

wwPDB X-ray Structure Validation Summary Report (i)

Oct 11, 2023 – 09:50 PM EDT

PDB ID	:	7LK1
Title	:	Ornithine Aminotransferase (OAT) with its potent inhibitor - (S)-3-amino-4,
		4-difluorocyclopent-1-enecarboxylic acid (SS-1-148) - 1 Hour Soaking
Authors	:	Butrin, A.; Shen, S.; Liu, D.; Silverman, R.
Deposited on	:	2021-02-01
Resolution	:	1.79 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 (i) 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 (1)) were used in the production of this report:

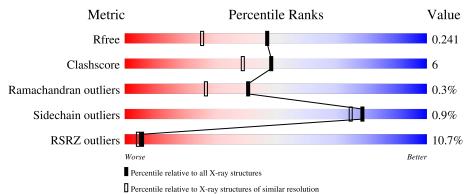
MolProbity	:	4.02b-467 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)		
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.79 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	$\begin{array}{c} \textbf{Whole archive} \\ \textbf{(\#Entries)} \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R _{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	А	404	82%	17%
1	В	404	85%	14% •
1	С	404	5% 91%	7% •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:



Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	Y37	А	501[B]	Х	-	-	-
2	Y37	С	501[A]	Х	-	-	-
2	Y37	С	501[B]	Х	-	-	-



2 Entry composition (i)

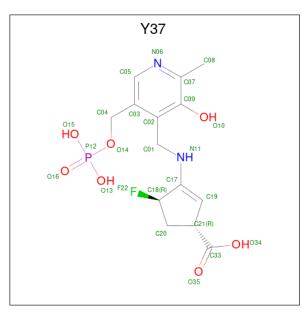
There are 3 unique types of molecules in this entry. The entry contains 10380 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Δ	404	Total	С	Ν	0	\mathbf{S}	0	1	0
	А	404	3166	2034	535	585	12	0		0
1	D	403	Total	С	Ν	0	S	0	0	0
	D	405	3157	2028	532	585	12	0	0	0
1	C	404	Total	С	Ν	0	S	0	1	0
	U	404	3166	2034	535	585	12	U		0

• Molecule 1 is a protein called Ornithine aminotransferase, mitochondrial.

• Molecule 2 is (1R,4R)-4-fluoro-3-[({3-hydroxy-2-methyl-5-[(phosphonooxy)methyl]pyridi n-4-yl}methyl)amino]cyclopent-2-ene-1-carboxylic acid (three-letter code: Y37) (formula: C₁₄H₁₈FN₂O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
9	۸	1	Total	С	F	Ν	Ο	Р	0	1
	A	1	50	28	2	4	14	2	0	T
9	В	1	Total	С	F	Ν	Ο	Р	0	0
	D	1	25	14	1	2	7	1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf			
2	С	1	Total 50	C 28	F 2	N 4	0 14	Р 2	0	1

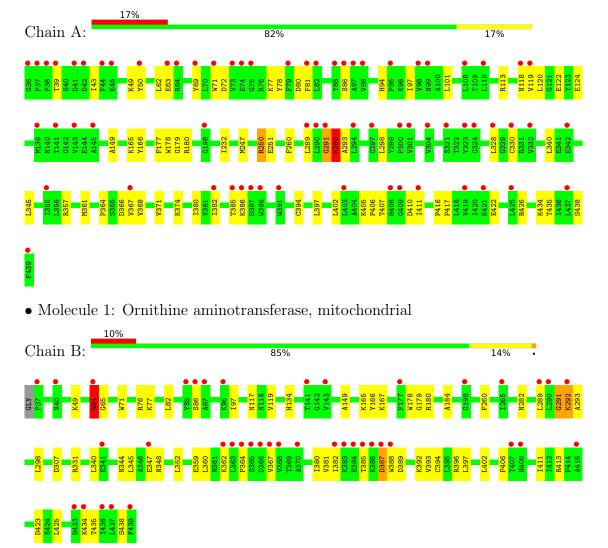
• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	243	Total O 243 243	0	0
3	В	224	Total O 224 224	0	0
3	С	299	Total O 299 299	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: Ornithine aminotransferase, mitochondrial

• Molecule 1: Ornithine aminotransferase, mitochondrial





V367 C36 V367 L52 L376 L62 L385 K66 T385 K66 T385 K66 T385 K107 H408 L110 T411 V119 T411 V119 T413 H139 K434 H139 K434 H139 K434 N139 K165 V139 K165 N139 K165 N139 K165 N139 K165 N139 K165 N139 K165 N139 N282 N282 N282 N282 N282 N358 N361 N361 N361 N361



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	115.65Å 115.65 Å 186.69 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	-
Resolution (Å)	36.32 - 1.79	Depositor
	36.32 $ 1.79$	EDS
% Data completeness	100.0 (36.32 - 1.79)	Depositor
(in resolution range)	100.0 (36.32 - 1.79)	EDS
R _{merge}	0.08	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.22 (at 1.79 Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
D D	0.216 , 0.241	Depositor
R, R_{free}	0.216 , 0.241	DCC
R_{free} test set	6743 reflections $(4.95%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	34.6	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34 , 42.2	EDS
L-test for twinning ²	$< L > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10380	wwPDB-VP
Average B, all atoms $(Å^2)$	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.70% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: $\mathbf{Y37}$

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	Chain Bond length		Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.37	0/3240	0.63	7/4399~(0.2%)	
1	В	0.34	0/3231	0.84	4/4387~(0.1%)	
1	С	0.36	0/3240	0.59	2/4399~(0.0%)	
All	All	0.36	0/9711	0.69	13/13185~(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	А	0	4
1	В	0	3
1	С	0	5
All	All	0	12

There are no bond length outliers.

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\mathbf{Ideal}(^{o})$
1	В	291	GLY	O-C-N	-41.07	56.99	122.70
1	А	292[A]	LYS	O-C-N	-11.76	103.89	122.70
1	А	292[B]	LYS	O-C-N	-11.76	103.89	122.70
1	С	291	GLY	O-C-N	-10.39	106.07	122.70
1	В	64	ARG	NE-CZ-NH2	-9.37	115.61	120.30

There are no chirality outliers.

5 of 12 planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	А	291	GLY	Mainchain
1	А	292[A]	LYS	Mainchain
1	А	292[B]	LYS	Mainchain
1	В	291	GLY	Peptide,Mainchain
1	В	292	LYS	Mainchain

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	А	3166	0	3169	48	0
1	В	3157	0	3160	46	0
1	С	3166	0	3170	22	0
2	А	50	0	0	3	0
2	В	25	0	0	2	0
2	С	50	0	0	2	0
3	А	243	0	0	5	0
3	В	224	0	0	5	0
3	С	299	0	0	3	0
All	All	10380	0	9499	115	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.

The worst 5 of 115 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:ARG:HE	1:B:65:GLY:H	1.17	0.90
1:B:64:ARG:HE	1:B:65:GLY:N	1.74	0.85
1:B:64:ARG:HB3	1:B:71:TRP:HB2	1.66	0.77
1:C:180:ARG:NH1	3:C:602:HOH:O	2.17	0.77
1:C:104:GLN:HA	1:C:107:LYS:HG2	1.67	0.74

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	Perce	ntiles
1	А	403/404~(100%)	380~(94%)	21 (5%)	2(0%)	29	15
1	В	401/404~(99%)	380~(95%)	21 (5%)	0	100	100
1	С	403/404~(100%)	381 (94%)	18 (4%)	4 (1%)	15	5
All	All	1207/1212~(100%)	1141 (94%)	60~(5%)	6 (0%)	41	15

 $5~{\rm of}~6$ Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	292[A]	LYS
1	А	292[B]	LYS
1	С	166	TYR
1	С	292[A]	LYS
1	С	292[B]	LYS

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	А	337/337~(100%)	335~(99%)	2(1%)	86	84	
1	В	337/337~(100%)	332~(98%)	5(2%)	65	56	
1	С	337/337~(100%)	335~(99%)	2(1%)	86	84	
All	All	1011/1011 (100%)	1002 (99%)	9 (1%)	78	75	

5 of 9 residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	С	260	PHE
1	С	387	ASP
1	В	82	LEU
1	В	260	PHE
1	В	387	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	С	344	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

5 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).

Mal	Mol Type Chain		Res	Link	Bond lengths			Bond angles		
10101	Type	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
2	Y37	С	501[B]	1	23,26,26	1.99	1 (4%)	$26,\!38,\!38$	1.75	2 (7%)
2	Y37	С	501[A]	1	23,26,26	1.97	1 (4%)	26,38,38	2.02	<mark>6 (23%)</mark>
2	Y37	В	501	1	23,26,26	1.86	1 (4%)	26,38,38	1.80	2 (7%)



Mal	Mol Type Chain Res		Link	Bond lengths			Bond angles			
10101	Type	Chain	Res	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	Y37	А	501[B]	1	23,26,26	2.10	1 (4%)	$26,\!38,\!38$	1.66	3 (11%)
2	Y37	А	501[A]	1	23,26,26	1.94	2 (8%)	$26,\!38,\!38$	1.92	3 (11%)

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	Y37	С	501[B]	1	1/1/5/7	0/14/27/27	0/2/2/2
2	Y37	С	501[A]	1	2/2/5/7	0/14/27/27	0/2/2/2
2	Y37	В	501	1	-	4/14/27/27	0/2/2/2
2	Y37	А	501[B]	1	2/2/5/7	3/14/27/27	0/2/2/2
2	Y37	А	501[A]	1	-	1/14/27/27	0/2/2/2

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
2	А	501[B]	Y37	C17-N11	8.55	1.45	1.34
2	С	501[B]	Y37	C17-N11	8.10	1.45	1.34
2	С	501[A]	Y37	C17-N11	7.93	1.45	1.34
2	А	501[A]	Y37	C17-N11	7.64	1.44	1.34
2	В	501	Y37	C17-N11	6.97	1.43	1.34

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	С	501[A]	Y37	F22-C18-C20	7.81	120.86	109.53
2	В	501	Y37	F22-C18-C20	7.72	120.72	109.53
2	А	501[A]	Y37	F22-C18-C20	7.55	120.49	109.53
2	С	501[B]	Y37	F22-C18-C20	7.27	120.07	109.53
2	А	501[B]	Y37	F22-C18-C20	6.82	119.42	109.53

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	А	501[B]	Y37	C21
2	А	501[B]	Y37	C18
2	С	501[A]	Y37	C21
2	С	501[A]	Y37	C18
2	С	501[B]	Y37	C21



Mol	Chain	Res	Type	Atoms
2	А	501[A]	Y37	C18-C17-N11-C01
2	В	501	Y37	C18-C17-N11-C01
2	А	501[B]	Y37	N11-C01-C02-C03
2	В	501	Y37	N11-C01-C02-C03
2	А	501[B]	Y37	N11-C01-C02-C09

5 of 8 torsion outliers are listed below:

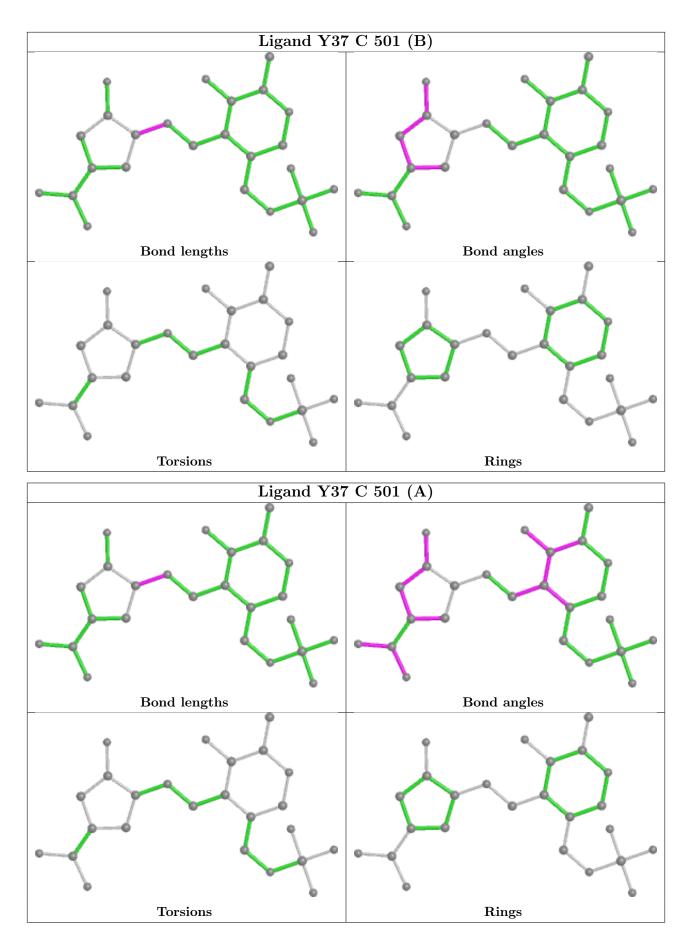
There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	501[B]	Y37	2	0
2	В	501	Y37	2	0
2	А	501[B]	Y37	1	0
2	А	501[A]	Y37	2	0

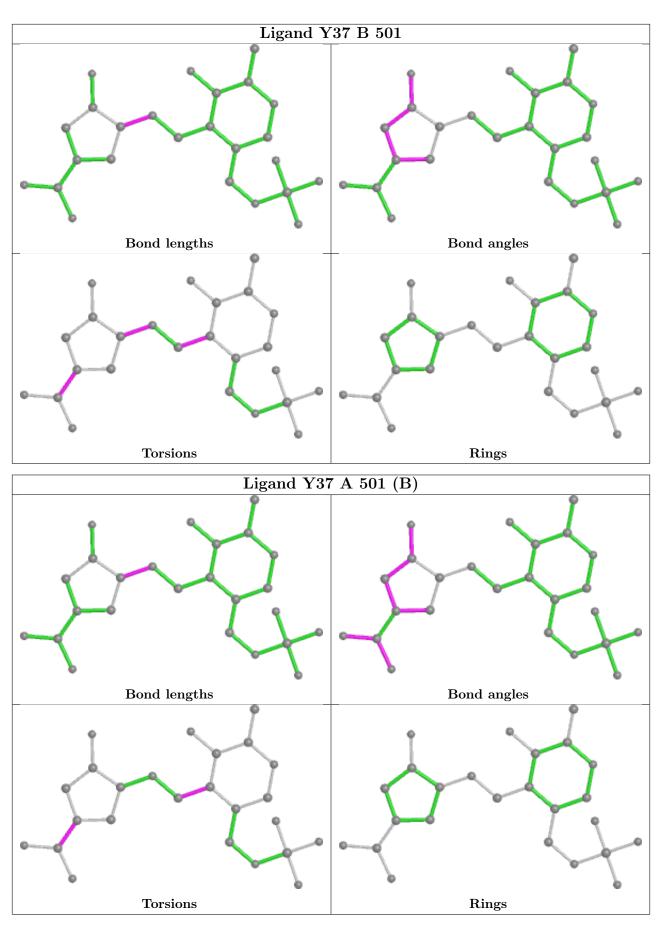
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.





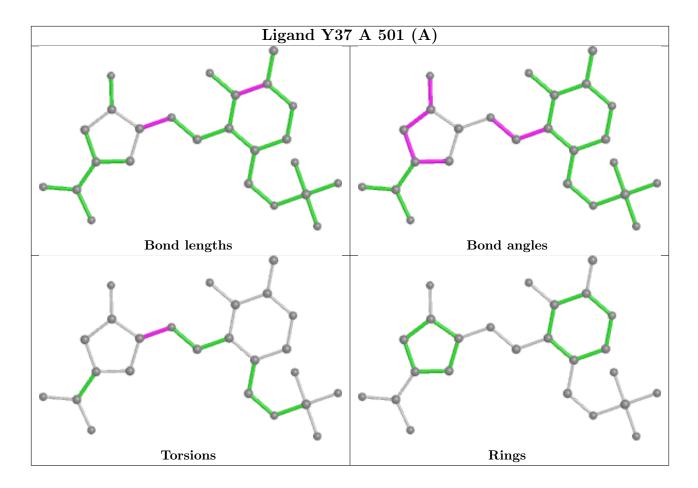












5.7 Other polymers (i)

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 (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, 95^{th} 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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	$\mathbf{Q}{<}0.9$
1	А	404/404~(100%)	0.82	67~(16%) 1 1	24, 45, 70, 109	0
1	В	403/404 (99%)	0.63	40 (9%) 7 5	25, 42, 79, 122	1 (0%)
1	С	404/404 (100%)	0.32	22 (5%) 25 20	25, 37, 57, 100	1 (0%)
All	All	$1211/1212 \ (99\%)$	0.59	129 (10%) 6 4	24, 40, 72, 122	2(0%)

The worst 5 of 129 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	388	TRP	8.9
1	В	367	VAL	7.9
1	А	36	GLY	6.9
1	А	73	VAL	6.8
1	С	107	LYS	6.6

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 monosaccharides in this entry.

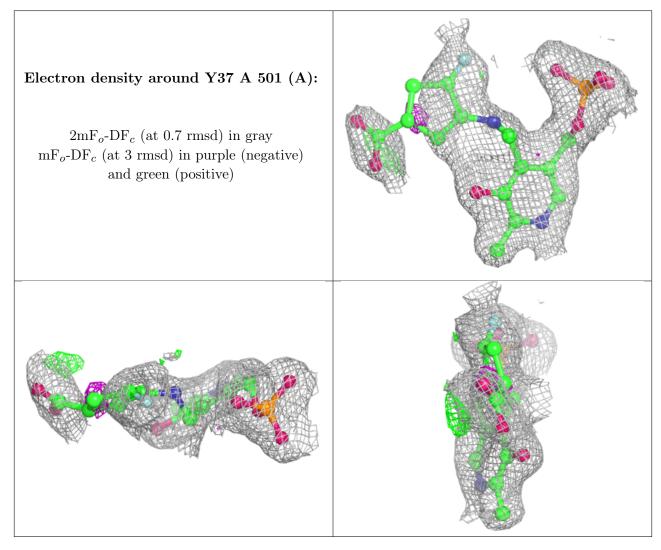
6.4 Ligands (i)

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, 95^{th} 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.

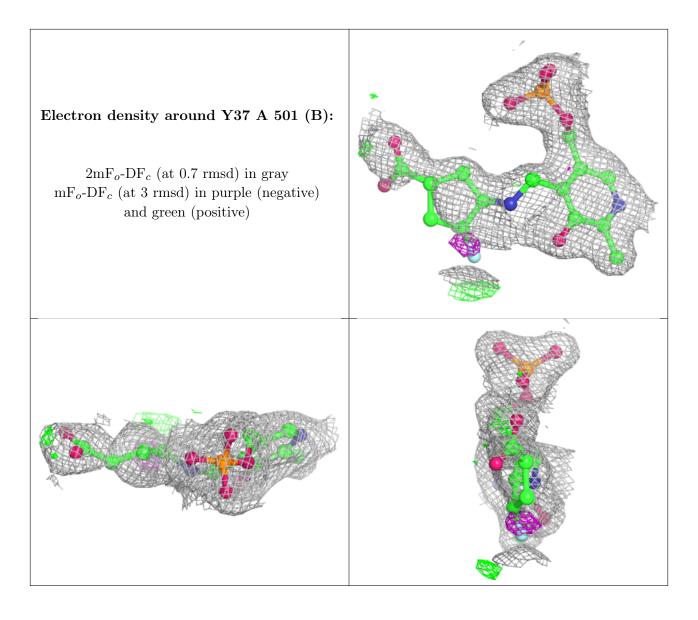


Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q < 0.9
2	Y37	А	501[A]	25/25	0.95	0.21	31,40,59,66	25
2	Y37	А	501[B]	25/25	0.95	0.21	32,38,65,78	25
2	Y37	С	501[A]	25/25	0.96	0.17	27,34,54,67	25
2	Y37	С	501[B]	25/25	0.96	0.17	28,35,61,64	25
2	Y37	В	501	25/25	0.97	0.22	30,47,80,83	0

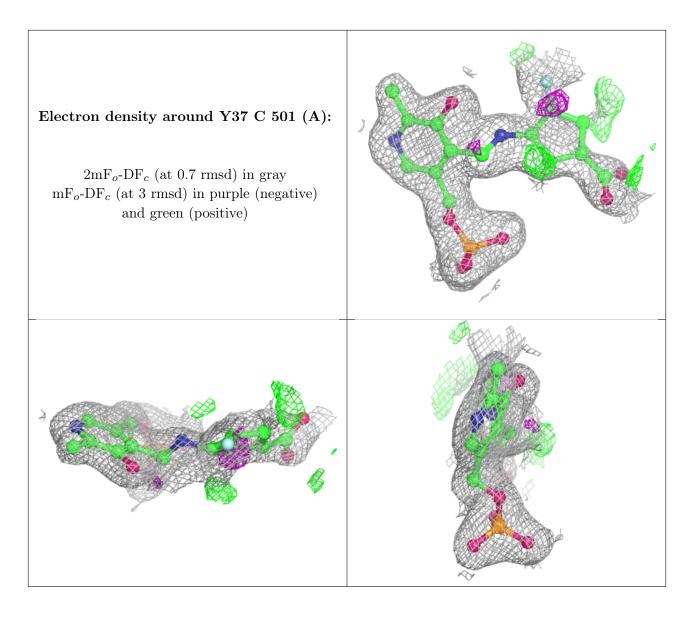
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



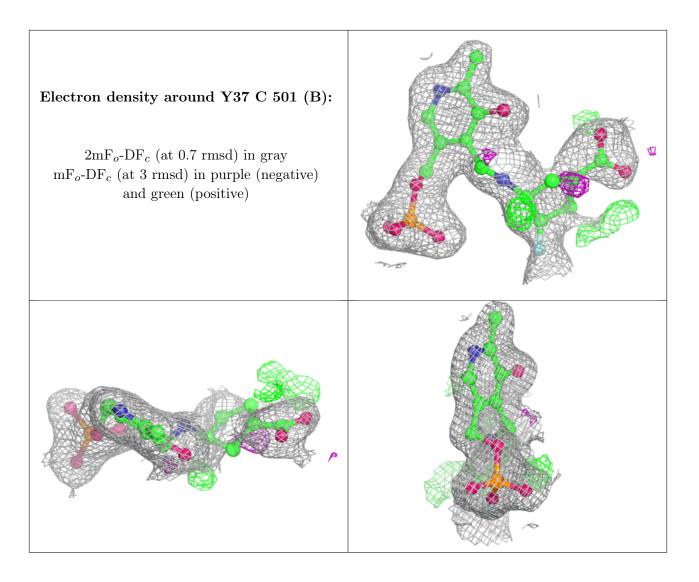




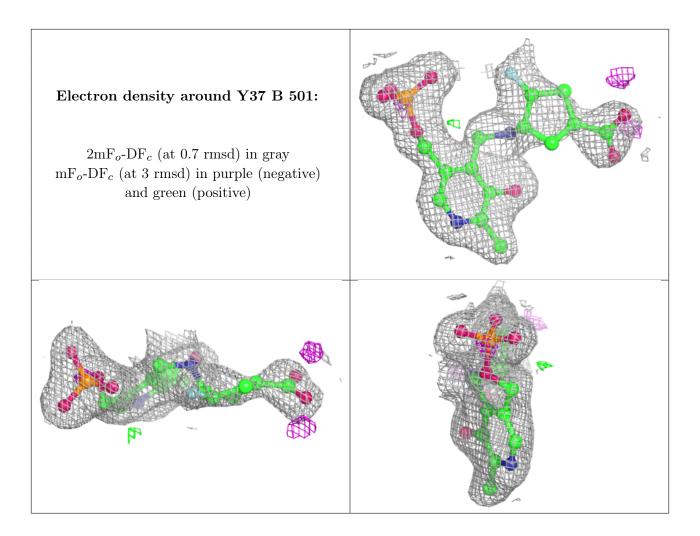












6.5 Other polymers (i)

There are no such residues in this entry.

