



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 03:07 PM UTC

PDB ID : 5HSJ / pdb_00005hsj
Title : Structure of tyrosine decarboxylase complex with PLP at 1.9 Angstroms resolution
Authors : Ni, Y.; Zhou, J.; Zhu, H.; Zhang, K.
Deposited on : 2016-01-25
Resolution : 1.90 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

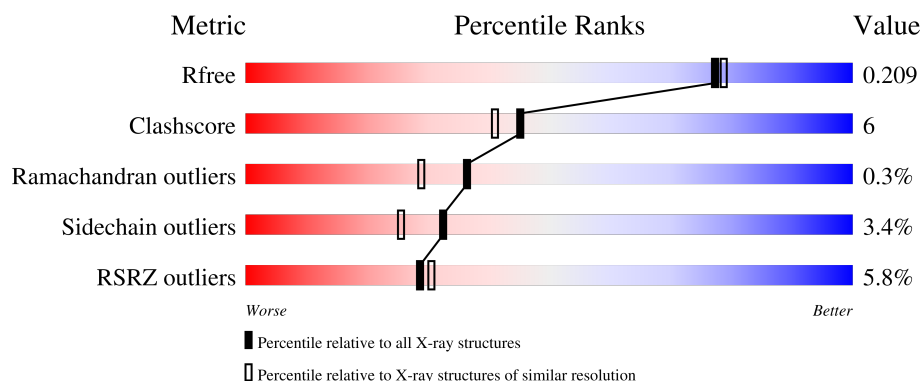
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	634	 7% 80% 13% . .
1	B	634	 4% 81% 11% . 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10136 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.

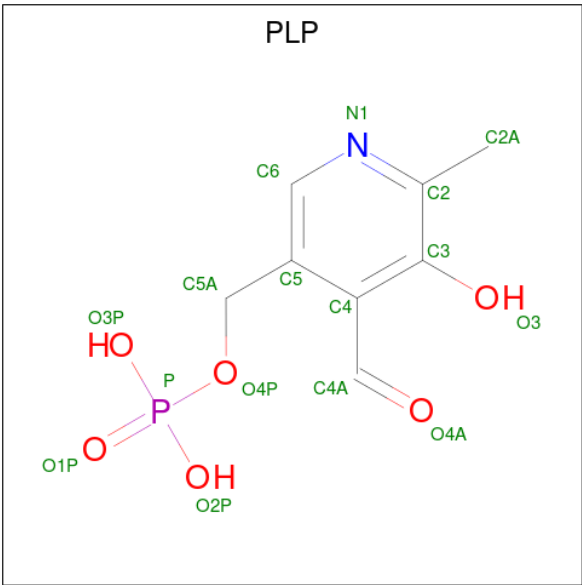
- Molecule 1 is a protein called Putative decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	606	Total	C	N	O	S	0	0	0
			4807	3078	805	906	18			
1	B	595	Total	C	N	O	S	0	0	0
			4732	3029	794	891	18			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	627	LEU	-	expression tag	UNP J7GQ11
A	628	GLU	-	expression tag	UNP J7GQ11
A	629	HIS	-	expression tag	UNP J7GQ11
A	630	HIS	-	expression tag	UNP J7GQ11
A	631	HIS	-	expression tag	UNP J7GQ11
A	632	HIS	-	expression tag	UNP J7GQ11
A	633	HIS	-	expression tag	UNP J7GQ11
A	634	HIS	-	expression tag	UNP J7GQ11
B	627	LEU	-	expression tag	UNP J7GQ11
B	628	GLU	-	expression tag	UNP J7GQ11
B	629	HIS	-	expression tag	UNP J7GQ11
B	630	HIS	-	expression tag	UNP J7GQ11
B	631	HIS	-	expression tag	UNP J7GQ11
B	632	HIS	-	expression tag	UNP J7GQ11
B	633	HIS	-	expression tag	UNP J7GQ11
B	634	HIS	-	expression tag	UNP J7GQ11

- 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		

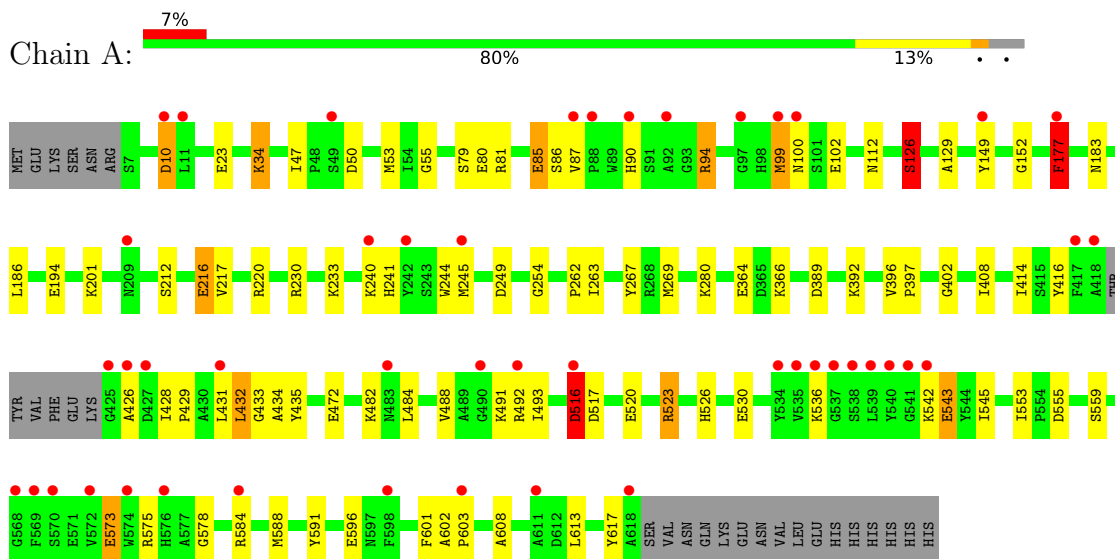
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	217	Total	O	0	0
			217	217		
3	B	350	Total	O	0	0
			350	350		

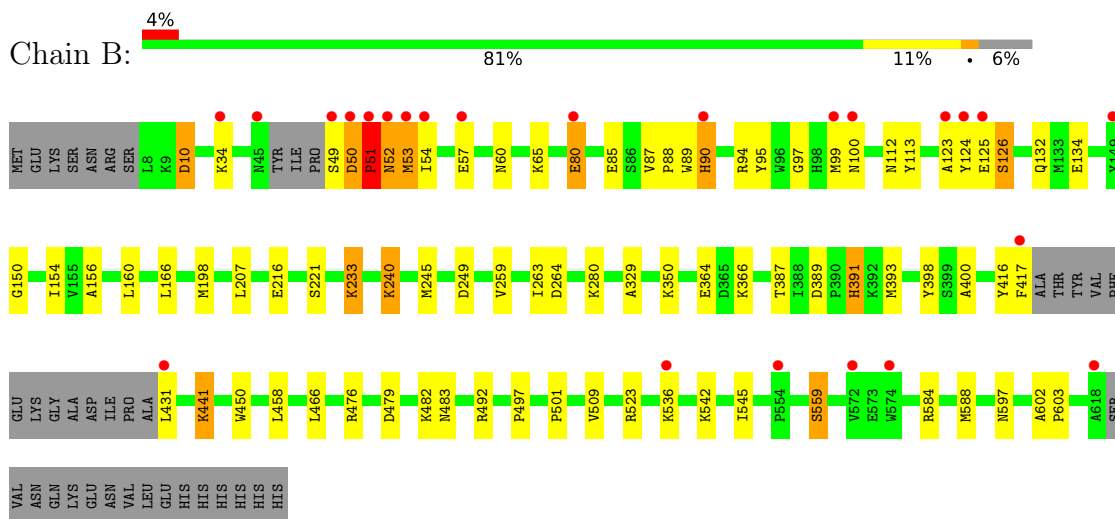
3 Residue-property plots [i](#)

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: Putative decarboxylase



• Molecule 1: Putative decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.87Å 126.83Å 82.88Å 90.00° 109.73° 90.00°	Depositor
Resolution (Å)	49.26 – 1.90 49.26 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.26-1.90) 99.7 (49.26-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.172 , 0.212 0.169 , 0.209	Depositor DCC
R_{free} test set	4732 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.592	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10136	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: 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	1.33	10/4925 (0.2%)	1.21	26/6673 (0.4%)
1	B	1.44	21/4846 (0.4%)	1.23	18/6561 (0.3%)
All	All	1.39	31/9771 (0.3%)	1.22	44/13234 (0.3%)

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	509	VAL	N-CA	7.59	1.54	1.46
1	A	402	GLY	N-CA	7.25	1.52	1.45
1	B	387	THR	CA-CB	7.03	1.63	1.53
1	A	244	TRP	N-CA	7.01	1.54	1.46
1	B	90	HIS	N-CA	-7.01	1.37	1.46
1	B	90	HIS	CB-CG	-6.68	1.40	1.50
1	B	221	SER	C-O	-6.41	1.17	1.23
1	B	233	LYS	C-O	6.11	1.31	1.23
1	A	262	PRO	C-O	6.10	1.31	1.23
1	B	94	ARG	N-CA	6.09	1.54	1.46
1	A	152	GLY	N-CA	6.07	1.52	1.45
1	B	150	GLY	C-O	-6.05	1.18	1.23
1	A	217	VAL	CA-C	5.91	1.60	1.52
1	B	559	SER	C-N	5.83	1.38	1.33
1	B	501	PRO	CA-C	5.72	1.59	1.52
1	B	90	HIS	CA-CB	-5.66	1.44	1.53
1	B	97	GLY	C-O	-5.54	1.19	1.24
1	B	87	VAL	CA-C	5.54	1.57	1.52
1	B	329	ALA	CA-CB	5.40	1.60	1.53
1	A	269	MET	SD-CE	-5.35	1.66	1.79
1	B	132	GLN	CD-NE2	-5.35	1.22	1.33
1	A	112	ASN	N-CA	5.28	1.52	1.46
1	A	233	LYS	C-O	5.24	1.29	1.23
1	B	154	ILE	N-CA	-5.13	1.40	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	450	TRP	N-CA	5.12	1.52	1.46
1	B	264	ASP	C-O	-5.12	1.18	1.23
1	B	466	LEU	N-CA	-5.10	1.40	1.46
1	B	398	TYR	CA-CB	5.09	1.62	1.53
1	A	516	ASP	N-CA	5.06	1.52	1.46
1	B	259	VAL	CA-CB	5.05	1.58	1.54
1	A	241	HIS	CA-C	5.04	1.59	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	516	ASP	CB-CA-C	-8.79	95.71	110.56
1	A	523	ARG	CG-CD-NE	-8.25	93.86	112.00
1	A	216	GLU	CA-CB-CG	-8.15	97.80	114.10
1	B	90	HIS	N-CA-CB	-8.07	97.34	109.69
1	B	523	ARG	CG-CD-NE	-7.91	94.60	112.00
1	B	280	LYS	CD-CE-NZ	-7.60	87.57	111.90
1	B	80	GLU	CA-CB-CG	-7.57	98.95	114.10
1	A	575	ARG	NE-CZ-NH2	7.54	125.99	119.20
1	A	596	GLU	CB-CA-C	-6.81	99.92	110.81
1	B	94	ARG	NE-CZ-NH2	6.79	125.31	119.20
1	A	389	ASP	CB-CA-C	-6.66	102.77	110.17
1	A	254	GLY	N-CA-C	6.63	120.51	112.48
1	A	80	GLU	CB-CA-C	-6.44	100.77	110.88
1	B	483	ASN	N-CA-CB	-6.33	102.06	111.43
1	B	90	HIS	N-CA-C	6.28	119.45	110.10
1	A	575	ARG	NE-CZ-NH1	-6.26	115.24	121.50
1	B	476	ARG	N-CA-C	-6.16	104.65	111.36
1	B	559	SER	CA-C-N	-6.15	114.55	120.83
1	B	559	SER	C-N-CA	-6.15	114.55	120.83
1	A	591	TYR	N-CA-C	6.05	120.50	113.18
1	B	350	LYS	CB-CA-C	-6.03	99.09	110.67
1	A	244	TRP	N-CA-C	5.99	117.89	111.36
1	B	492	ARG	CB-CG-CD	5.94	124.96	111.30
1	A	50	ASP	C-N-CD	5.82	133.41	120.60
1	A	194	GLU	N-CA-C	5.78	117.25	111.07
1	A	90	HIS	CA-C-N	5.69	130.24	122.84
1	A	90	HIS	C-N-CA	5.69	130.24	122.84
1	A	34	LYS	CB-CA-C	-5.59	102.11	110.88
1	A	472	GLU	N-CA-C	-5.58	104.83	111.69
1	B	52	ASN	N-CA-C	5.56	122.65	110.80
1	A	126	SER	N-CA-C	5.52	118.22	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	PHE	N-CA-C	5.45	117.22	111.28
1	A	201	LYS	CG-CD-CE	-5.38	98.93	111.30
1	B	94	ARG	NE-CZ-NH1	-5.34	116.16	121.50
1	A	484	LEU	CD1-CG-CD2	-5.33	99.06	110.80
1	B	350	LYS	CB-CG-CD	-5.27	99.18	111.30
1	A	482	LYS	CA-C-N	-5.12	111.77	121.54
1	A	482	LYS	C-N-CA	-5.12	111.77	121.54
1	A	555	ASP	N-CA-C	-5.09	107.20	113.41
1	A	55	GLY	CA-C-N	5.08	124.86	119.28
1	A	55	GLY	C-N-CA	5.08	124.86	119.28
1	B	389	ASP	CB-CA-C	-5.07	103.04	110.13
1	B	458	LEU	CD1-CG-CD2	-5.03	99.73	110.80
1	B	89	TRP	CA-C-O	5.00	127.01	121.56

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4807	0	4700	67	0
1	B	4732	0	4626	55	0
2	A	15	0	7	4	0
2	B	15	0	7	1	0
3	A	217	0	0	3	0
3	B	350	0	0	8	0
All	All	10136	0	9340	110	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 6.

All (110) 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:392:LYS:HZ1	2:A:701:PLP:C4A	1.26	1.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:491:LYS:HD2	1:A:617:TYR:HE2	1.16	1.05
1:A:94:ARG:NH2	1:B:50:ASP:O	1.89	1.04
1:A:126:SER:HB2	1:B:99:MET:HE1	1.42	1.02
1:B:588:MET:SD	3:B:1096:HOH:O	2.19	1.00
1:B:90:HIS:CE1	1:B:95:TYR:CD2	2.50	1.00
1:B:53:MET:HA	1:B:53:MET:HE2	1.45	0.95
1:A:10:ASP:N	1:A:10:ASP:OD1	2.01	0.94
1:A:53:MET:HE3	1:A:53:MET:HA	1.48	0.93
1:A:99:MET:HE1	1:B:126:SER:HB2	1.50	0.93
1:A:491:LYS:HD2	1:A:617:TYR:CE2	2.04	0.93
1:A:392:LYS:HZ3	2:A:701:PLP:C4A	1.65	0.92
1:A:416:TYR:O	1:A:434:ALA:HB1	1.69	0.92
1:A:177:PHE:CE2	1:A:230:ARG:CZ	2.53	0.91
1:A:177:PHE:CZ	1:A:230:ARG:NH2	2.46	0.83
1:A:408:ILE:HD13	1:A:426:ALA:HB1	1.62	0.79
1:B:391:HIS:CD2	1:B:400:ALA:H	2.01	0.79
1:B:53:MET:HA	1:B:53:MET:CE	2.11	0.79
1:A:94:ARG:HH11	1:A:94:ARG:CG	1.97	0.78
1:A:177:PHE:HE2	1:A:230:ARG:CZ	1.97	0.78
1:A:47:ILE:HD12	1:B:88:PRO:HB3	1.67	0.76
1:B:90:HIS:CE1	1:B:95:TYR:HD2	2.02	0.76
1:A:94:ARG:HH11	1:A:94:ARG:HG2	1.50	0.75
1:A:392:LYS:HZ1	2:A:701:PLP:C4	1.99	0.75
1:A:392:LYS:NZ	2:A:701:PLP:C4	2.50	0.73
1:B:391:HIS:HD2	1:B:400:ALA:H	1.36	0.72
1:A:126:SER:HB2	1:B:99:MET:CE	2.19	0.71
1:A:177:PHE:HZ	1:A:230:ARG:NH2	1.89	0.71
1:A:542:LYS:HB2	1:A:545:ILE:HD11	1.72	0.70
1:B:431:LEU:N	3:B:803:HOH:O	2.25	0.69
1:A:177:PHE:CE2	1:A:230:ARG:NH1	2.61	0.69
1:A:99:MET:HE1	1:B:126:SER:CB	2.23	0.67
1:B:49:SER:C	1:B:50:ASP:OD1	2.37	0.67
1:A:126:SER:CB	1:B:99:MET:HE1	2.24	0.66
1:B:90:HIS:HE1	1:B:95:TYR:CD2	2.13	0.65
1:A:99:MET:CE	1:B:126:SER:HB2	2.25	0.62
1:B:584:ARG:HB3	1:B:584:ARG:CZ	2.28	0.62
1:B:10:ASP:N	1:B:10:ASP:OD1	2.33	0.61
1:B:90:HIS:C	1:B:90:HIS:CD2	2.77	0.61
1:A:53:MET:HA	1:A:53:MET:CE	2.29	0.59
1:B:134:GLU:OE2	1:B:441:LYS:NZ	2.35	0.59
1:B:240:LYS:HE2	1:B:245:MET:HG3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:ASN:O	1:B:65:LYS:HE3	2.03	0.58
1:A:53:MET:HE3	1:A:53:MET:CA	2.27	0.57
1:A:94:ARG:NH1	3:A:808:HOH:O	2.37	0.56
1:A:100:ASN:OD1	1:A:102:GLU:N	2.38	0.56
1:B:364:GLU:OE1	1:B:366:LYS:HE2	2.05	0.56
1:B:100:ASN:ND2	3:B:802:HOH:O	2.09	0.55
1:A:516:ASP:N	1:A:516:ASP:OD1	2.33	0.55
1:B:417:PHE:N	1:B:417:PHE:CD1	2.74	0.55
1:A:177:PHE:CZ	1:A:230:ARG:CZ	2.87	0.55
1:A:129:ALA:HB2	1:B:85:GLU:HB2	1.89	0.54
1:B:90:HIS:HE1	1:B:95:TYR:HD2	1.50	0.54
1:B:417:PHE:N	1:B:417:PHE:HD1	2.05	0.54
1:B:49:SER:O	1:B:50:ASP:OD1	2.24	0.54
1:B:216:GLU:HG2	3:B:997:HOH:O	2.07	0.53
1:A:364:GLU:OE1	1:A:366:LYS:CE	2.57	0.53
1:A:364:GLU:OE1	1:A:366:LYS:HE3	2.09	0.53
1:B:393:MET:SD	1:B:588:MET:HG2	2.49	0.52
1:B:542:LYS:HB2	1:B:545:ILE:HD11	1.91	0.52
1:A:542:LYS:HB2	1:A:545:ILE:CD1	2.40	0.52
1:A:429:PRO:HG3	1:A:435:TYR:CD2	2.44	0.52
1:A:396:VAL:HG13	1:A:397:PRO:HD2	1.92	0.51
1:B:391:HIS:HE1	2:B:702:PLP:O3P	1.93	0.51
1:A:23:GLU:HG2	1:B:54:ILE:O	2.10	0.51
1:A:34:LYS:HD2	3:A:1006:HOH:O	2.10	0.51
1:A:431:LEU:C	1:A:433:GLY:N	2.67	0.50
1:B:156:ALA:HB3	1:B:160:LEU:HD12	1.92	0.50
1:A:573:GLU:HA	1:A:573:GLU:OE1	2.12	0.50
1:B:416:TYR:C	1:B:417:PHE:HD1	2.19	0.50
1:A:416:TYR:O	1:A:434:ALA:CB	2.52	0.49
1:A:94:ARG:CG	1:A:94:ARG:NH1	2.67	0.49
1:A:584:ARG:HB3	1:A:584:ARG:CZ	2.43	0.48
1:A:488:VAL:HG21	1:A:613:LEU:HB3	1.98	0.46
1:A:53:MET:HE1	1:B:597:ASN:CG	2.41	0.46
1:A:79:SER:HA	1:B:113:TYR:CE2	2.51	0.46
1:B:263:ILE:HG13	1:B:559:SER:HB3	1.98	0.45
1:B:198:MET:HB2	1:B:198:MET:HE3	1.72	0.44
1:B:99:MET:HE2	1:B:99:MET:HB3	1.74	0.44
1:B:207:LEU:HD12	1:B:207:LEU:HA	1.87	0.44
1:A:526:HIS:O	1:A:530:GLU:HG3	2.18	0.44
1:A:429:PRO:HG3	1:A:435:TYR:HD2	1.81	0.44
1:B:112:ASN:CG	3:B:812:HOH:O	2.60	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:429:PRO:CG	1:A:435:TYR:HD2	2.30	0.44
1:A:263:ILE:HD12	1:A:267:TYR:HA	1.99	0.44
1:A:428:ILE:HA	1:A:429:PRO:HD3	1.81	0.44
1:B:100:ASN:HB3	3:B:1096:HOH:O	2.18	0.43
1:A:249:ASP:OD1	1:A:249:ASP:C	2.61	0.43
1:A:216:GLU:HG3	3:A:966:HOH:O	2.19	0.43
1:A:149:TYR:CD1	1:A:149:TYR:N	2.87	0.42
1:B:416:TYR:C	1:B:417:PHE:CD1	2.97	0.42
1:B:240:LYS:HE2	1:B:245:MET:CG	2.49	0.42
1:B:479:ASP:HA	1:B:482:LYS:HE2	2.00	0.42
1:A:183:ASN:HB3	1:A:186:LEU:HD12	2.02	0.42
1:A:429:PRO:HB3	1:A:435:TYR:CD2	2.55	0.42
1:A:602:ALA:HB3	1:A:603:PRO:HD3	2.02	0.42
1:B:233:LYS:NZ	3:B:816:HOH:O	2.41	0.42
1:B:482:LYS:HB3	1:B:497:PRO:HG2	2.01	0.42
1:B:123:ALA:O	1:B:124:TYR:C	2.62	0.41
1:A:81:ARG:O	1:A:85:GLU:CG	2.68	0.41
1:A:553:ILE:HD12	1:A:578:GLY:HA2	2.01	0.41
1:A:280:LYS:HE2	1:A:280:LYS:HB3	1.85	0.41
1:A:94:ARG:NH1	1:A:601:PHE:HE1	2.19	0.41
1:A:263:ILE:HG13	1:A:559:SER:HB3	2.02	0.41
1:A:543:GLU:HB3	1:A:608:ALA:HB2	2.03	0.41
1:A:431:LEU:O	1:A:432:LEU:C	2.62	0.41
1:B:80:GLU:CD	3:B:819:HOH:O	2.63	0.40
1:A:517:ASP:HB3	1:A:520:GLU:HB2	2.03	0.40
1:B:602:ALA:HB3	1:B:603:PRO:HD3	2.03	0.40
1:B:51:PRO:O	1:B:51:PRO:CD	2.69	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	602/634 (95%)	584 (97%)	16 (3%)	2 (0%)	36	29
1	B	589/634 (93%)	576 (98%)	11 (2%)	2 (0%)	36	29
All	All	1191/1268 (94%)	1160 (97%)	27 (2%)	4 (0%)	36	29

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	LYS
1	B	52	ASN
1	A	432	LEU
1	B	51	PRO

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	502/530 (95%)	482 (96%)	20 (4%)	28	20
1	B	495/530 (93%)	481 (97%)	14 (3%)	38	32
All	All	997/1060 (94%)	963 (97%)	34 (3%)	32	25

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

Mol	Chain	Res	Type
1	A	10	ASP
1	A	85	GLU
1	A	86	SER
1	A	87	VAL
1	A	94	ARG
1	A	99	MET
1	A	126	SER
1	A	177	PHE
1	A	212	SER
1	A	220	ARG
1	A	245	MET
1	A	414	ILE

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Mol	Chain	Res	Type
1	A	492	ARG
1	A	493	ILE
1	A	516	ASP
1	A	523	ARG
1	A	536	LYS
1	A	543	GLU
1	A	573	GLU
1	A	588	MET
1	B	10	ASP
1	B	34	LYS
1	B	50	ASP
1	B	51	PRO
1	B	53	MET
1	B	57	GLU
1	B	125	GLU
1	B	126	SER
1	B	166	LEU
1	B	240	LYS
1	B	249	ASP
1	B	391	HIS
1	B	441	LYS
1	B	536	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	353	GLN
1	A	548	HIS
1	A	576	HIS
1	B	33	ASN
1	B	90	HIS
1	B	100	ASN
1	B	317	GLN
1	B	391	HIS
1	B	526	HIS

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](#)

2 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	$\# Z > 2$
2	PLP	A	701	1	15,15,16	3.05	4 (26%)	21,22,23	1.99	7 (33%)
2	PLP	B	702	1	15,15,16	1.60	3 (20%)	21,22,23	2.19	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	701	1	-	0/6/6/8	0/1/1/1
2	PLP	B	702	1	-	0/6/6/8	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	PLP	C3-C2	7.58	1.48	1.41
2	A	701	PLP	C5-C4	6.39	1.47	1.40
2	A	701	PLP	C3-C4	4.94	1.49	1.40
2	B	702	PLP	C5-C4	4.16	1.45	1.40
2	B	702	PLP	C3-C2	2.35	1.43	1.41
2	A	701	PLP	C6-C5	2.33	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	702	PLP	C3-C4	2.25	1.44	1.40

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	702	PLP	C4A-C4-C5	6.62	127.76	120.94
2	A	701	PLP	O3-C3-C2	4.15	126.19	117.58
2	A	701	PLP	C6-N1-C2	3.92	126.31	119.20
2	A	701	PLP	C5A-C5-C4	-3.23	116.26	122.64
2	B	702	PLP	C4A-C4-C3	-3.14	115.29	120.52
2	A	701	PLP	C5A-C5-C6	2.95	124.16	119.36
2	A	701	PLP	C3-C2-N1	-2.90	117.31	120.96
2	B	702	PLP	O3P-P-O4P	-2.69	99.65	106.67
2	B	702	PLP	C6-N1-C2	2.48	123.70	119.20
2	B	702	PLP	C6-C5-C4	2.32	120.00	118.10
2	A	701	PLP	C5-C6-N1	-2.24	120.19	123.83
2	A	701	PLP	C4A-C4-C5	2.20	123.20	120.94
2	B	702	PLP	C2A-C2-N1	2.12	121.64	117.64
2	B	702	PLP	C5A-C5-C6	-2.10	115.93	119.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	PLP	4	0
2	B	702	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 ⓘ

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	606/634 (95%)	0.33	46 (7%)	20 21	18, 31, 64, 150	0
1	B	595/634 (93%)	-0.06	24 (4%)	42 45	15, 25, 52, 144	0
All	All	1201/1268 (94%)	0.14	70 (5%)	29 30	15, 27, 58, 150	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	535	VAL	8.7
1	B	431	LEU	6.0
1	A	87	VAL	5.3
1	B	51	PRO	5.2
1	A	537	GLY	5.0
1	A	568	GLY	4.8
1	B	54	ILE	4.5
1	A	534	TYR	4.2
1	B	123	ALA	3.9
1	A	418	ALA	3.9
1	A	177	PHE	3.8
1	A	417	PHE	3.8
1	A	569	PHE	3.8
1	A	539	LEU	3.6
1	A	540	TYR	3.5
1	B	49	SER	3.4
1	A	538	SER	3.4
1	A	483	ASN	3.3
1	A	92	ALA	3.3
1	B	53	MET	3.3
1	B	99	MET	3.3
1	A	10	ASP	3.3
1	B	100	ASN	3.2
1	B	90	HIS	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	618	ALA	3.1
1	A	88	PRO	3.1
1	A	245	MET	3.1
1	B	45	ASN	3.0
1	A	536	LYS	3.0
1	B	536	LYS	3.0
1	B	149	TYR	2.9
1	A	431	LEU	2.9
1	A	541	GLY	2.9
1	A	242	TYR	2.8
1	A	576	HIS	2.8
1	A	574	TRP	2.8
1	A	572	VAL	2.8
1	A	611	ALA	2.6
1	B	124	TYR	2.6
1	A	427	ASP	2.6
1	A	209	ASN	2.6
1	B	52	ASN	2.6
1	B	125	GLU	2.5
1	A	570	SER	2.5
1	A	492	ARG	2.5
1	B	50	ASP	2.5
1	A	97	GLY	2.4
1	B	572	VAL	2.4
1	A	240	LYS	2.4
1	A	516	ASP	2.3
1	A	425	GLY	2.3
1	A	49	SER	2.3
1	A	598	PHE	2.3
1	B	57	GLU	2.3
1	B	34	LYS	2.3
1	B	618	ALA	2.2
1	A	542	LYS	2.2
1	A	90	HIS	2.2
1	A	100	ASN	2.2
1	A	99	MET	2.2
1	B	417	PHE	2.1
1	A	490	GLY	2.1
1	B	574	TRP	2.1
1	A	426	ALA	2.1
1	B	80	GLU	2.0
1	A	603	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	2.0
1	A	584	ARG	2.0
1	A	149	TYR	2.0
1	B	554	PRO	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	701	15/16	0.96	0.07	26,31,39,46	0
2	PLP	B	702	15/16	0.98	0.05	19,22,25,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.