



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 6, 2021 – 07:17 am BST

PDB ID : 7NB2
Title : Crystal structure of human choline alpha in complex with an inhibitor
Authors : Casale, E.; Fasolini, M.
Deposited on : 2021-01-25
Resolution : 2.40 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.23.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

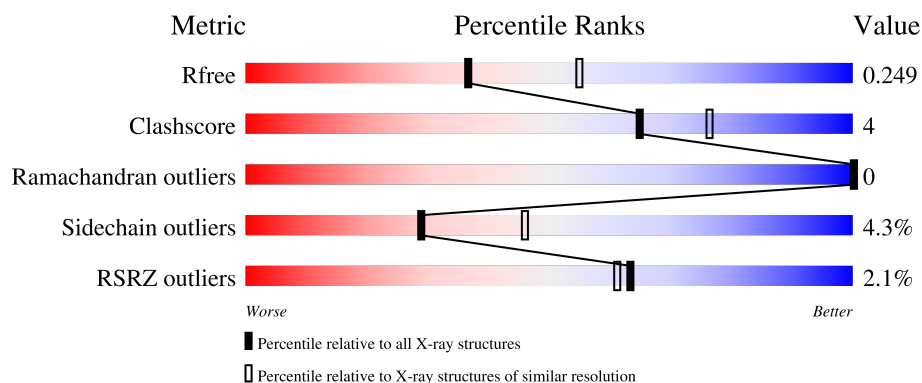
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	AAA	384	
1	BBB	384	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5870 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.

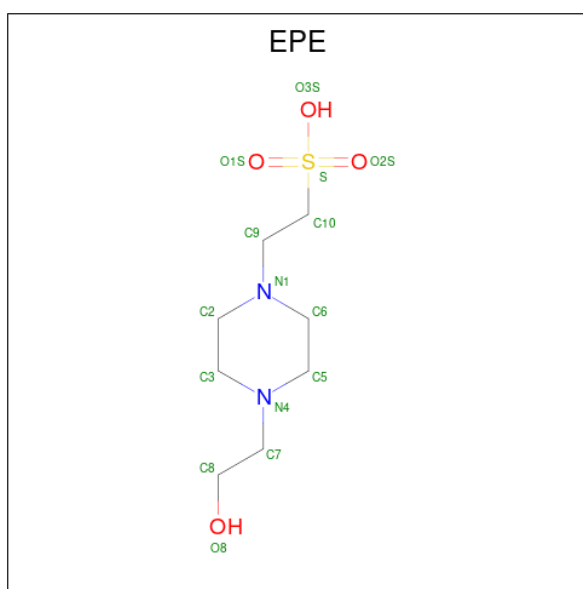
- Molecule 1 is a protein called Choline kinase alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	347	Total	C	N	O	S	0	0	0
			2829	1835	466	512	16			
1	BBB	352	Total	C	N	O	S	0	0	0
			2873	1863	481	513	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	74	GLY	-	expression tag	UNP P35790
BBB	74	GLY	-	expression tag	UNP P35790

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



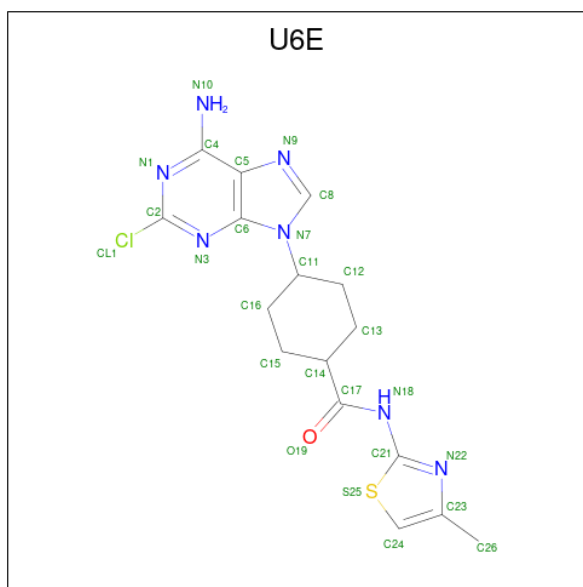
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	BBB	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 3 is 4-(6-azanyl-2-chloranyl-purin-9-yl)- {N}-(4-methyl-1,3-thiazol-2-yl)cyclohexan e-1-carboxamide (three-letter code: U6E) (formula: C₁₆H₁₈ClN₇OS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	AAA	1	Total	C	Cl	N	O	S	0
			26	16	1	7	1	1	
3	BBB	1	Total	C	Cl	N	O	S	0
			26	16	1	7	1	1	

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AAA	1	Total	Mg	0	0
			1	1		
4	BBB	1	Total	Mg	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	36	Total 36	O 36	0	0
5	BBB	33	Total 33	O 33	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.59Å 120.93Å 130.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.20 – 2.40 51.14 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (51.20-2.40) 99.6 (51.14-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.202 , 0.248 0.203 , 0.249	Depositor DCC
R_{free} test set	1718 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	52.0	Xtrriage
Anisotropy	0.667	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5870	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% 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: EPE, MG, U6E

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	AAA	0.37	0/2904	0.71	0/3920
1	BBB	0.35	0/2949	0.68	0/3981
All	All	0.36	0/5853	0.69	0/7901

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	AAA	2829	0	2744	26	0
1	BBB	2873	0	2808	25	0
2	AAA	30	0	35	2	0
2	BBB	15	0	17	0	0
3	AAA	26	0	0	1	0
3	BBB	26	0	0	0	0
4	AAA	1	0	0	0	0
4	BBB	1	0	0	0	0
5	AAA	36	0	0	1	0
5	BBB	33	0	0	0	0
All	All	5870	0	5604	47	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 4.

All (47) 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:AAA:292:LEU:HD22	1:AAA:457:VAL:HG11	1.74	0.67
1:BBB:252:THR:HG22	1:BBB:256:TYR:CE2	2.30	0.67
1:BBB:200:PHE:HB2	1:BBB:201:PRO:CD	2.26	0.66
1:AAA:267:GLU:H	1:AAA:267:GLU:CD	1.99	0.66
1:AAA:200:PHE:HB2	1:AAA:201:PRO:CD	2.30	0.62
1:AAA:200:PHE:HB2	1:AAA:201:PRO:HD2	1.81	0.62
1:AAA:192:LEU:HD13	1:AAA:238:MET:SD	2.43	0.59
1:BBB:451:GLN:HE21	1:BBB:451:GLN:HA	1.67	0.59
1:AAA:122:ASN:ND2	1:AAA:146:ARG:HH11	2.02	0.57
1:AAA:124:LEU:HD22	1:AAA:144:LEU:HD21	1.87	0.56
1:BBB:195:LYS:H	1:BBB:207:GLN:HE21	1.54	0.56
1:AAA:245:GLU:CD	1:BBB:104:ARG:HH12	2.12	0.53
1:BBB:149:GLY:O	1:BBB:150:ALA:CB	2.57	0.52
1:BBB:97:GLU:HA	1:BBB:104:ARG:HH22	1.75	0.51
1:AAA:282:LEU:N	1:AAA:283:PRO:CD	2.74	0.51
1:AAA:97:GLU:HG2	1:BBB:245:GLU:CG	2.42	0.50
1:BBB:389:ASP:OD2	1:BBB:389:ASP:N	2.45	0.50
1:BBB:252:THR:CG2	1:BBB:256:TYR:CE2	2.95	0.50
1:BBB:371:THR:O	1:BBB:375:GLN:HG3	2.12	0.49
1:AAA:97:GLU:HG2	1:BBB:245:GLU:HG3	1.94	0.48
1:BBB:220:SER:HB3	1:BBB:350:TRP:CZ3	2.49	0.48
1:AAA:145:LEU:HD11	1:AAA:147:LEU:HD23	1.94	0.48
1:BBB:321:SER:O	1:BBB:325:LYS:HD3	2.13	0.48
1:AAA:146:ARG:O	1:AAA:203:GLY:HA3	2.15	0.47
1:AAA:227:GLU:HB3	1:AAA:326:LEU:HD12	1.96	0.47
1:BBB:359:TYR:CD1	1:BBB:360:PRO:HA	2.48	0.47
1:BBB:411:ARG:O	1:BBB:414:LEU:HB2	2.15	0.47
1:AAA:431:SER:O	2:AAA:502:EPE:H51	2.14	0.46
1:BBB:282:LEU:N	1:BBB:283:PRO:CD	2.78	0.46
1:AAA:231:LYS:HE2	1:AAA:326:LEU:O	2.16	0.45
1:BBB:200:PHE:HB2	1:BBB:201:PRO:HD2	1.96	0.45
1:AAA:360:PRO:O	1:AAA:361:PHE:HB2	2.16	0.45
1:BBB:423:TRP:O	1:BBB:427:GLN:HG2	2.17	0.45
1:AAA:145:LEU:HD11	1:AAA:147:LEU:CD2	2.49	0.43
1:AAA:453:ARG:NH2	5:AAA:603:HOH:O	2.48	0.43
1:BBB:359:TYR:CG	1:BBB:360:PRO:HA	2.54	0.43
1:BBB:369:TYR:CG	1:BBB:370:PRO:HD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:245:GLU:OE2	1:BBB:104:ARG:NH1	2.51	0.42
1:BBB:262:ARG:HH11	1:BBB:262:ARG:HG3	1.84	0.42
1:AAA:375:GLN:O	1:AAA:379:ILE:HG23	2.20	0.42
1:BBB:262:ARG:HG3	1:BBB:262:ARG:NH1	2.35	0.42
1:AAA:87:THR:HG23	1:AAA:90:ARG:HD3	2.03	0.41
1:AAA:433:ILE:HG21	2:AAA:501:EPE:H71	2.01	0.41
1:AAA:265:PHE:O	1:AAA:271:ILE:HD11	2.21	0.41
1:AAA:87:THR:HA	1:AAA:90:ARG:HG3	2.03	0.41
1:AAA:144:LEU:HD13	3:AAA:503:U6E:C4	2.52	0.40
1:BBB:427:GLN:O	1:BBB:431:SER:HB3	2.21	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	AAA	341/384 (89%)	324 (95%)	17 (5%)	0	100	100
1	BBB	348/384 (91%)	330 (95%)	18 (5%)	0	100	100
All	All	689/768 (90%)	654 (95%)	35 (5%)	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	AAA	300/342 (88%)	289 (96%)	11 (4%)	34	53
1	BBB	304/342 (89%)	289 (95%)	15 (5%)	25	40
All	All	604/684 (88%)	578 (96%)	26 (4%)	29	46

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	87	THR
1	AAA	110	GLU
1	AAA	113	ILE
1	AAA	114	SER
1	AAA	121	SER
1	AAA	123	MET
1	AAA	199	ILE
1	AAA	200	PHE
1	AAA	267	GLU
1	AAA	331	PHE
1	AAA	366	ILE
1	BBB	82	GLN
1	BBB	120	LEU
1	BBB	121	SER
1	BBB	123	MET
1	BBB	131	ASP
1	BBB	239	LYS
1	BBB	258	LYS
1	BBB	268	GLU
1	BBB	291	SER
1	BBB	329	ILE
1	BBB	330	ASP
1	BBB	342	ASP
1	BBB	357	GLU
1	BBB	399	SER
1	BBB	451	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	U6E	BBB	502	-	22,29,29	1.99	5 (22%)	23,42,42	3.99	5 (21%)
3	U6E	AAA	503	-	22,29,29	1.83	3 (13%)	23,42,42	3.66	8 (34%)
2	EPE	BBB	501	-	15,15,15	1.11	1 (6%)	18,20,20	1.17	1 (5%)
2	EPE	AAA	501	-	15,15,15	0.68	1 (6%)	18,20,20	0.65	0
2	EPE	AAA	502	-	15,15,15	1.12	1 (6%)	18,20,20	1.02	1 (5%)

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	U6E	BBB	502	-	-	0/6/22/22	0/4/4/4
3	U6E	AAA	503	-	-	3/6/22/22	0/4/4/4
2	EPE	BBB	501	-	-	3/9/19/19	0/1/1/1
2	EPE	AAA	501	-	-	3/9/19/19	0/1/1/1
2	EPE	AAA	502	-	-	4/9/19/19	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	BBB	502	U6E	C2-N3	6.22	1.35	1.30
3	AAA	503	U6E	C2-N3	6.21	1.35	1.30
2	AAA	502	EPE	O1S-S	3.97	1.56	1.45
3	BBB	502	U6E	C2-CL1	3.88	1.84	1.73
2	BBB	501	EPE	O2S-S	3.67	1.55	1.45
3	AAA	503	U6E	C24-S25	-2.93	1.66	1.70
3	BBB	502	U6E	C5-C6	-2.92	1.33	1.40
3	AAA	503	U6E	C5-C6	-2.65	1.33	1.40
3	BBB	502	U6E	C24-S25	-2.63	1.66	1.70
2	AAA	501	EPE	O3S-S	2.33	1.55	1.47
3	BBB	502	U6E	C4-C5	-2.05	1.35	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	BBB	502	U6E	N3-C2-N1	-11.93	118.61	130.62
3	AAA	503	U6E	N3-C2-N1	-11.04	119.51	130.62
3	BBB	502	U6E	CL1-C2-N3	9.90	124.15	115.70
3	AAA	503	U6E	C2-N3-C6	9.71	121.78	114.09
3	BBB	502	U6E	C2-N3-C6	9.10	121.29	114.09
3	AAA	503	U6E	CL1-C2-N3	4.86	119.85	115.70
3	AAA	503	U6E	CL1-C2-N1	3.84	120.64	115.15
3	AAA	503	U6E	C21-N18-C17	-3.71	120.76	129.02
3	AAA	503	U6E	C26-C23-C24	-3.60	119.54	129.06
2	BBB	501	EPE	O3S-S-O1S	3.33	119.42	111.27
3	BBB	502	U6E	C26-C23-C24	-3.03	121.04	129.06
3	AAA	503	U6E	C6-C5-N9	-2.98	106.30	109.40
3	AAA	503	U6E	O19-C17-C14	-2.69	118.61	122.12
3	BBB	502	U6E	C2-N1-C4	2.68	119.96	116.64
2	AAA	502	EPE	O3S-S-O2S	2.41	117.17	111.27

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	502	EPE	S-C10-C9-N1
2	BBB	501	EPE	N4-C7-C8-O8
2	AAA	502	EPE	N4-C7-C8-O8
2	AAA	501	EPE	C9-C10-S-O3S
2	AAA	502	EPE	C10-C9-N1-C2
2	AAA	502	EPE	C10-C9-N1-C6
2	AAA	501	EPE	C9-C10-S-O1S

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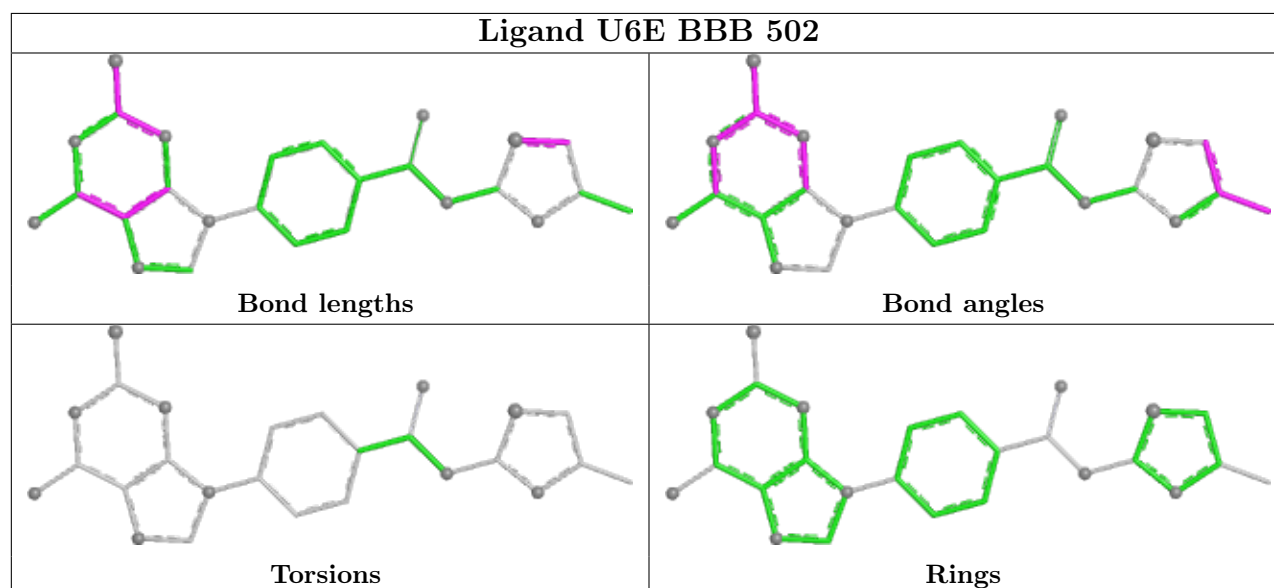
Mol	Chain	Res	Type	Atoms
2	AAA	501	EPE	C9-C10-S-O2S
3	AAA	503	U6E	C15-C14-C17-O19
2	BBB	501	EPE	C10-C9-N1-C2
2	BBB	501	EPE	C10-C9-N1-C6
3	AAA	503	U6E	C15-C14-C17-N18
3	AAA	503	U6E	C13-C14-C17-O19

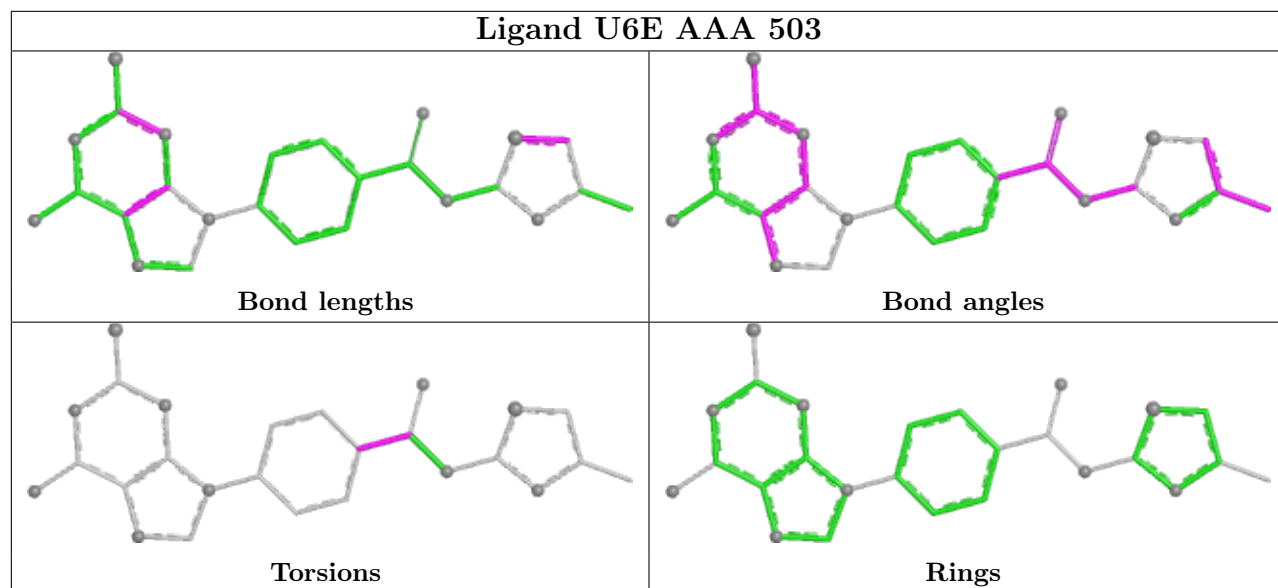
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	AAA	503	U6E	1	0
2	AAA	501	EPE	1	0
2	AAA	502	EPE	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.





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, 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	AAA	347/384 (90%)	0.28	8 (2%) 60 58	38, 56, 107, 140	1 (0%)
1	BBB	352/384 (91%)	0.17	7 (1%) 65 63	42, 61, 98, 121	0
All	All	699/768 (91%)	0.23	15 (2%) 63 61	38, 59, 102, 140	1 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	111	PHE	5.5
1	AAA	129	LEU	4.4
1	AAA	100	PRO	3.7
1	BBB	87	THR	3.7
1	AAA	105	GLY	3.3
1	AAA	132	THR	3.3
1	AAA	130	PRO	3.1
1	AAA	107	ARG	3.0
1	BBB	390	PHE	2.7
1	AAA	147	LEU	2.4
1	BBB	383	LEU	2.4
1	BBB	393	LEU	2.4
1	BBB	82	GLN	2.1
1	BBB	388	ASN	2.0
1	BBB	400	ILE	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 monosaccharides in this entry.

6.4 Ligands ⓘ

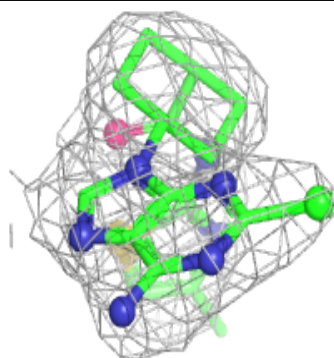
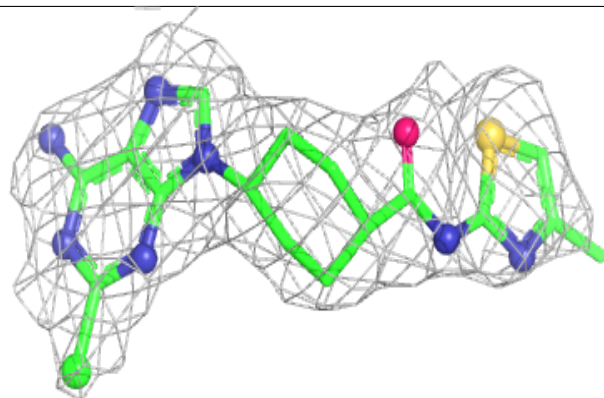
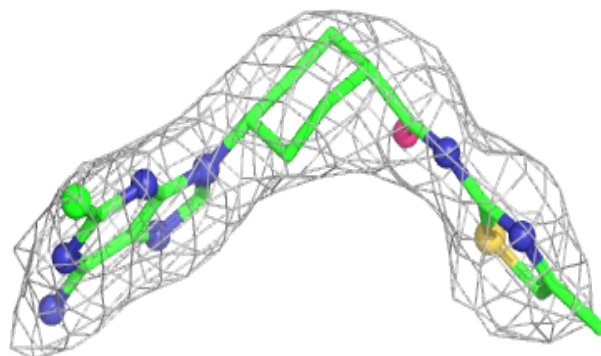
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	AAA	504	1/1	0.77	0.12	48,48,48,48	0
2	EPE	AAA	501	15/15	0.79	0.23	84,100,111,114	0
2	EPE	BBB	501	15/15	0.81	0.22	75,89,124,127	0
4	MG	BBB	503	1/1	0.83	0.24	53,53,53,53	0
2	EPE	AAA	502	15/15	0.87	0.23	74,89,108,115	0
3	U6E	AAA	503	26/26	0.92	0.18	54,62,92,95	0
3	U6E	BBB	502	26/26	0.93	0.19	51,61,81,104	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.

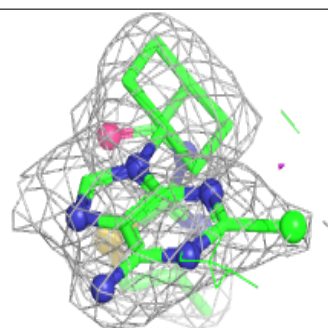
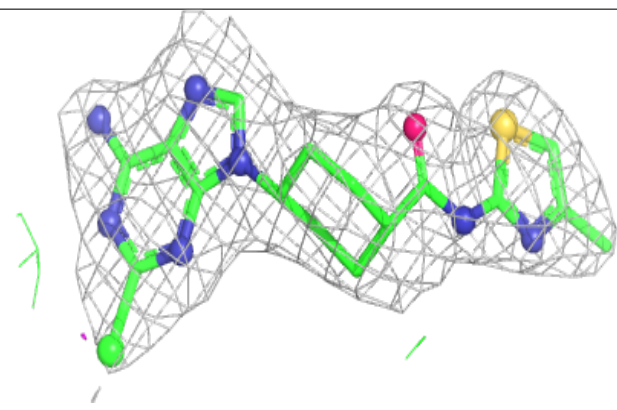
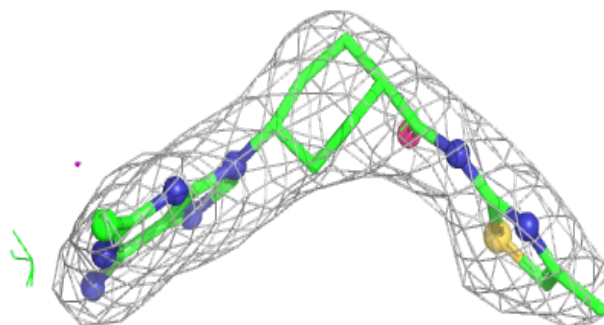
Electron density around U6E AAA 503:

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 U6E BBB 502:

$2mF_o-DF_c$ (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.