



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:42 AM UTC

PDB ID : 1L0C / pdb_0000110c
Title : Investigation of the Roles of Catalytic Residues in Serotonin N-Acetyltransferase
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Deposited on : 2002-02-08
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

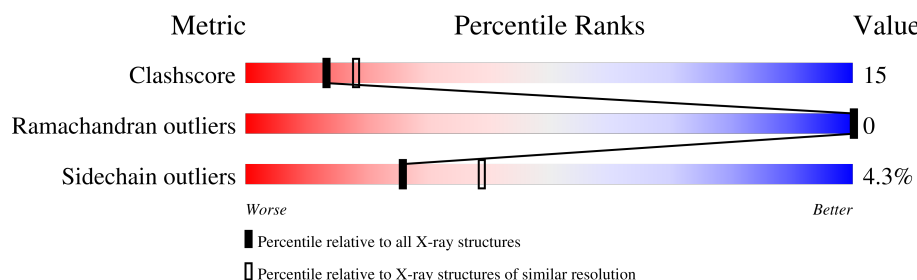
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	207	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1464 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.

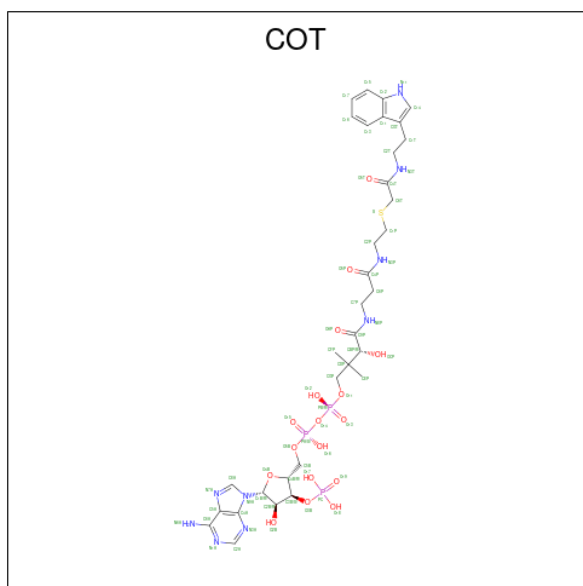
- Molecule 1 is a protein called Serotonin N-acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	166	1311	840	237	226	8	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	168	PHE	TYR	engineered mutation	UNP Q29495

- Molecule 2 is COA-S-ACETYL TRYPTAMINE (CCD ID: COT) (formula: $C_{33}H_{48}N_9O_{17}P_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P	S	
2	A	1	63	33	9	17	3	1	0

- Molecule 3 is water.

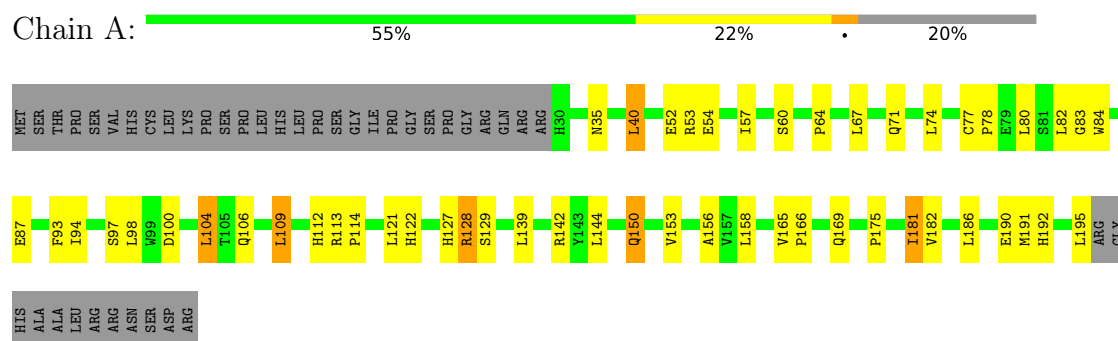
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	90	Total 90	O 90	0	0

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: Serotonin N-acetyltransferase



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	53.21Å 68.68Å 89.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.30)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.193 , 0.246	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1464	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: COT

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	0.40	0/1349	0.92	3/1833 (0.2%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	150	GLN	CA-C-N	5.67	126.92	119.84
1	A	150	GLN	C-N-CA	5.67	126.92	119.84
1	A	181	ILE	N-CA-C	5.02	116.86	109.63

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	1311	0	1280	40	0
2	A	63	0	41	0	0
3	A	90	0	0	3	0
All	All	1464	0	1321	40	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 15.

All (40) 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:128:ARG:HH11	1:A:128:ARG:HB3	1.19	1.01
1:A:35:ASN:HD22	1:A:84:TRP:HE1	1.07	0.94
1:A:74:LEU:O	1:A:78:PRO:HG3	1.77	0.84
1:A:128:ARG:HB3	1:A:128:ARG:NH1	1.97	0.80
1:A:166:PRO:HA	1:A:169:GLN:HE21	1.50	0.76
1:A:67:LEU:O	1:A:71:GLN:HG3	1.96	0.66
1:A:106:GLN:HE22	1:A:182:VAL:H	1.43	0.66
1:A:128:ARG:HH11	1:A:128:ARG:CB	2.03	0.65
1:A:127:HIS:CD2	1:A:129:SER:H	2.15	0.65
1:A:53:ARG:HA	1:A:57:ILE:HD12	1.79	0.64
1:A:127:HIS:HD2	1:A:129:SER:H	1.47	0.61
1:A:35:ASN:ND2	1:A:84:TRP:HE1	1.90	0.61
1:A:97:SER:OG	1:A:122:HIS:HE1	1.84	0.60
1:A:166:PRO:HG2	3:A:976:HOH:O	2.02	0.60
1:A:35:ASN:HD22	1:A:84:TRP:NE1	1.91	0.60
1:A:158:LEU:C	1:A:158:LEU:HD12	2.29	0.57
1:A:190:GLU:OE2	1:A:192:HIS:HE1	1.87	0.56
1:A:122:HIS:HD2	3:A:902:HOH:O	1.87	0.56
1:A:175:PRO:HA	1:A:191:MET:HG2	1.90	0.54
1:A:60:SER:CB	1:A:186:LEU:HD11	2.37	0.53
1:A:112:HIS:O	1:A:114:PRO:HD3	2.09	0.53
1:A:100:ASP:OD1	1:A:113:ARG:NH1	2.42	0.53
1:A:150:GLN:NE2	3:A:922:HOH:O	2.45	0.49
1:A:156:ALA:HB2	1:A:195:LEU:HD11	1.94	0.49
1:A:106:GLN:O	1:A:109:LEU:HB2	2.13	0.49
1:A:153:VAL:HG12	1:A:195:LEU:HD12	1.95	0.48
1:A:82:LEU:HD21	1:A:142:ARG:HG2	1.95	0.47
1:A:80:LEU:HD11	1:A:98:LEU:HG	1.95	0.47
1:A:60:SER:HB2	1:A:186:LEU:HD11	1.99	0.45
1:A:94:ILE:HG13	1:A:139:LEU:HD23	1.97	0.45
1:A:109:LEU:HD13	1:A:181:ILE:HD13	1.98	0.44
1:A:165:VAL:O	1:A:169:GLN:HG3	2.17	0.44
1:A:93:PHE:C	1:A:93:PHE:CD1	2.97	0.42
1:A:52:GLU:OE1	1:A:64:PRO:HG2	2.20	0.42
1:A:104:LEU:HD12	1:A:104:LEU:HA	1.86	0.42
1:A:54:GLU:CD	1:A:54:GLU:C	2.88	0.41
1:A:77:CYS:N	1:A:78:PRO:HD3	2.35	0.41
1:A:40:LEU:HD11	1:A:83:GLY:HA3	2.01	0.41
1:A:144:LEU:HD22	1:A:195:LEU:HD21	2.02	0.40
1:A:166:PRO:HA	1:A:169:GLN:NE2	2.29	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	164/207 (79%)	160 (98%)	4 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	139/174 (80%)	133 (96%)	6 (4%)	26	39

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LEU
1	A	87	GLU
1	A	104	LEU
1	A	109	LEU
1	A	121	LEU
1	A	128	ARG

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	35	ASN
1	A	71	GLN
1	A	72	HIS
1	A	106	GLN
1	A	122	HIS
1	A	127	HIS
1	A	133	GLN
1	A	150	GLN
1	A	169	GLN
1	A	192	HIS

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	COT	A	900	-	65,67,67	2.24	20 (30%)	92,98,98	1.38	6 (6%)

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	COT	A	900	-	-	10/58/74/74	0/5/5/5

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	COT	PA-O14	-6.83	1.52	1.59
2	A	900	COT	PB-O14	-5.86	1.53	1.59
2	A	900	COT	C1P-C2P	-5.79	1.27	1.51
2	A	900	COT	C4A-N3A	5.73	1.45	1.34
2	A	900	COT	C4P-N3P	4.83	1.44	1.33
2	A	900	COT	C9P-N8P	4.41	1.44	1.33
2	A	900	COT	O2B-C2B	-3.60	1.34	1.43
2	A	900	COT	C11-C12	3.37	1.45	1.41
2	A	900	COT	C5A-C6A	3.14	1.49	1.41
2	A	900	COT	C17-C16	2.85	1.44	1.38
2	A	900	COT	C16-C13	2.64	1.43	1.38
2	A	900	COT	O4B-C1B	2.50	1.47	1.42
2	A	900	COT	C13-C11	2.49	1.43	1.39
2	A	900	COT	C1B-N9A	2.43	1.53	1.46
2	A	900	COT	C2A-N3A	2.39	1.38	1.33
2	A	900	COT	CGP-CDP	2.34	1.56	1.52
2	A	900	COT	OCP-CBP	2.30	1.46	1.42
2	A	900	COT	PC-O3B	-2.16	1.55	1.59
2	A	900	COT	C8A-N9A	2.08	1.41	1.37
2	A	900	COT	C6A-N1A	2.01	1.41	1.35

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	COT	C1P-C2P-N3P	6.99	127.01	112.41
2	A	900	COT	C1P-S-C6T	5.48	110.75	101.72
2	A	900	COT	O4B-C1B-N9A	3.02	113.89	108.09
2	A	900	COT	O11-CGP-CDP	3.00	115.36	110.55
2	A	900	COT	C3B-C2B-C1B	2.88	106.23	99.89
2	A	900	COT	CEP-CDP-CBP	2.57	113.15	108.77

There are no chirality outliers.

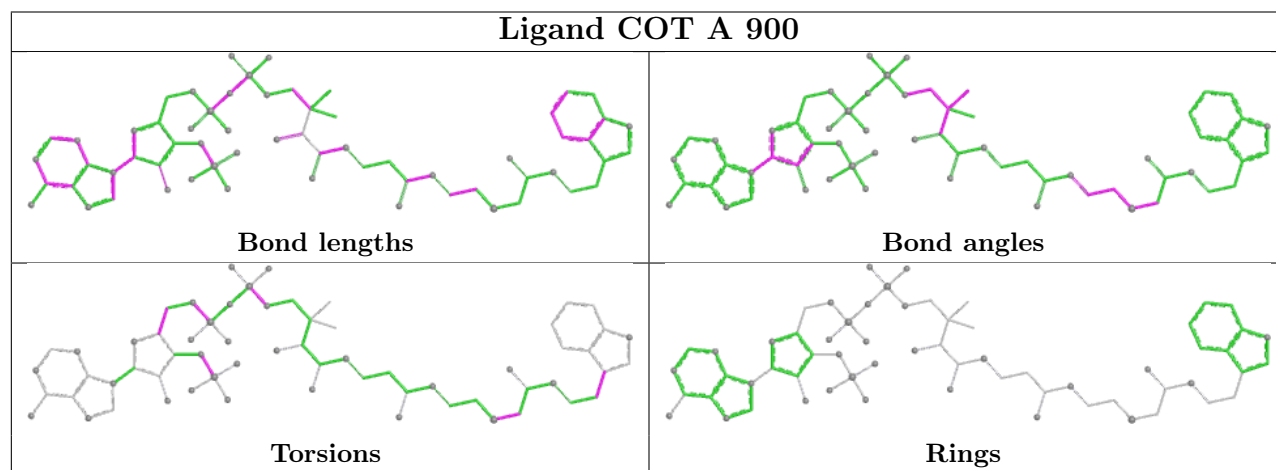
All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	900	COT	CGP-O11-PB-O13
2	A	900	COT	C3B-C4B-C5B-O5B
2	A	900	COT	O4B-C4B-C5B-O5B
2	A	900	COT	C4T-C6T-S-C1P
2	A	900	COT	CGP-O11-PB-O14
2	A	900	COT	C5B-O5B-PA-O15
2	A	900	COT	C3B-O3B-PC-O18
2	A	900	COT	C3B-O3B-PC-O19
2	A	900	COT	C3B-O3B-PC-O17
2	A	900	COT	C2T-C1T-CGT-C11

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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.