	0
1	B	3247	0	3252	71	0
2	A	15	0	6	2	0
2	B	15	0	6	2	0
3	A	1	0	0	0	0
4	A	199	0	0	13	0
4	B	204	0	0	16	0
All	All	6927	0	6516	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 11.

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:A:62:GLN:NE2	4:A:2039:HOH:O	1.66	1.23
1:B:-1:GLU:C	1:B:2:ASN:N	1.98	1.20
1:A:240:LYS:HD3	1:A:240:LYS:H	1.10	1.14
1:B:8:THR:HG21	4:B:2078:HOH:O	1.47	1.12
1:A:-1:GLU:C	1:A:2:ASN:HA	1.80	1.05
1:A:315:PHE:CD2	1:A:316:MET:HE2	1.99	0.97
1:B:399:ARG:HG2	1:B:399:ARG:HH11	1.31	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:369:LEU:HD11	1:B:420:LEU:HB3	1.44	0.95
1:A:8:THR:HG21	4:A:2076:HOH:O	1.69	0.91
1:B:56:GLU:OE1	1:B:57:ASN:N	2.02	0.90
1:A:88:LYS:HG2	1:A:111:LEU:HD13	1.52	0.89
1:A:99:THR:HG21	4:A:2071:HOH:O	1.75	0.87
1:A:8:THR:HG22	1:A:11:SER:H	1.41	0.85
1:A:-1:GLU:C	1:A:2:ASN:CA	2.51	0.83
1:A:-1:GLU:O	1:A:2:ASN:HA	1.81	0.81
1:B:56:GLU:CD	1:B:57:ASN:H	1.90	0.80
1:A:36:LEU:HD12	1:B:381:MET:HE3	1.64	0.80
1:A:315:PHE:HD2	1:A:316:MET:HE2	1.47	0.79
1:A:240:LYS:H	1:A:240:LYS:CD	1.93	0.77
1:A:240:LYS:HD3	1:A:240:LYS:N	1.94	0.77
1:A:99:THR:HG22	1:A:109:MET:HB2	1.68	0.75
1:A:-1:GLU:C	1:A:2:ASN:N	2.46	0.74
1:A:36:LEU:HD12	1:B:381:MET:CE	2.17	0.74
1:B:178:TRP:CZ2	4:B:2105:HOH:O	2.41	0.73
1:B:265:ILE:HG21	1:B:303:ILE:HD12	1.72	0.71
1:B:339:LYS:HD2	4:B:2175:HOH:O	1.92	0.69
1:B:67:MET:SD	1:B:305:GLN:NE2	2.66	0.69
1:B:315:PHE:CD2	1:B:316:MET:HE2	2.27	0.69
1:A:294:HIS:CE1	1:B:270:ARG:HD3	2.28	0.68
1:B:399:ARG:HG2	1:B:399:ARG:NH1	2.08	0.68
1:A:179:LYS:HD3	4:A:2107:HOH:O	1.92	0.67
1:A:349:VAL:HG12	4:A:2179:HOH:O	1.96	0.66
1:A:315:PHE:CD2	1:A:316:MET:CE	2.78	0.66
1:A:264:ILE:HG22	1:A:265:ILE:HG12	1.77	0.65
1:B:67:MET:HE1	1:B:305:GLN:HE22	1.60	0.65
1:B:56:GLU:CG	1:B:57:ASN:H	2.08	0.65
1:A:8:THR:CG2	1:A:11:SER:H	2.10	0.63
1:B:343:GLY:O	1:B:361:LYS:HE3	1.99	0.63
1:B:8:THR:HG22	1:B:11:SER:H	1.64	0.63
1:A:139:GLU:OE2	1:A:389:VAL:N	2.20	0.62
1:B:381:MET:HE1	1:B:414:PHE:HE1	1.64	0.62
1:B:178:TRP:HZ2	4:B:2105:HOH:O	1.80	0.62
1:B:315:PHE:HD2	1:B:316:MET:HE2	1.62	0.62
1:B:178:TRP:NE1	4:B:2105:HOH:O	2.34	0.60
1:B:263:LYS:NZ	2:B:1263:PLP:H5A2	2.16	0.60
1:B:264:ILE:O	1:B:318:HIS:HE1	1.84	0.59
1:B:369:LEU:CD1	1:B:420:LEU:HB3	2.26	0.59
1:A:131:PRO:HB2	4:A:2083:HOH:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:LEU:O	1:A:88:LYS:HG3	2.03	0.58
1:B:5:ARG:HD3	4:B:2005:HOH:O	2.04	0.58
1:B:95:HIS:HE1	4:B:2065:HOH:O	1.87	0.57
1:A:263:LYS:NZ	2:A:1263:PLP:H5A2	2.19	0.57
1:A:381:MET:HE3	1:A:400:ALA:HB2	1.87	0.57
1:A:374:ALA:HB2	1:A:420:LEU:CD1	2.35	0.56
1:A:382:LEU:HD23	1:A:383:PRO:HD2	1.87	0.56
1:B:340:TRP:CZ2	1:B:415:GLN:HG2	2.42	0.55
1:A:42:ASN:OD1	1:A:44:ASN:HB2	2.07	0.55
1:A:315:PHE:CE2	1:A:316:MET:HE2	2.39	0.55
1:B:178:TRP:CE2	4:B:2105:HOH:O	2.60	0.55
1:A:315:PHE:CE2	1:A:316:MET:CE	2.89	0.55
1:A:49:LYS:HB2	1:B:54:THR:HG22	1.89	0.54
1:A:419:GLN:HA	1:A:422:LYS:HD2	1.89	0.54
1:B:168:PRO:HD3	1:B:212:ARG:HD3	1.90	0.54
1:B:7:ILE:HD12	1:B:127:MET:HG2	1.90	0.54
1:B:56:GLU:CG	1:B:57:ASN:N	2.68	0.53
1:B:340:TRP:CE2	1:B:415:GLN:HG2	2.45	0.52
1:A:256:ILE:HD11	1:A:280:LEU:HD13	1.92	0.52
1:A:263:LYS:NZ	2:A:1263:PLP:C5A	2.72	0.52
1:A:3:TYR:O	1:A:7:ILE:HG12	2.10	0.52
1:A:140:PRO:HG2	1:A:204:THR:CG2	2.40	0.51
1:B:18:PRO:O	4:B:2018:HOH:O	2.19	0.51
1:A:77:SER:OG	1:A:289:GLN:HG2	2.11	0.50
1:A:315:PHE:HD2	1:A:316:MET:CE	2.22	0.50
1:A:365:ASP:HB2	1:A:394:PRO:HB3	1.93	0.50
1:B:366:VAL:HG12	1:B:370:ILE:HD12	1.93	0.50
1:B:7:ILE:HD11	1:B:280:LEU:HD22	1.94	0.50
1:B:8:THR:CG2	1:B:11:SER:H	2.25	0.49
1:B:240:LYS:HA	4:B:2143:HOH:O	2.11	0.49
1:A:7:ILE:HD12	1:A:283:ARG:HG3	1.95	0.49
1:A:326:TYR:CE1	1:A:354:MET:HE2	2.48	0.48
1:B:1:GLU:C	1:B:2:ASN:CA	2.83	0.48
1:A:377:MET:HE2	1:A:416:VAL:HG22	1.95	0.48
1:A:385:ASN:H	1:A:385:ASN:HD22	1.60	0.48
1:A:7:ILE:HA	4:A:2006:HOH:O	2.13	0.48
1:B:373:LYS:O	1:B:376:LYS:HB3	2.13	0.48
1:B:56:GLU:HG2	1:B:57:ASN:N	2.29	0.48
1:B:203:PRO:HB3	1:B:399:ARG:HD3	1.95	0.47
1:A:8:THR:HG22	1:A:11:SER:N	2.20	0.47
1:A:104:PRO:HD2	4:A:2070:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:ALA:O	1:B:84:LEU:HD21	2.14	0.47
1:B:67:MET:CE	1:B:305:GLN:HE22	2.25	0.47
1:A:334:LEU:HD11	1:A:349:VAL:HG23	1.97	0.47
1:B:326:TYR:CE1	1:B:354:MET:HE2	2.50	0.47
1:B:332:ALA:CB	1:B:410:MET:HE1	2.45	0.47
1:A:6:PHE:C	4:A:2006:HOH:O	2.57	0.46
1:A:184:LYS:NZ	4:A:2111:HOH:O	2.45	0.46
1:A:75:SER:HB2	1:A:76:PRO:CD	2.45	0.46
1:A:385:ASN:H	1:A:385:ASN:ND2	2.13	0.46
1:B:295:PRO:O	1:B:296:SER:C	2.58	0.46
1:B:56:GLU:HG2	1:B:57:ASN:H	1.78	0.46
1:B:59:LYS:HD2	1:B:309:GLU:CD	2.41	0.46
1:B:118:GLN:HE22	1:B:142:TYR:HE2	1.64	0.46
1:B:283:ARG:HD2	1:B:283:ARG:N	2.31	0.46
1:B:381:MET:HA	4:B:2195:HOH:O	2.15	0.46
1:B:233:TYR:OH	2:B:1263:PLP:O3	2.23	0.46
1:A:99:THR:CG2	1:A:109:MET:HB2	2.41	0.45
1:A:32:SER:HB2	1:B:380:LEU:HD22	1.98	0.45
1:A:99:THR:HG23	1:A:109:MET:N	2.31	0.45
1:A:304:SER:O	1:A:308:HIS:HD2	1.99	0.45
1:A:358:ILE:HD12	1:A:381:MET:HE1	1.98	0.45
1:A:139:GLU:HA	1:A:140:PRO:C	2.42	0.44
1:A:172:ARG:HD2	4:A:2101:HOH:O	2.17	0.44
1:B:239:ASN:HB3	4:B:2142:HOH:O	2.17	0.44
1:B:218:GLU:HG2	4:B:2127:HOH:O	2.17	0.44
1:B:84:LEU:O	1:B:88:LYS:HG3	2.18	0.44
1:B:97:PRO:HG2	1:B:109:MET:SD	2.58	0.44
1:B:382:LEU:HA	1:B:383:PRO:HD3	1.92	0.43
1:A:2:ASN:N	4:A:2003:HOH:O	2.52	0.43
1:A:2:ASN:N	1:A:2:ASN:OD1	2.51	0.43
1:B:197:VAL:HB	1:B:201:ASN:HD22	1.84	0.42
1:B:385:ASN:ND2	1:B:385:ASN:H	2.17	0.42
1:A:181:GLU:CD	4:A:2107:HOH:O	2.62	0.42
1:B:137:LEU:C	1:B:137:LEU:HD12	2.44	0.42
1:A:374:ALA:HB2	1:A:420:LEU:HD13	2.02	0.42
1:B:283:ARG:HD3	4:B:2153:HOH:O	2.20	0.42
1:B:8:THR:HG22	1:B:11:SER:HB2	2.01	0.42
1:A:211:GLU:H	1:A:211:GLU:CD	2.28	0.42
1:A:16:PRO:HA	1:A:286:LEU:HD22	2.02	0.42
1:A:137:LEU:HD22	1:A:156:ILE:CG2	2.50	0.42
1:A:381:MET:HE2	1:A:398:LEU:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:LEU:O	1:A:421:ILE:HG13	2.19	0.41
1:B:400:ALA:HA	4:B:2195:HOH:O	2.21	0.41
1:B:67:MET:HE1	1:B:305:GLN:NE2	2.30	0.41
1:A:134:ASN:HD22	1:A:155:ASN:HB2	1.85	0.40
1:A:147:GLN:O	1:A:151:PRO:HD3	2.21	0.40
1:B:129:ILE:HD13	1:B:129:ILE:HA	1.85	0.40
1:A:282:GLU:CD	1:A:283:ARG:HH21	2.29	0.40
1:B:107:GLY:HA2	1:B:277:PRO:HB3	2.02	0.40
1:A:137:LEU:CD2	1:A:156:ILE:CG2	2.99	0.40
1:B:313:GLU:HB2	4:B:2163:HOH:O	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	408/427 (96%)	398 (98%)	8 (2%)	2 (0%)	24	31
1	B	410/427 (96%)	401 (98%)	8 (2%)	1 (0%)	43	55
All	All	818/854 (96%)	799 (98%)	16 (2%)	3 (0%)	30	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	56	GLU
1	A	423	GLU
1	A	266	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	360/372 (97%)	343 (95%)	17 (5%)	23	35
1	B	360/372 (97%)	342 (95%)	18 (5%)	22	33
All	All	720/744 (97%)	685 (95%)	35 (5%)	22	34

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ARG
1	A	8	THR
1	A	32	SER
1	A	99	THR
1	A	111	LEU
1	A	137	LEU
1	A	156	ILE
1	A	196	THR
1	A	240	LYS
1	A	282	GLU
1	A	320	ASP
1	A	363	ILE
1	A	368	GLU
1	A	370	ILE
1	A	375	VAL
1	A	382	LEU
1	A	417	LEU
1	B	7	ILE
1	B	8	THR
1	B	32	SER
1	B	54	THR
1	B	56	GLU
1	B	62	GLN
1	B	135	VAL
1	B	137	LEU
1	B	169	ASP
1	B	181	GLU

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Mol	Chain	Res	Type
1	B	283	ARG
1	B	313	GLU
1	B	320	ASP
1	B	339	LYS
1	B	368	GLU
1	B	376	LYS
1	B	398	LEU
1	B	399	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	44	ASN
1	A	101	HIS
1	A	134	ASN
1	A	150	HIS
1	A	308	HIS
1	A	328	ASN
1	A	348	HIS
1	A	385	ASN
1	B	44	ASN
1	B	95	HIS
1	B	101	HIS
1	B	106	GLN
1	B	118	GLN
1	B	134	ASN
1	B	150	HIS
1	B	155	ASN
1	B	201	ASN
1	B	305	GLN
1	B	318	HIS
1	B	385	ASN

5.3.3 RNA ⓘ

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 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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$
2	PLP	A	1263	1	15,15,16	2.01	4 (26%)	21,22,23	1.69	4 (19%)
2	PLP	B	1263	1	15,15,16	2.11	4 (26%)	21,22,23	1.80	4 (19%)

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	1263	1	-	0/6/6/8	0/1/1/1
2	PLP	B	1263	1	-	3/6/6/8	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1263	PLP	O3-C3	-5.61	1.24	1.36
2	B	1263	PLP	O3-C3	-5.42	1.24	1.36
2	B	1263	PLP	C2-N1	3.29	1.39	1.33
2	B	1263	PLP	C3-C2	-3.06	1.37	1.41
2	A	1263	PLP	C3-C2	-2.97	1.37	1.41
2	A	1263	PLP	C2-N1	2.97	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1263	PLP	C6-N1	2.13	1.38	1.34
2	A	1263	PLP	P-O3P	-2.10	1.47	1.54

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1263	PLP	O4P-C5A-C5	4.76	118.28	109.36
2	A	1263	PLP	O4P-C5A-C5	4.74	118.25	109.36
2	B	1263	PLP	C4A-C4-C3	-2.77	115.90	120.52
2	A	1263	PLP	O3P-P-O4P	-2.66	99.73	106.67
2	B	1263	PLP	C5-C6-N1	-2.51	119.75	123.83
2	A	1263	PLP	O3P-P-O2P	2.47	117.06	107.80
2	A	1263	PLP	C5-C6-N1	-2.17	120.30	123.83
2	B	1263	PLP	C3-C4-C5	2.11	121.12	118.59

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1263	PLP	C5A-O4P-P-O2P
2	B	1263	PLP	C5A-O4P-P-O3P
2	B	1263	PLP	C5A-O4P-P-O1P

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1263	PLP	2	0
2	B	1263	PLP	2	0

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
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	-1:GLU	C	2:ASN	N	2.46
1	B	-1:GLU	C	2:ASN	N	1.98

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	414/427 (96%)	-0.07	6 (1%) 73 75	26, 38, 70, 87	0
1	B	414/427 (96%)	-0.06	6 (1%) 73 75	27, 38, 63, 79	0
All	All	828/854 (96%)	-0.06	12 (1%) 73 75	26, 38, 66, 87	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	374	ALA	2.9
1	B	425	LEU	2.9
1	A	241	PHE	2.8
1	A	33	MET	2.7
1	B	370	ILE	2.6
1	A	374	ALA	2.5
1	B	366	VAL	2.5
1	A	366	VAL	2.3
1	B	241	PHE	2.3
1	B	376	LYS	2.1
1	A	-1	GLU	2.0
1	A	425	LEU	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 [i](#)

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($\text{\AA}^2$)	Q<0.9
2	PLP	A	1263	15/16	0.97	0.07	37,47,49,50	0
2	PLP	B	1263	15/16	0.97	0.07	40,43,45,45	0
3	IOD	A	1426	1/1	1.00	0.06	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.