



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 20, 2026 – 11:32 AM UTC

PDB ID : 8X4S / pdb_00008x4s
Title : The L-tryptophan specific decarboxylase PsiD covalent bonding with tryptamine
Authors : Meng, C.Y.; Wen, Y.; Guo, W.T.; Wu, B.X.
Deposited on : 2023-11-15
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

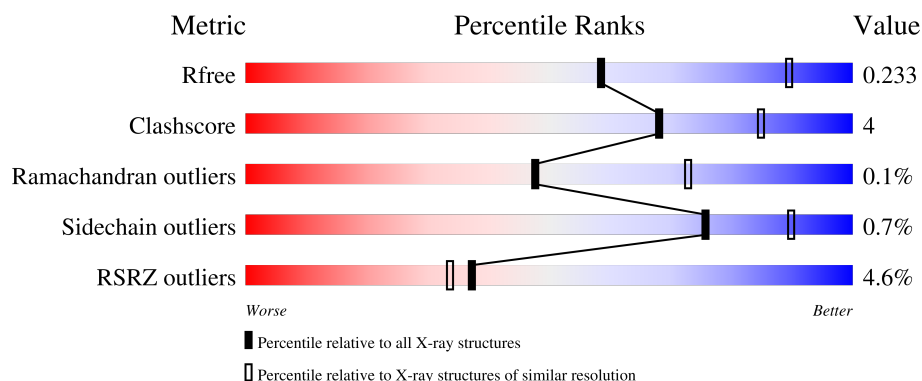
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	402	 2% 79% 7% 13%
1	C	402	 % 78% 7% 14%
1	E	402	 4% 78% 8% 14%
1	G	402	 3% 79% 6% 14%
1	I	402	 2% 79% 7% 14%

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Mol	Chain	Length	Quality of chain
1	K	402	13% 77% 9% 14%
2	B	36	92% 6%
2	D	36	92% 6%
2	F	36	92% 6%
2	H	36	92% 6%
2	J	36	92% 6%
2	L	36	11% 86% 8% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18594 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.

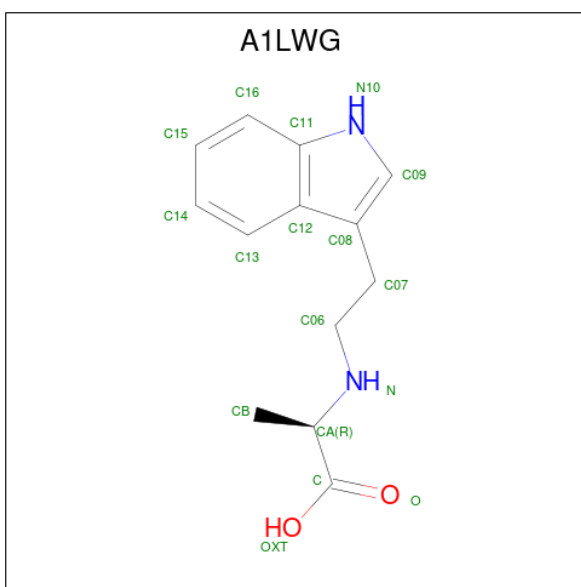
- Molecule 1 is a protein called L-tryptophan decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	I	346	Total	C	N	O	S	0	0	0
			2747	1766	461	505	15			
1	G	346	Total	C	N	O	S	0	0	0
			2754	1768	460	511	15			
1	A	350	Total	C	N	O	S	0	0	0
			2792	1792	469	516	15			
1	C	346	Total	C	N	O	S	0	0	0
			2763	1776	464	508	15			
1	E	346	Total	C	N	O	S	0	0	0
			2753	1769	459	510	15			
1	K	347	Total	C	N	O	S	0	0	0
			2762	1773	463	511	15			

- Molecule 2 is a protein called L-tryptophan decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	34	Total	C	N	O	S	0	0	0
			240	155	42	42	1			
2	H	34	Total	C	N	O	S	0	0	0
			244	158	43	42	1			
2	B	34	Total	C	N	O	S	0	0	0
			240	155	42	42	1			
2	D	34	Total	C	N	O	S	0	0	0
			244	157	42	44	1			
2	F	34	Total	C	N	O	S	0	0	0
			248	159	42	46	1			
2	L	34	Total	C	N	O	S	0	0	0
			244	157	42	44	1			

- Molecule 3 is (2 {R})-2-[2-(1 {H}-indol-3-yl)ethylamino]propanoic acid (CCD ID: A1LWG) (formula: C₁₃H₁₆N₂O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	J	1	Total	C	N	O	0	0
			16	13	2	1		
3	H	1	Total	C	N	O	0	0
			16	13	2	1		
3	B	1	Total	C	N	O	0	0
			16	13	2	1		
3	D	1	Total	C	N	O	0	0
			16	13	2	1		
3	F	1	Total	C	N	O	0	0
			16	13	2	1		
3	L	1	Total	C	N	O	0	0
			16	13	2	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	64	Total	O	0	0
			64	64		
4	J	2	Total	O	0	0
			2	2		
4	G	99	Total	O	0	0
			99	99		
4	H	8	Total	O	0	0
			8	8		
4	A	90	Total	O	0	0
			90	90		
4	B	9	Total	O	0	0
			9	9		

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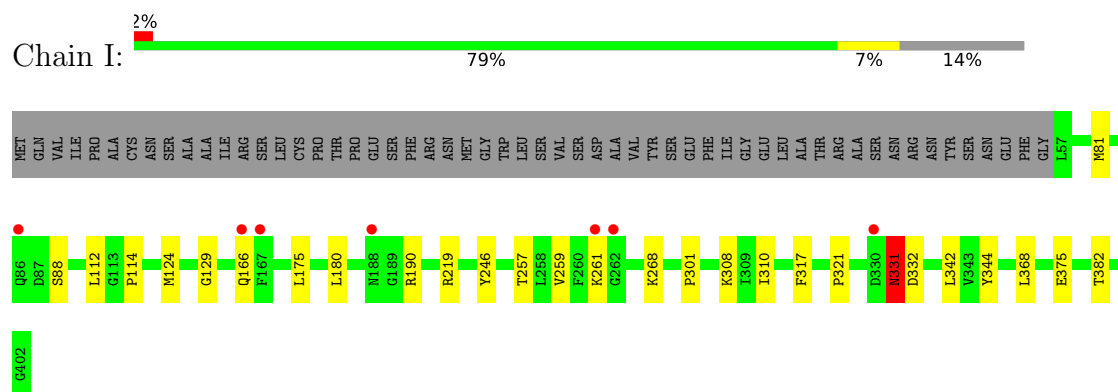
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	74	Total 74	O 74	0	0
4	D	9	Total 9	O 9	0	0
4	E	73	Total 73	O 73	0	0
4	F	4	Total 4	O 4	0	0
4	K	34	Total 34	O 34	0	0
4	L	1	Total 1	O 1	0	0

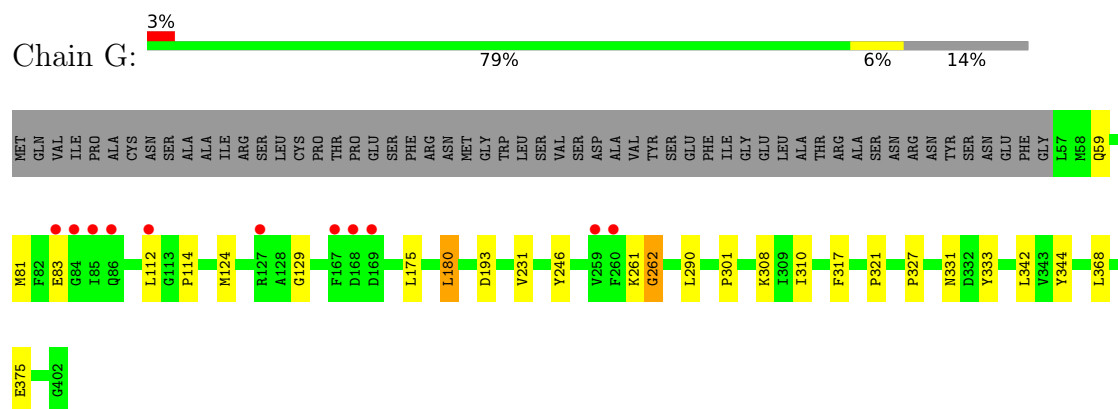
3 Residue-property plots [i](#)

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.

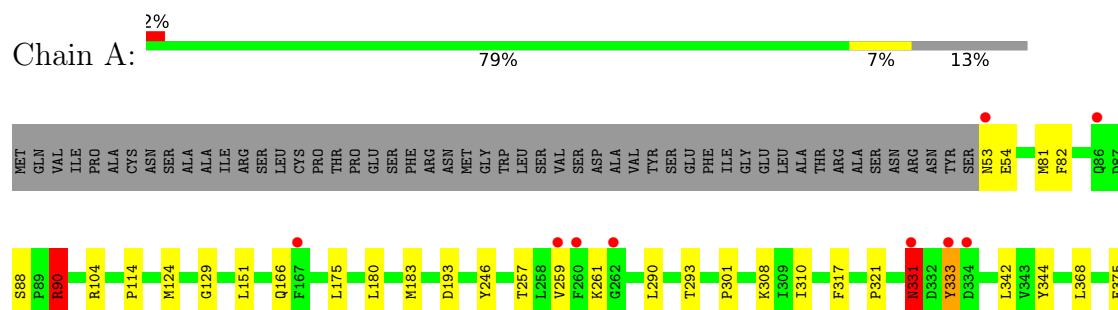
- Molecule 1: L-tryptophan decarboxylase



- Molecule 1: L-tryptophan decarboxylase

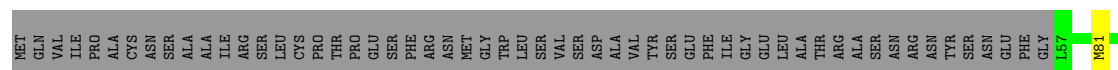
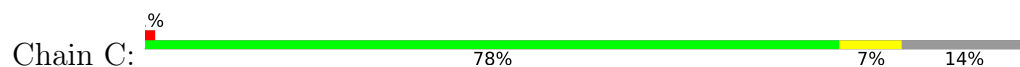


- Molecule 1: L-tryptophan decarboxylase

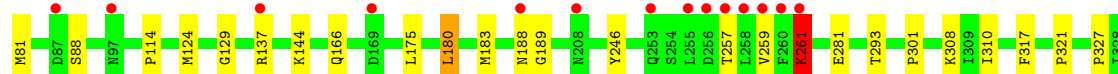
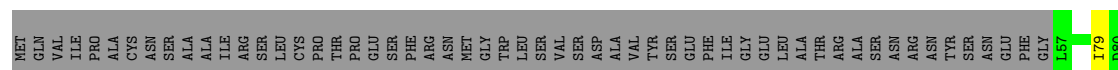
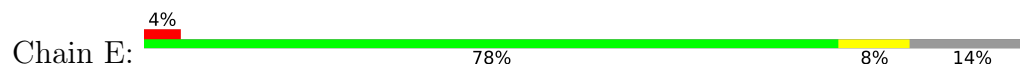




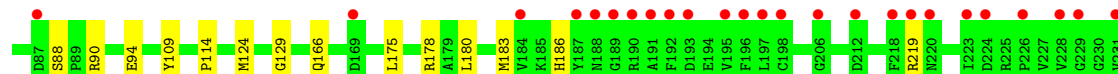
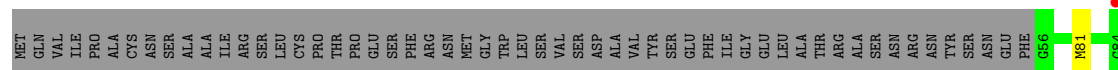
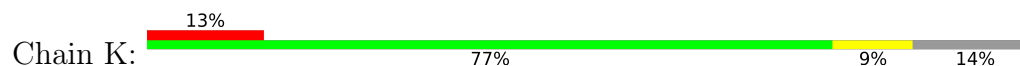
- Molecule 1: L-tryptophan decarboxylase



- Molecule 1: L-tryptophan decarboxylase

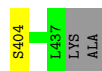


- Molecule 1: L-tryptophan decarboxylase



- Molecule 2: L-tryptophan decarboxylase

Chain J:  92% • 6%



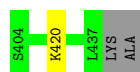
- Molecule 2: L-tryptophan decarboxylase

Chain H:  92% • 6%



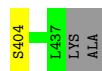
- Molecule 2: L-tryptophan decarboxylase

Chain B:  92% • 6%



- Molecule 2: L-tryptophan decarboxylase

Chain D:  92% • 6%



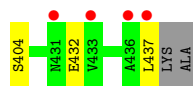
- Molecule 2: L-tryptophan decarboxylase

Chain F:  92% • 6%



- Molecule 2: L-tryptophan decarboxylase

Chain L:  11% 86% 8% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.68Å 122.50Å 128.72Å 90.00° 99.77° 90.00°	Depositor
Resolution (Å)	29.77 – 2.70 29.77 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.77-2.70) 99.6 (29.77-2.70)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.68Å)	Xtriage
Refinement program	PHENIX 1.21rc1_5156	Depositor
R, R_{free}	0.198 , 0.233 0.200 , 0.233	Depositor DCC
R_{free} test set	3487 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	40.6	Xtriage
Anisotropy	0.372	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18594	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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: A1LWG

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	0.46	2/2870 (0.1%)	0.62	2/3896 (0.1%)
1	C	0.40	0/2840	0.60	2/3854 (0.1%)
1	E	0.43	1/2830 (0.0%)	0.62	1/3844 (0.0%)
1	G	0.42	0/2831	0.62	1/3846 (0.0%)
1	I	0.43	1/2824 (0.0%)	0.63	2/3836 (0.1%)
1	K	0.45	1/2839 (0.0%)	0.62	2/3855 (0.1%)
2	B	0.40	0/242	0.75	0/329
2	D	0.40	0/246	0.79	1/334 (0.3%)
2	F	0.44	0/250	1.13	1/339 (0.3%)
2	H	0.41	0/246	0.96	1/333 (0.3%)
2	J	0.40	0/242	0.96	1/329 (0.3%)
2	L	0.40	0/246	1.33	1/334 (0.3%)
All	All	0.43	5/18506 (0.0%)	0.66	15/25129 (0.1%)

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	A	0	2
1	I	0	1
1	K	0	2
All	All	0	5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	330	ASP	CB-CG	6.99	1.69	1.52
1	A	333	TYR	C-N	6.10	1.47	1.33
1	I	331	ASN	C-N	6.03	1.42	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	331	ASN	C-N	-5.64	1.25	1.33
1	E	261	LYS	C-N	5.03	1.37	1.33

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	404	SER	O-C-N	-21.52	88.56	123.00
2	F	404	SER	O-C-N	-17.35	95.23	123.00
2	J	404	SER	O-C-N	-13.29	101.73	123.00
2	H	404	SER	O-C-N	-13.22	101.85	123.00
1	G	83	GLU	CA-CB-CG	-9.98	94.15	114.10

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	331	ASN	Mainchain
1	A	90	ARG	Sidechain
1	I	331	ASN	Mainchain
1	K	178	ARG	Sidechain
1	K	331	ASN	Peptide

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	2792	0	2643	19	2
1	C	2763	0	2631	20	0
1	E	2753	0	2602	29	2
1	G	2754	0	2601	18	0
1	I	2747	0	2602	19	1
1	K	2762	0	2617	33	0
2	B	240	0	241	0	0
2	D	244	0	245	0	0
2	F	248	0	249	0	0
2	H	244	0	252	0	0
2	J	240	0	241	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	244	0	245	2	0
3	B	16	0	0	1	0
3	D	16	0	0	1	0
3	F	16	0	0	0	0
3	H	16	0	0	1	0
3	J	16	0	0	0	0
3	L	16	0	0	1	0
4	A	90	0	0	1	1
4	B	9	0	0	0	0
4	C	74	0	0	2	0
4	D	9	0	0	0	0
4	E	73	0	0	6	0
4	F	4	0	0	0	0
4	G	99	0	0	3	0
4	H	8	0	0	0	0
4	I	64	0	0	1	0
4	J	2	0	0	0	0
4	K	34	0	0	3	0
4	L	1	0	0	0	0
All	All	18594	0	17169	128	3

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 4.

The worst 5 of 128 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:K:331:ASN:OD1	1:K:333:TYR:N	2.13	0.81
1:K:186:HIS:O	4:K:501:HOH:O	1.99	0.79
1:E:166:GLN:HB3	1:K:232:ASN:CG	2.07	0.79
1:C:388:HIS:O	4:C:501:HOH:O	2.02	0.78
1:A:53:ASN:O	1:A:104:ARG:NH1	2.22	0.72

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:PHE:O	1:E:188:ASN:ND2[2_656]	2.03	0.17
1:I:268:LYS:NZ	4:A:588:HOH:O[2_645]	2.04	0.16
1:A:90:ARG:O	1:E:188:ASN:OD1[2_656]	2.07	0.13

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	348/402 (87%)	335 (96%)	13 (4%)	0	100	100
1	C	344/402 (86%)	331 (96%)	12 (4%)	1 (0%)	36	60
1	E	344/402 (86%)	331 (96%)	13 (4%)	0	100	100
1	G	344/402 (86%)	331 (96%)	12 (4%)	1 (0%)	36	60
1	I	344/402 (86%)	330 (96%)	14 (4%)	0	100	100
1	K	345/402 (86%)	332 (96%)	13 (4%)	0	100	100
2	B	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
2	D	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
2	F	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
2	H	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
2	J	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
2	L	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
All	All	2261/2628 (86%)	2176 (96%)	83 (4%)	2 (0%)	48	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	262	GLY
1	C	262	GLY

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	299/350 (85%)	296 (99%)	3 (1%)	68	86
1	C	296/350 (85%)	295 (100%)	1 (0%)	86	94
1	E	294/350 (84%)	291 (99%)	3 (1%)	68	86
1	G	295/350 (84%)	294 (100%)	1 (0%)	86	94
1	I	293/350 (84%)	292 (100%)	1 (0%)	86	94
1	K	296/350 (85%)	293 (99%)	3 (1%)	68	86
2	B	23/29 (79%)	22 (96%)	1 (4%)	26	54
2	D	24/29 (83%)	24 (100%)	0	100	100
2	F	25/29 (86%)	25 (100%)	0	100	100
2	H	24/29 (83%)	24 (100%)	0	100	100
2	J	23/29 (79%)	23 (100%)	0	100	100
2	L	24/29 (83%)	24 (100%)	0	100	100
All	All	1916/2274 (84%)	1903 (99%)	13 (1%)	76	90

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	88	SER
1	E	180	LEU
1	K	330	ASP
1	K	88	SER
1	K	292	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 39 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	166	GLN
1	K	166	GLN
1	E	186	HIS
1	E	319	GLN
1	K	253	GLN

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](#)

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	A1LWG	H	501	2	16,17,18	2.26	6 (37%)	20,22,24	1.35	1 (5%)
3	A1LWG	B	501	2	16,17,18	1.89	6 (37%)	20,22,24	1.57	2 (10%)
3	A1LWG	F	501	2	16,17,18	1.94	5 (31%)	20,22,24	1.19	2 (10%)
3	A1LWG	J	501	2	16,17,18	2.06	4 (25%)	20,22,24	1.00	0
3	A1LWG	L	501	2	16,17,18	1.96	6 (37%)	20,22,24	1.01	1 (5%)
3	A1LWG	D	501	2	16,17,18	2.16	9 (56%)	20,22,24	1.34	3 (15%)

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	A1LWG	H	501	2	-	1/7/8/10	0/2/2/2
3	A1LWG	B	501	2	-	1/7/8/10	0/2/2/2
3	A1LWG	F	501	2	-	2/7/8/10	0/2/2/2
3	A1LWG	J	501	2	-	1/7/8/10	0/2/2/2
3	A1LWG	L	501	2	-	1/7/8/10	0/2/2/2
3	A1LWG	D	501	2	-	1/7/8/10	0/2/2/2

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	501	A1LWG	C07-C08	4.84	1.59	1.50
3	F	501	A1LWG	CA-N	4.20	1.54	1.47
3	J	501	A1LWG	C07-C08	4.14	1.58	1.50
3	J	501	A1LWG	CA-N	4.01	1.54	1.47
3	H	501	A1LWG	CA-N	3.67	1.53	1.47

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	A1LWG	CB-CA-N	4.77	117.51	108.80
3	H	501	A1LWG	C06-N-CA	3.88	119.46	113.42
3	F	501	A1LWG	CB-CA-N	3.79	115.72	108.80
3	D	501	A1LWG	C06-N-CA	3.23	118.45	113.42
3	D	501	A1LWG	O-C-CA	-3.00	114.61	123.94

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	J	501	A1LWG	CB-CA-N-C06
3	H	501	A1LWG	CB-CA-N-C06
3	B	501	A1LWG	N-C06-C07-C08
3	F	501	A1LWG	CB-CA-N-C06
3	F	501	A1LWG	N-C06-C07-C08

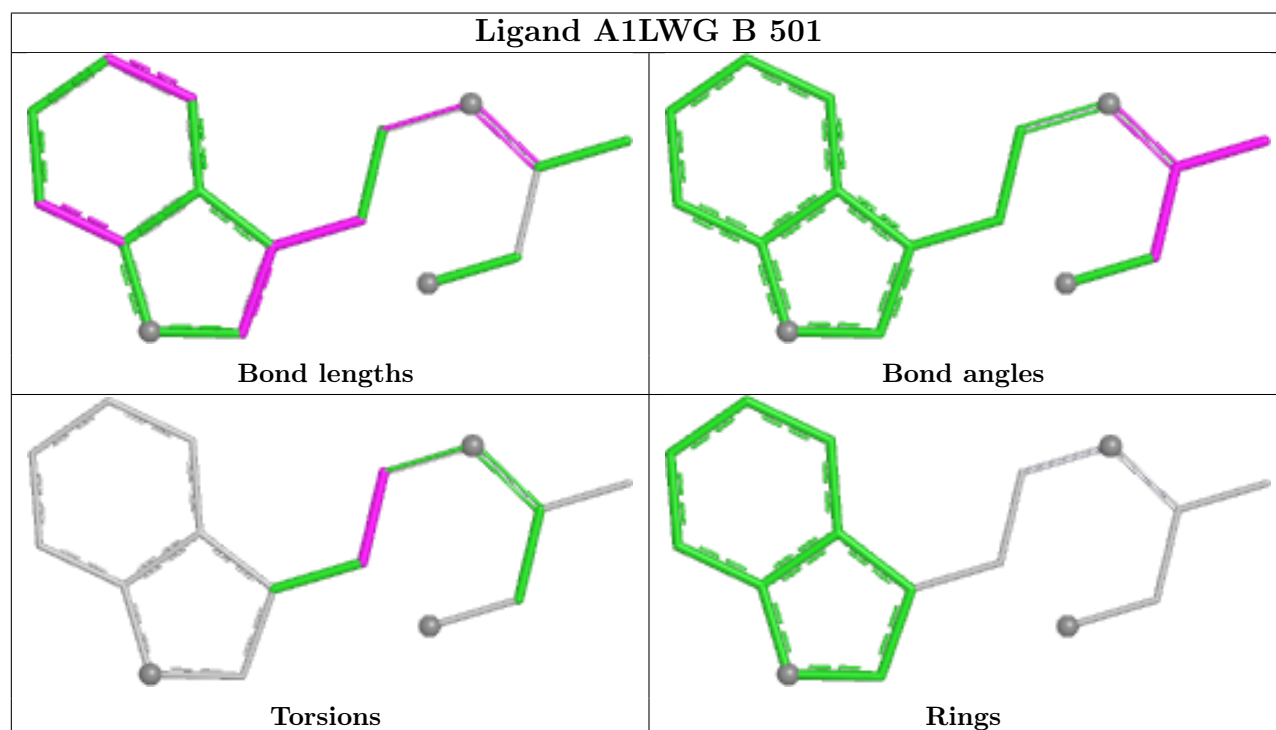
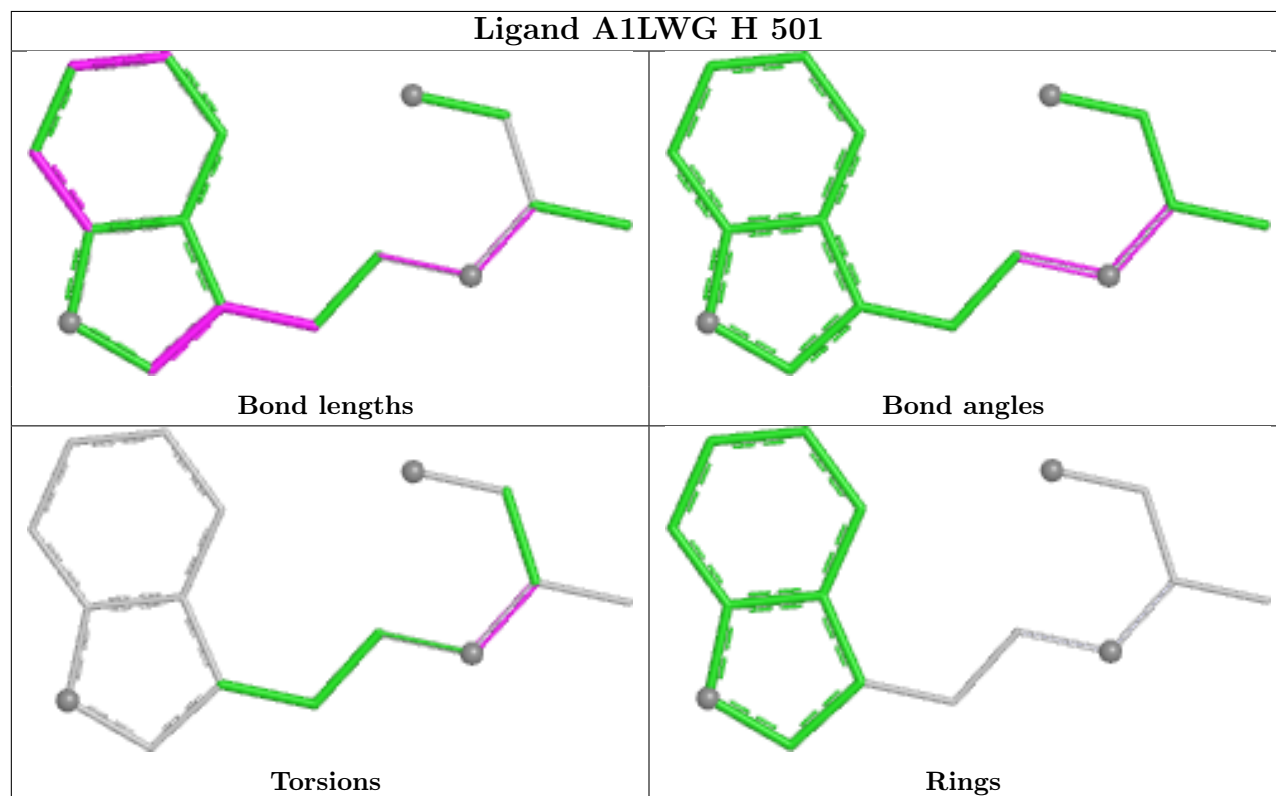
There are no ring outliers.

4 monomers are involved in 4 short contacts:

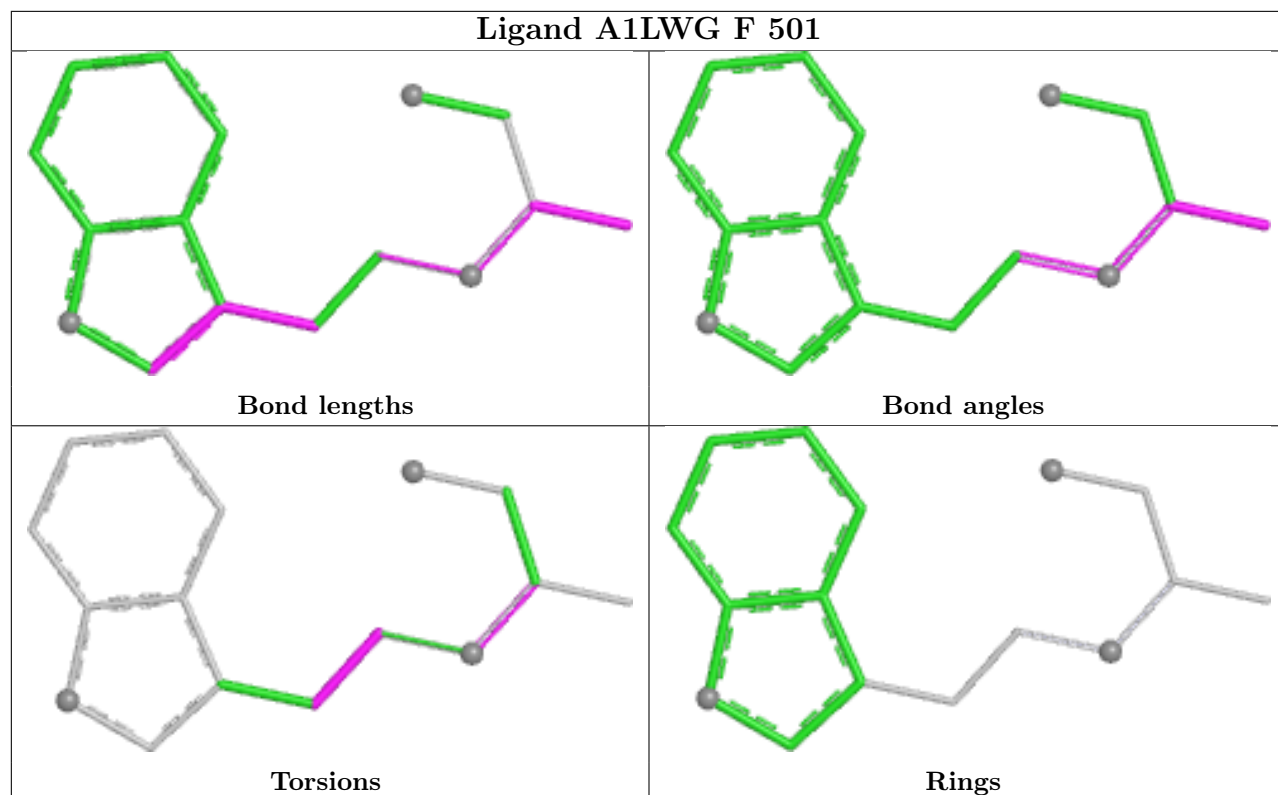
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	501	A1LWG	1	0
3	B	501	A1LWG	1	0
3	L	501	A1LWG	1	0
3	D	501	A1LWG	1	0

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

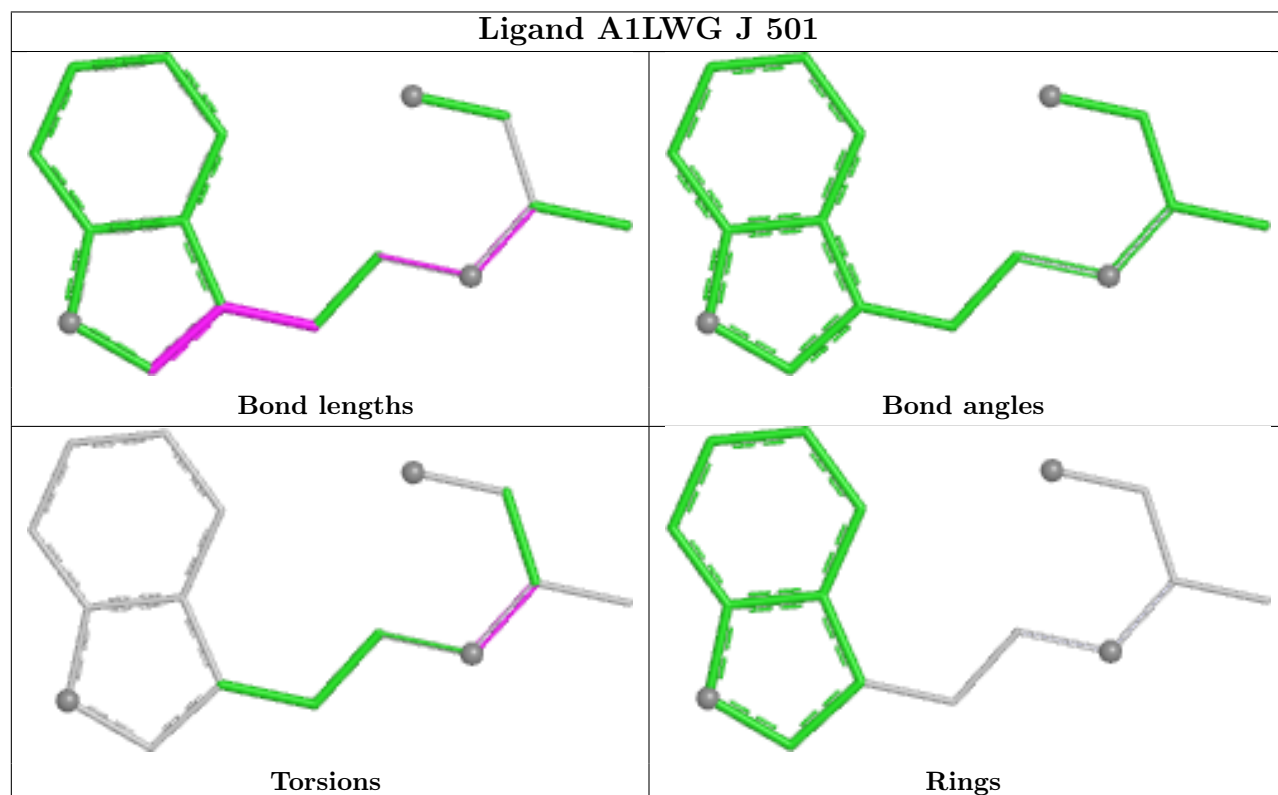
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

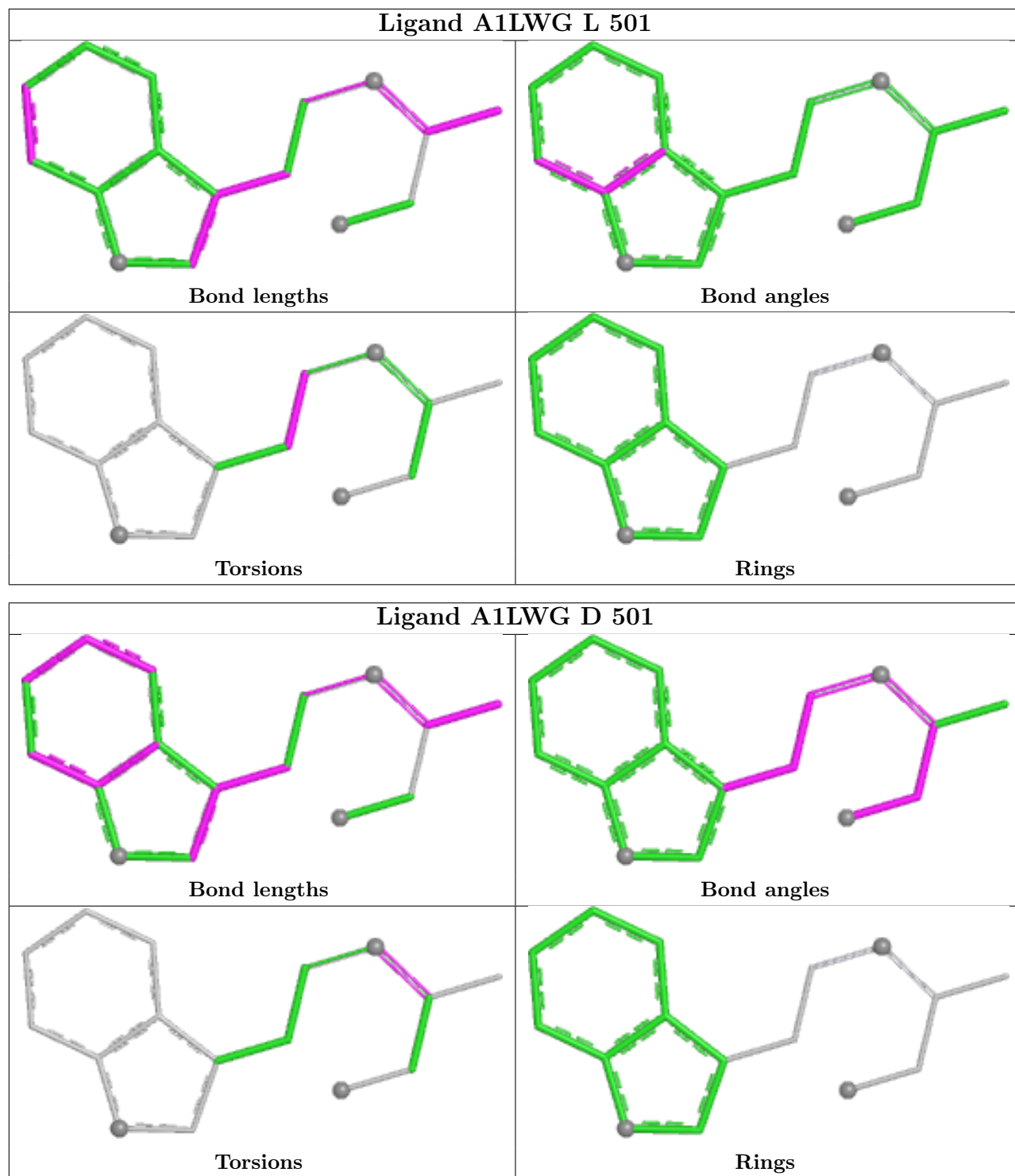


Ligand A1LWG F 501



Ligand A1LWG J 501





5.7 Other polymers [i](#)

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 [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	350/402 (87%)	-0.10	9 (2%) 57 54	20, 35, 70, 102	0
1	C	346/402 (86%)	-0.09	5 (1%) 73 72	20, 40, 71, 111	0
1	E	346/402 (86%)	0.12	15 (4%) 40 36	24, 41, 76, 113	0
1	G	346/402 (86%)	-0.04	11 (3%) 50 46	19, 34, 69, 108	0
1	I	346/402 (86%)	0.13	7 (2%) 65 62	23, 43, 75, 97	0
1	K	347/402 (86%)	0.81	54 (15%) 5 4	25, 57, 97, 132	0
2	B	34/36 (94%)	-0.35	0 100 100	20, 37, 62, 78	0
2	D	34/36 (94%)	-0.31	0 100 100	23, 36, 61, 73	0
2	F	34/36 (94%)	-0.15	0 100 100	23, 37, 55, 69	0
2	H	34/36 (94%)	-0.46	0 100 100	23, 33, 58, 70	0
2	J	34/36 (94%)	-0.11	0 100 100	26, 39, 68, 75	0
2	L	34/36 (94%)	1.24	4 (11%) 9 7	37, 69, 94, 111	0
All	All	2285/2628 (86%)	0.12	105 (4%) 37 34	19, 40, 83, 132	0

The worst 5 of 105 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	259	VAL	5.5
1	K	256	ASP	5.5
1	G	167	PHE	4.9
1	E	188	ASN	4.7
1	K	257	THR	4.2

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](#)

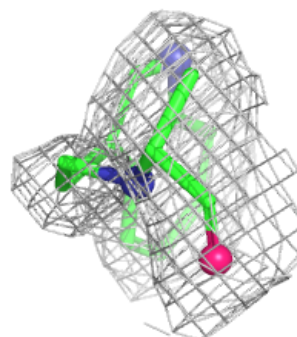
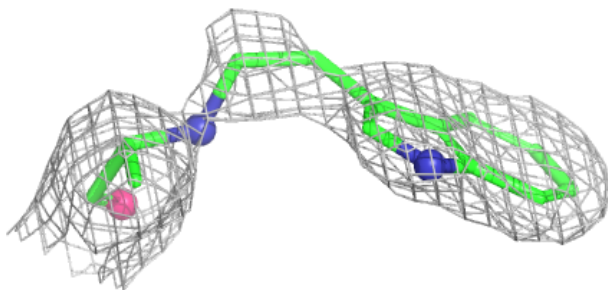
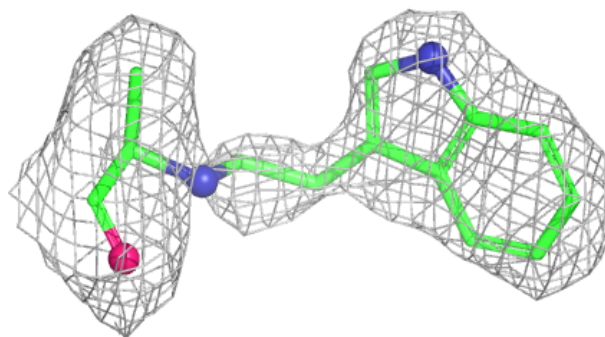
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	A1LWG	D	501	16/17	0.93	0.14	17,48,123,128	0
3	A1LWG	L	501	16/17	0.94	0.14	23,59,127,130	0
3	A1LWG	B	501	16/17	0.95	0.14	16,60,122,124	0
3	A1LWG	J	501	16/17	0.95	0.10	17,54,92,96	0
3	A1LWG	H	501	16/17	0.95	0.10	16,39,89,90	0
3	A1LWG	F	501	16/17	0.96	0.10	24,38,126,146	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

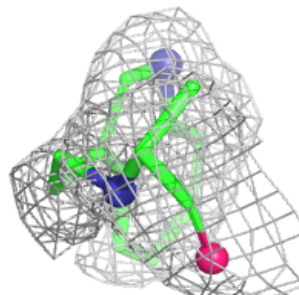
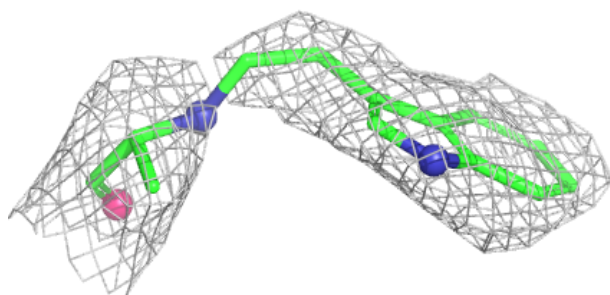
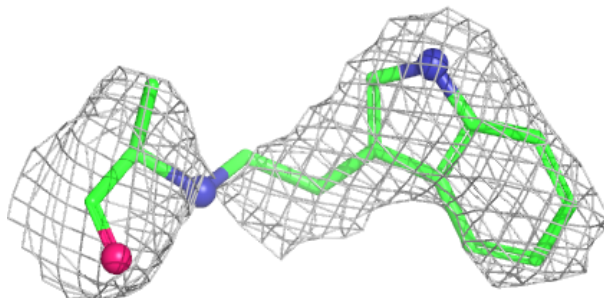
Electron density around A1LWG D 501:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

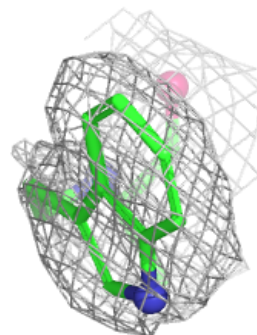
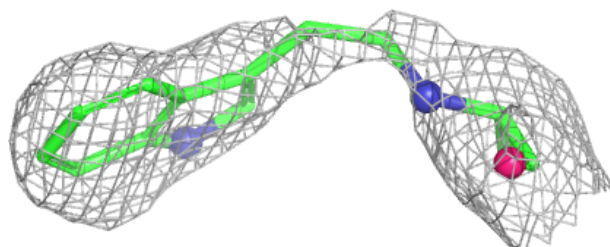
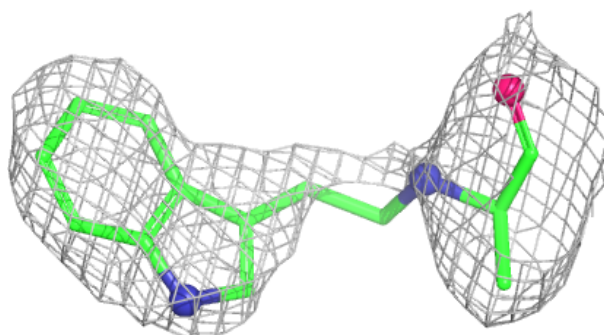


Electron density around A1LWG L 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

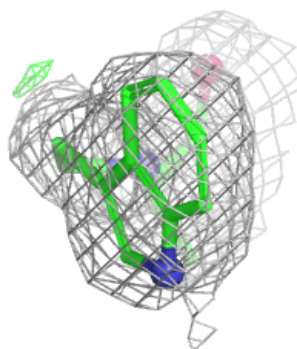
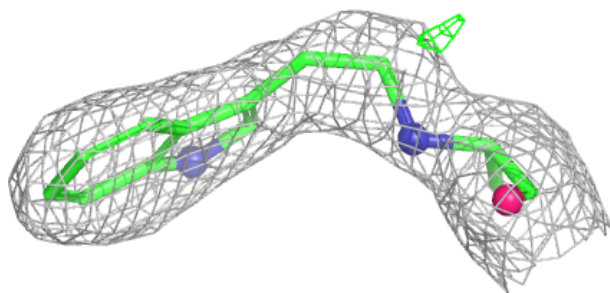
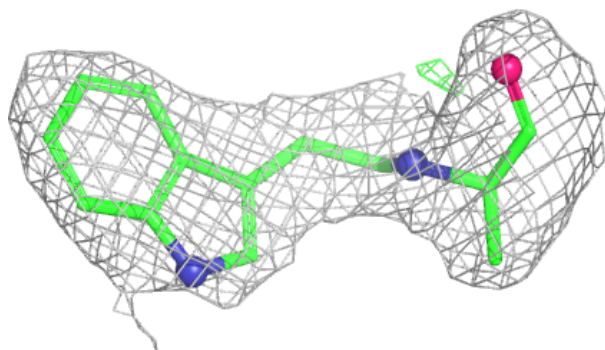
**Electron density around A1LWG B 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

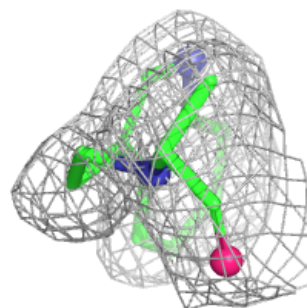
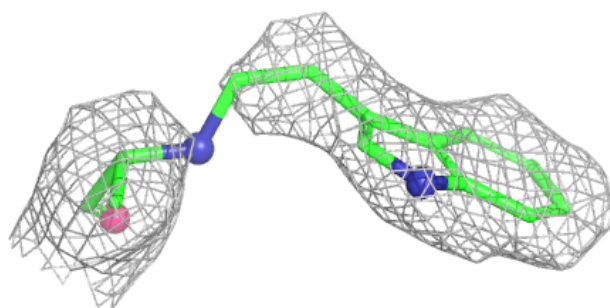
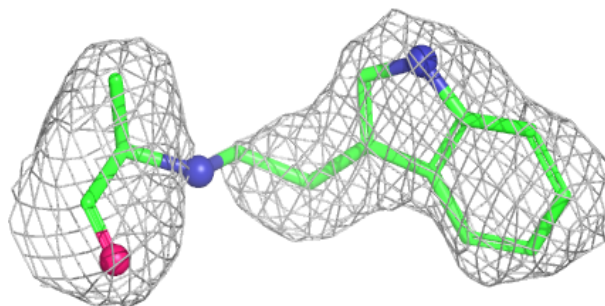


Electron density around A1LWG J 501:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)

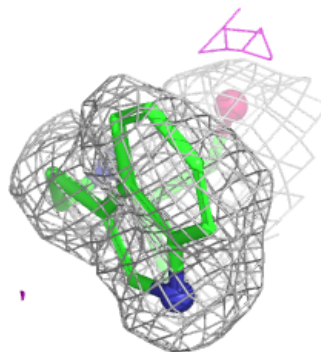
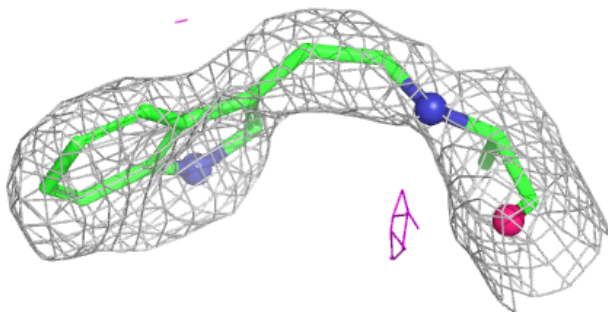
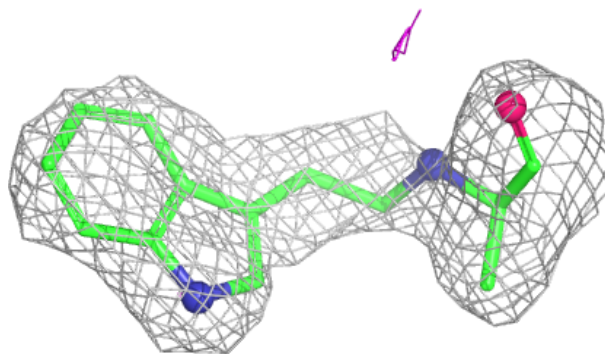
**Electron density around A1LWG H 501:**

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around A1LWG F 501:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.