

Full wwPDB X-ray Structure Validation Report (i)

Oct 15, 2023 – 11:48 AM EDT

PDB ID : 8D8Y

Title : Crystal structure of ChoE N147A mutant in complex with acetylthiocholine

Authors: Pham, V.D.; Shi, R.

Deposited on : 2022-06-09

Resolution : 1.54 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 (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

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.36

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

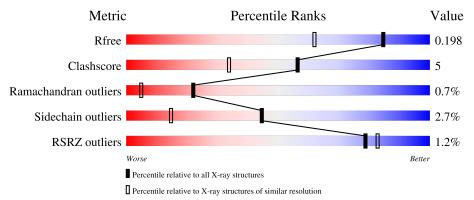
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.54 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	2556 (1.56-1.52)
Clashscore	141614	2634 (1.56-1.52)
Ramachandran outliers	138981	2580 (1.56-1.52)
Sidechain outliers	138945	2577 (1.56-1.52)
RSRZ outliers	127900	2524 (1.56-1.52)

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	A	287	91%	8%	-
1	В	287	90%	9%	-

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
ſ	2	AT3	В	401[A]	-	-	X	-
	4	CL	В	403	-	_	X	-



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5170 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.

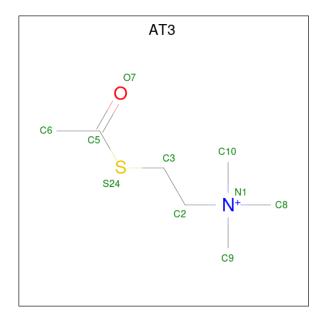
• Molecule 1 is a protein called ChoE.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	287	Total 2265	C 1425	11	O 411	S 5	0	6	0
1	В	287	Total 2265	C 1427		O 413	S 5	0	8	0

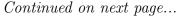
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	ALA	ASN	engineered mutation	UNP Q9HUP2
В	147	ALA	ASN	engineered mutation	UNP Q9HUP2

• Molecule 2 is ACETYLTHIOCHOLINE (three-letter code: AT3) (formula: C₇H₁₆NOS) (labeled as "Ligand of Interest" by depositor).



\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf		
2	A	1	Total 10	C 7	N 1	O 1	S 1	0	1

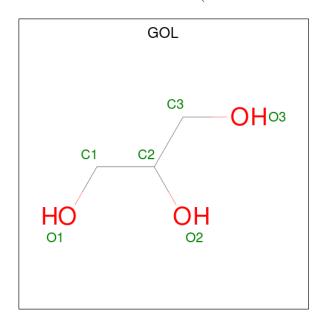




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Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf
2	D	1	Total	С	N	О	S	0	1
	Б	1	10	7	1	1	1		1

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



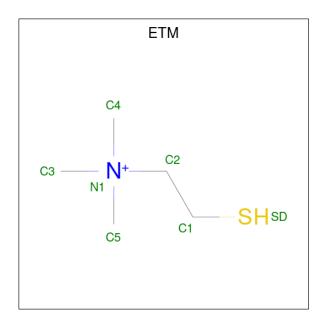
\mathbf{Mol}	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	В	1	Total C O 6 3 3	0	0

• Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0
4	В	1	Total Cl 1 1	0	0

• Molecule 5 is 2-(TRIMETHYLAMMONIUM)ETHYL THIOL (three-letter code: ETM) (formula: $C_5H_{14}NS$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total 7		N 1		0	1
5	В	1	Total 7	C 5	-1	S 1	0	1

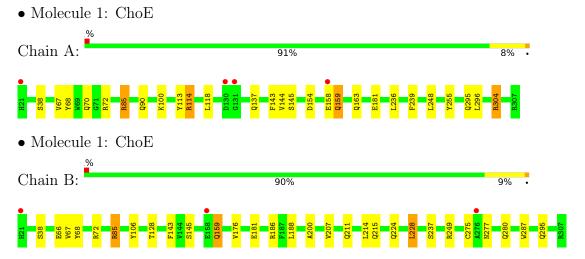
• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	334	Total O 334 334	0	0
6	В	258	Total O 258 258	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.





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	45.63Å 81.79Å 81.26Å	Donositor
a, b, c, α , β , γ	90.00° 99.83° 90.00°	Depositor
Resolution (Å)	45.00 - 1.54	Depositor
Resolution (A)	44.96 - 1.54	EDS
% Data completeness	95.8 (45.00-1.54)	Depositor
(in resolution range)	95.8 (44.96-1.54)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) > 1$	2.08 (at 1.54Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
D D.	0.154 , 0.186	Depositor
R, R_{free}	0.165 , 0.198	DCC
R_{free} test set	4076 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.466	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.37, 48.0	EDS
L-test for twinning ²	$ < L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5170	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.97% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ 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: GOL, AT3, ETM, CL

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	
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.80	0/2337	0.95	1/3175 (0.0%)
1	В	0.79	0/2346	0.92	3/3190 (0.1%)
All	All	0.79	0/4683	0.94	4/6365 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	В	249	ARG	NE-CZ-NH1	7.05	123.83	120.30
1	В	85	ARG	NE-CZ-NH1	5.59	123.09	120.30
1	A	85	ARG	NE-CZ-NH1	5.33	122.96	120.30
1	В	186	ARG	NE-CZ-NH2	-5.26	117.67	120.30

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	2265	0	2210	20	0
1	В	2265	0	2214	16	0
2	A	10	0	16	3	0
2	В	10	0	16	8	0
3	A	6	0	8	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	В	6	0	8	1	0
4	A	1	0	0	0	0
4	В	1	0	0	5	0
5	A	7	0	14	0	0
5	В	7	0	14	2	0
6	A	334	0	0	7	0
6	В	258	0	0	5	0
All	All	5170	0	4500	45	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 5.

All (45) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	${\rm distance}(\mathring{\rm A})$	overlap (Å)
2:B:401[A]:AT3:C6	4:B:403:CL:CL	2.11	1.33
2:B:401[A]:AT3:S24	4:B:403:CL:CL	2.31	1.26
2:B:401[A]:AT3:C5	4:B:403:CL:CL	0.92	0.84
1:A:70:GLN:HG2	6:A:761:HOH:O	1.81	0.81
1:A:239:PHE:HB2	1:A:304[A]:ARG:NH1	2.00	0.77
2:A:401[A]:AT3:H9C3	2:A:401[A]:AT3:S24	2.26	0.75
1:A:159:GLN:HG2	1:A:163:GLN:HB3	1.72	0.72
1:B:277:ASN:ND2	1:B:280:GLN:OE1	2.19	0.71
1:A:159:GLN:HB2	6:A:526:HOH:O	1.91	0.69
1:B:275:CYS:HB3	6:B:714:HOH:O	1.98	0.62
1:B:277:ASN:OD1	1:B:277:ASN:N	2.33	0.61
1:A:154:ASP:OD1	3:A:402:GOL:H11	2.03	0.59
1:A:90:GLN:HG3	6:A:697:HOH:O	2.05	0.57
1:A:304[A]:ARG:HG3	1:A:304[A]:ARG:NH2	2.19	0.57
1:A:248[A]:LEU:HD23	1:A:255:TYR:CD2	2.41	0.55
1:A:85:ARG:NH2	1:A:295[A]:GLN:OE1	2.37	0.55
2:A:401[A]:AT3:S24	2:A:401[A]:AT3:C9	2.95	0.54
1:A:255:TYR:OH	6:A:501:HOH:O	2.18	0.53
2:B:401[A]:AT3:O7	4:B:403:CL:CL	0.34	0.52
1:B:38:SER:OG	2:B:401[A]:AT3:C5	2.57	0.52
1:B:188[B]:LEU:HD13	1:B:237[B]:SER:HB3	1.92	0.52
1:A:236:LEU:C	1:A:236:LEU:HD23	2.32	0.50
1:B:85:ARG:NH2	1:B:295[A]:GLN:OE1	2.44	0.50
1:A:70:GLN:CG	6:A:761:HOH:O	2.48	0.49
1:B:159:GLN:HB2	6:B:552:HOH:O	2.12	0.48
1:B:200:ALA:HB1	3:B:402:GOL:H11	1.96	0.47
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Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance } (\text{\AA}) \end{array}$	Clash overlap (Å)
1:A:137:GLN:HG3	6:B:698:HOH:O	2.14	0.46
1:A:248[A]:LEU:HD21	1:A:296:LEU:HD13	1.97	0.46
1:A:100[A]:LYS:NZ	6:A:508:HOH:O	2.47	0.46
1:A:113:TYR:O	1:A:114[A]:ARG:HD3	2.15	0.46
5:B:404[B]:ETM:H32	5:B:404[B]:ETM:H12	1.64	0.46
1:A:114[B]:ARG:NH2	6:A:503:HOH:O	2.40	0.45
1:A:68:TYR:HB3	1:A:72:ARG:O	2.17	0.44
1:B:215[B]:GLN:NE2	6:B:512:HOH:O	2.52	0.43
1:A:38:SER:OG	2:A:401[A]:AT3:C5	2.67	0.43
1:B:68:TYR:HB3	1:B:72:ARG:O	2.19	0.43
2:B:401[A]:AT3:S24	2:B:401[A]:AT3:C8	3.07	0.43
1:B:106:TYR:CE2	2:B:401[A]:AT3:H103	2.54	0.42
1:A:118:LEU:HD22	1:A:144:VAL:HG11	2.02	0.42
1:B:128:THR:HB	6:B:515:HOH:O	2.19	0.41
1:B:207:VAL:O	1:B:211:GLN:HG3	2.21	0.41
1:B:287:TRP:CE2	5:B:404[B]:ETM:H31	2.55	0.41
2:B:401[A]:AT3:H6C3	4:B:403:CL:CL	2.40	0.41
1:B:176:VAL:HG12	1:B:228:LEU:HD13	2.03	0.40
1:B:214:LEU:HD12	1:B:214:LEU:C	2.42	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	291/287 (101%)	284 (98%)	5 (2%)	2 (1%)	22 5
1	В	293/287 (102%)	283 (97%)	8 (3%)	2 (1%)	22 5
All	All	584/574~(102%)	567 (97%)	13 (2%)	4 (1%)	22 5

All (4) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	145	SER
1	В	145	SER
1	A	67	VAL
1	В	67	VAL

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	Perce	ntiles
1	A	226/220 (103%)	218 (96%)	8 (4%)	36	8
1	В	228/220 (104%)	222 (97%)	6 (3%)	46	16
All	All	454/440 (103%)	440 (97%)	14 (3%)	44	11

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114[A]	ARG
1	A	114[B]	ARG
1	A	143	PHE
1	A	158	GLU
1	A	159	GLN
1	A	181	GLU
1	A	304[A]	ARG
1	A	304[B]	ARG
1	В	66	GLU
1	В	143	PHE
1	В	159	GLN
1	В	181	GLU
1	В	224	GLN
1	В	228	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

\mathbf{Mol}	Chain	Res	Type
1	A	159	GLN

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Mol	Chain	Res	Type
1	В	159	GLN
1	В	163	GLN

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)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	Trino	Type Chain		Link	В	Bond lengths			Bond angles		
IVIOI	$oxed{Mol Type}$	Chain	Res	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
2	AT3	В	401[A]	-	9,9,9	1.04	0	12,12,12	1.39	2 (16%)	
2	AT3	A	401[A]	-	9,9,9	1.09	0	12,12,12	1.70	2 (16%)	
3	GOL	A	402	-	5,5,5	0.23	0	5,5,5	0.38	0	
5	ETM	A	404[B]	-	6,6,6	0.35	0	7,8,8	0.37	0	
3	GOL	В	402	-	5,5,5	0.21	0	5,5,5	0.47	0	
5	ETM	В	404[B]	-	6,6,6	0.22	0	7,8,8	0.31	0	

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	AT3	В	401[A]	-	-	4/7/7/7	-
2	AT3	A	401[A]	-	-	4/7/7/7	-
3	GOL	A	402	-	-	2/4/4/4	-
5	ETM	A	404[B]	-	-	0/4/4/4	-
3	GOL	В	402	-	-	0/4/4/4	-
5	ETM	В	404[B]	_	-	1/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
2	A	401[A]	AT3	C2-C3-S24	-3.69	100.54	109.75
2	A	401[A]	AT3	C3-S24-C5	3.50	120.10	101.68
2	В	401[A]	AT3	C2-C3-S24	-3.43	101.17	109.75
2	В	401[A]	AT3	C3-S24-C5	2.91	116.97	101.68

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401[A]	AT3	O7-C5-S24-C3
5	В	404[B]	ETM	SD-C1-C2-N1
2	A	401[A]	AT3	C6-C5-S24-C3
2	В	401[A]	AT3	N1-C2-C3-S24
3	A	402	GOL	O1-C1-C2-O2
2	A	401[A]	AT3	C2-C3-S24-C5
2	В	401[A]	AT3	C2-C3-S24-C5
2	В	401[A]	AT3	O7-C5-S24-C3
3	A	402	GOL	O1-C1-C2-C3
2	A	401[A]	AT3	N1-C2-C3-S24
2	В	401[A]	AT3	C6-C5-S24-C3

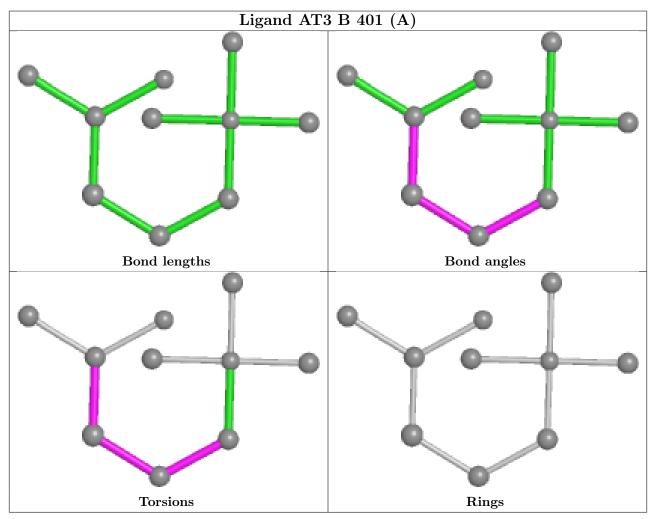
There are no ring outliers.

5 monomers are involved in 15 short contacts:

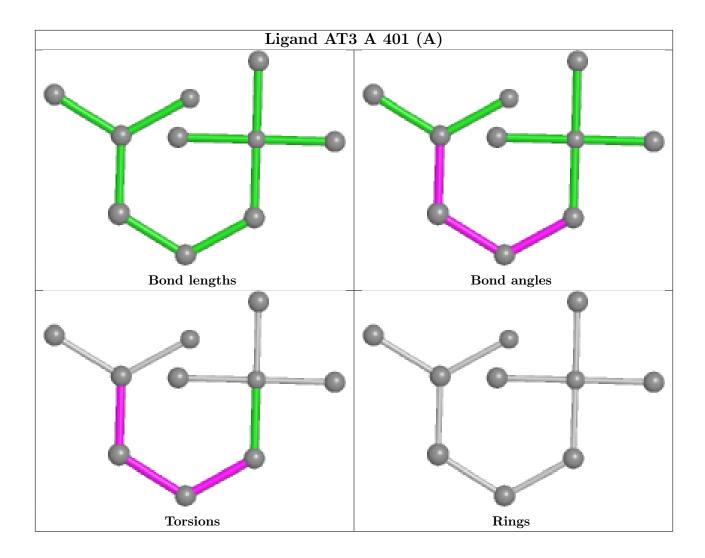
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	401[A]	AT3	8	0
2	A	401[A]	AT3	3	0
3	A	402	GOL	1	0
3	В	402	GOL	1	0
5	В	404[B]	ETM	2	0



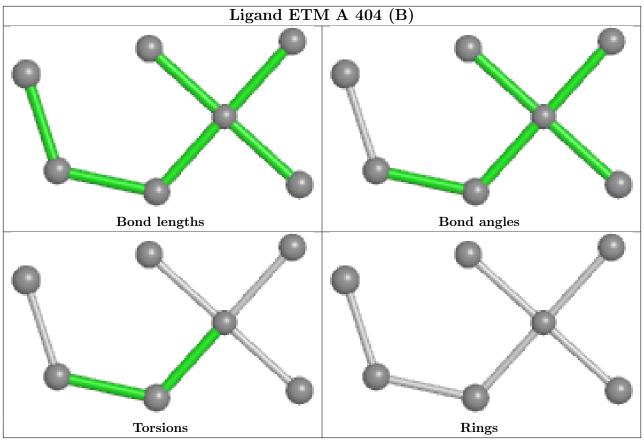
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

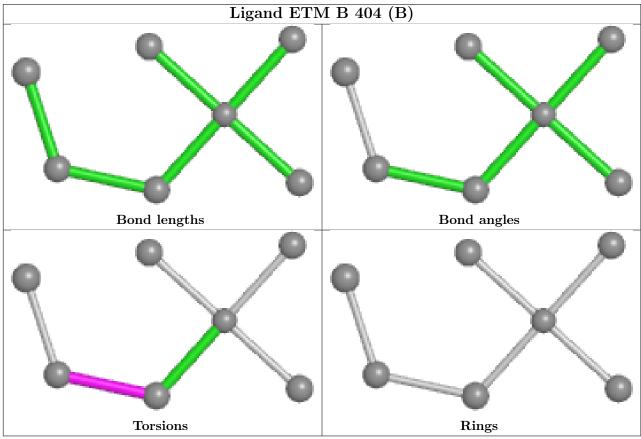














5.7 Other polymers (i)

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	<RSRZ $>$	# RSRZ > 2		$OWAB(A^2)$	Q < 0.9
1	A	287/287 (100%)	-0.37	4 (1%) 75 79	9	11, 17, 34, 62	0
1	В	287/287 (100%)	-0.34	3 (1%) 82 85	5	15, 22, 44, 68	0
All	All	574/574 (100%)	-0.36	7 (1%) 79 83	3	11, 20, 41, 68	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	GLY	4.2
1	В	21	HIS	3.6
1	A	158	GLU	2.2
1	В	276	ALA	2.1
1	В	158	GLU	2.1
1	A	21	HIS	2.1
1	A	130	ASP	2.1

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.

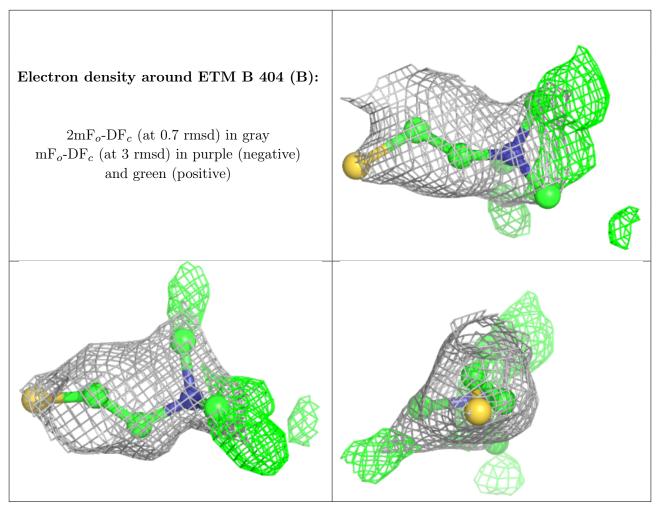
6.4 Ligands (i)

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	GOL	A	402	6/6	0.77	0.21	38,45,46,48	0
3	GOL	В	402	6/6	0.80	0.15	41,49,57,59	0
5	ETM	В	404[B]	7/7	0.84	0.20	33,35,36,39	7
5	ETM	A	404[B]	7/7	0.87	0.17	27,27,28,29	7
2	AT3	A	401[A]	10/10	0.94	0.11	24,42,44,46	10
2	AT3	В	401[A]	10/10	0.95	0.10	29,42,45,46	10
4	CL	A	403	1/1	0.99	0.05	13,13,13,13	1
4	CL	В	403	1/1	0.99	0.03	18,18,18,18	1

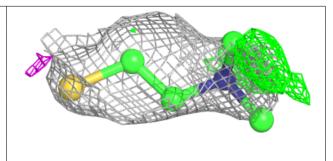
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

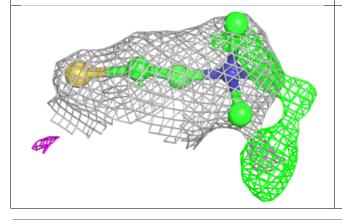


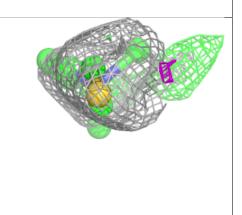


Electron density around ETM A 404 (B):

 $2 {\rm mF}_o\text{-}{\rm DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

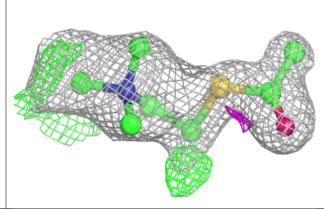


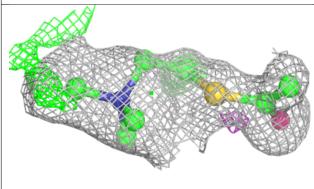


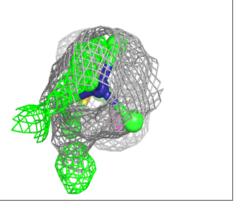


Electron density around AT3 A 401 (A):

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



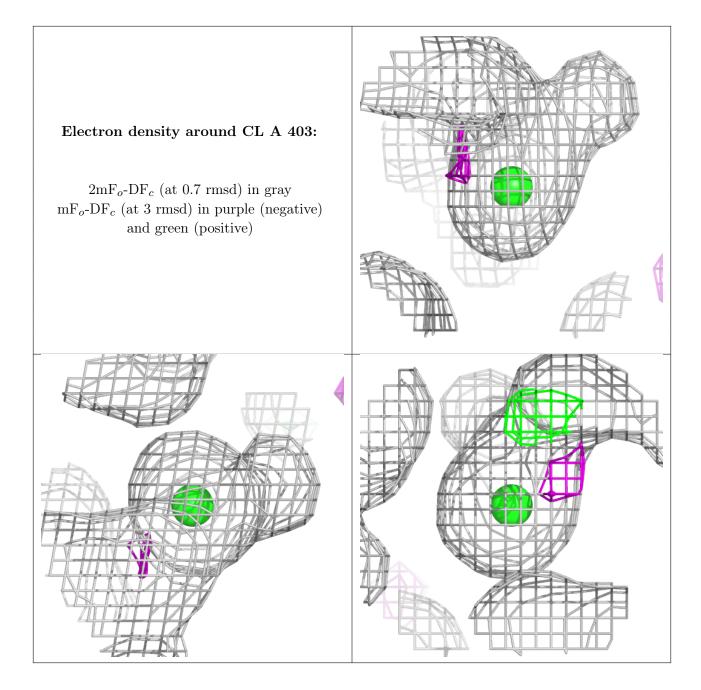




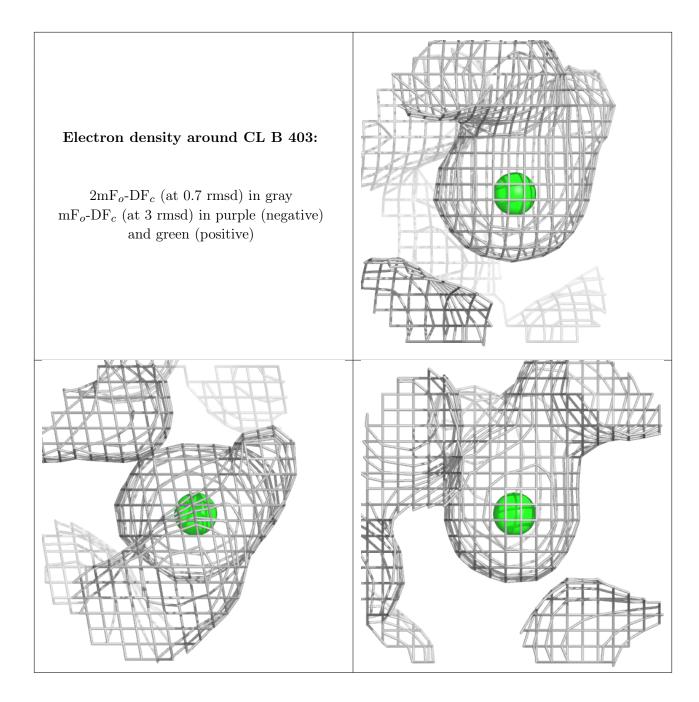


Electron density around AT3 B 401 (A): 2mF_o-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)









6.5 Other polymers (i)

There are no such residues in this entry.

