



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2026 – 09:42 PM EDT

PDB ID : 1Z9Y / pdb_00001z9y
Title : carbonic anhydrase II in complex with furosemide as sulfonamide inhibitor
Authors : Honndorf, V.S.; Heine, A.; Klebe, G.; Supuran, C.T.
Deposited on : 2005-04-05
Resolution : 1.66 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

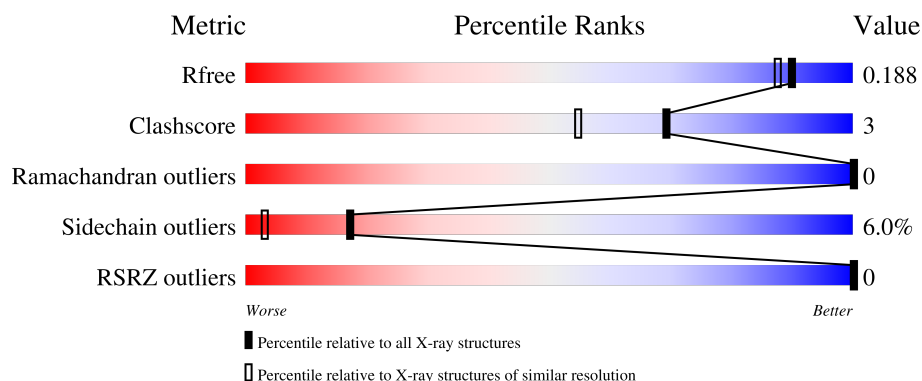
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.66 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	259	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2294 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.

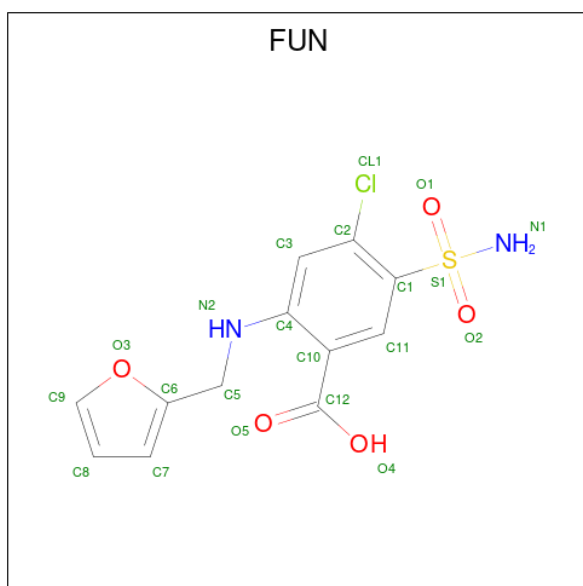
- Molecule 1 is a protein called Carbonic anhydrase II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	0	0	0
			2020	1296	346	376	2			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

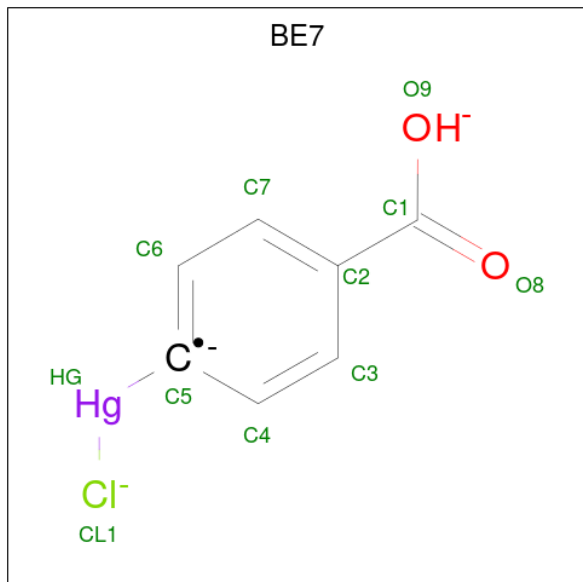
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is 5-(AMINOSULFONYL)-4-CHLORO-2-[(2-FURYL METHYL)AMINO]BENZ OIC ACID (CCD ID: FUN) (formula: C₁₂H₁₁ClN₂O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			21	12	1	2	5		

- Molecule 4 is (4-CARBOXYPHENYL)(CHLORO)MERCURY (CCD ID: BE7) (formula: $C_7H_5ClHgO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	Hg	O	0	0
			10	7	1	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	241	Total	O	0	0
			241	241		

- Molecule 1: Carbonic anhydrase II

C206	K213	K215	S219	Q222	K225	F231	V253	K254	Q255	I256	K257	F260	LYS	SER	HIS	H4	K9	R27	A38	K39	Y40	I41	P42	S43	L44	K45	D82	Q83	A84	L57	H64	E69	D75	K76	A77	K80	D85	G86	R89	Q92	F93	W97	H107	A116	L148	K149	Q158	K159	V160	V161	D162	V163	L198	L202
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.31Å 41.95Å 72.15Å 90.00° 104.57° 90.00°	Depositor
Resolution (Å)	10.00 – 1.66 10.00 – 1.66	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-1.66) 85.3 (10.00-1.66)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.14 (at 1.66Å)	Xtriage
Refinement program	SHELXL-97, CNS	Depositor
R, R_{free}	0.172 , 0.245 0.186 , 0.188	Depositor DCC
R_{free} test set	1390 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	2.8	Xtriage
Anisotropy	0.841	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.30$, $\langle L^2 \rangle = 0.13$	Xtriage
Estimated twinning fraction	0.219 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	2294	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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: BE7, FUN, ZN

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	0.52	0/2081	1.41	21/2831 (0.7%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	CYS	CB-CA-C	8.40	124.55	110.44
1	A	64	HIS	CA-CB-CG	8.15	121.95	113.80
1	A	206	CYS	O-C-N	-7.46	112.80	122.20
1	A	226	PHE	CA-CB-CG	7.40	121.20	113.80
1	A	231	PHE	CA-CB-CG	6.73	120.53	113.80
1	A	9	LYS	CA-C-N	-6.69	111.50	122.54
1	A	9	LYS	C-N-CA	-6.69	111.50	122.54
1	A	207	VAL	N-CA-C	6.52	117.30	108.17
1	A	27	ARG	NE-CZ-NH2	6.41	124.97	119.20
1	A	107	HIS	CA-CB-CG	-6.20	107.60	113.80
1	A	198	LEU	CB-CA-C	-6.14	98.78	109.64
1	A	93	PHE	CA-CB-CG	6.02	119.82	113.80
1	A	198	LEU	CA-C-O	-5.89	115.08	121.44
1	A	85	ASP	CA-CB-CG	-5.83	106.77	112.60
1	A	206	CYS	N-CA-C	-5.41	104.95	113.02
1	A	206	CYS	CA-CB-SG	5.29	126.56	114.40
1	A	97	TRP	CA-C-N	-5.20	116.48	122.77
1	A	97	TRP	C-N-CA	-5.20	116.48	122.77
1	A	86	GLY	CA-C-N	5.16	130.27	122.99
1	A	86	GLY	C-N-CA	5.16	130.27	122.99
1	A	89	ARG	NE-CZ-NH1	-5.03	116.47	121.50

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	A	2020	0	1943	14	0
2	A	2	0	0	0	0
3	A	21	0	9	0	0
4	A	10	0	4	0	0
5	A	241	0	0	4	0
All	All	2294	0	1956	14	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 3.

All (14) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LYS:NZ	1:A:215:PRO:HG2	2.07	0.69
1:A:158:GLN:HA	1:A:161:VAL:HG22	1.88	0.55
1:A:54:ALA:HB2	1:A:77:ALA:HB2	1.89	0.54
1:A:38:ALA:HA	5:A:2222:HOH:O	2.11	0.50
1:A:219:SER:OG	1:A:222:GLN:HG3	2.12	0.50
1:A:40:TYR:CE2	1:A:42:PRO:HG3	2.48	0.49
1:A:198:LEU:HB2	1:A:203:LEU:HA	1.95	0.48
1:A:257:LYS:HG3	5:A:2228:HOH:O	2.13	0.48
1:A:69:GLU:HB3	5:A:2072:HOH:O	2.14	0.46
1:A:160:VAL:O	1:A:163:VAL:HG12	2.16	0.45
1:A:149:LYS:HZ2	1:A:215:PRO:HG2	1.79	0.45
1:A:213:LYS:NZ	5:A:2069:HOH:O	2.51	0.42
1:A:213:LYS:HD2	1:A:260:PHE:CZ	2.55	0.42
1:A:116:ALA:HB3	1:A:148:LEU:HB2	2.02	0.41

There are no symmetry-related clashes.

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	254/259 (98%)	241 (95%)	13 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	216/224 (96%)	203 (94%)	13 (6%)	17	3

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	43	SER
1	A	45	LYS
1	A	52	ASP
1	A	57	LEU
1	A	75	ASP
1	A	80	LYS
1	A	85	ASP
1	A	92	GLN
1	A	149	LYS
1	A	225	LYS
1	A	253	ASN
1	A	255	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	ASN
1	A	230	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
4	BE7	A	700	1	7,10,11	1.69	2 (28%)	8,13,14	1.09	0
3	FUN	A	500	2	22,22,22	1.91	4 (18%)	32,32,32	2.12	11 (34%)

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
4	BE7	A	700	1	-	0/4/14/16	0/1/1/1
3	FUN	A	500	2	-	2/15/15/15	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	FUN	S1-N1	-6.39	1.48	1.60
3	A	500	FUN	O5-C12	3.30	1.32	1.22
4	A	700	BE7	C4-C3	3.22	1.43	1.35
3	A	500	FUN	O4-C12	-3.11	1.21	1.30
3	A	500	FUN	O1-S1	2.69	1.48	1.43
4	A	700	BE7	C2-C1	2.44	1.53	1.48

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	FUN	C11-C10-C4	4.72	123.34	119.09
3	A	500	FUN	O3-C6-C7	4.54	117.91	109.61
3	A	500	FUN	O3-C9-C8	3.57	116.81	110.44
3	A	500	FUN	C9-O3-C6	-3.21	99.82	106.22
3	A	500	FUN	O2-S1-N1	3.07	111.78	107.35
3	A	500	FUN	C5-N2-C4	-2.77	119.83	123.94
3	A	500	FUN	C3-C2-C1	2.67	123.94	121.36
3	A	500	FUN	C10-C11-C1	-2.64	118.72	122.04
3	A	500	FUN	C8-C7-C6	-2.51	102.02	107.05
3	A	500	FUN	C5-C6-C7	-2.27	125.30	133.07
3	A	500	FUN	C7-C8-C9	-2.22	103.22	106.94

There are no chirality outliers.

All (2) torsion outliers are listed below:

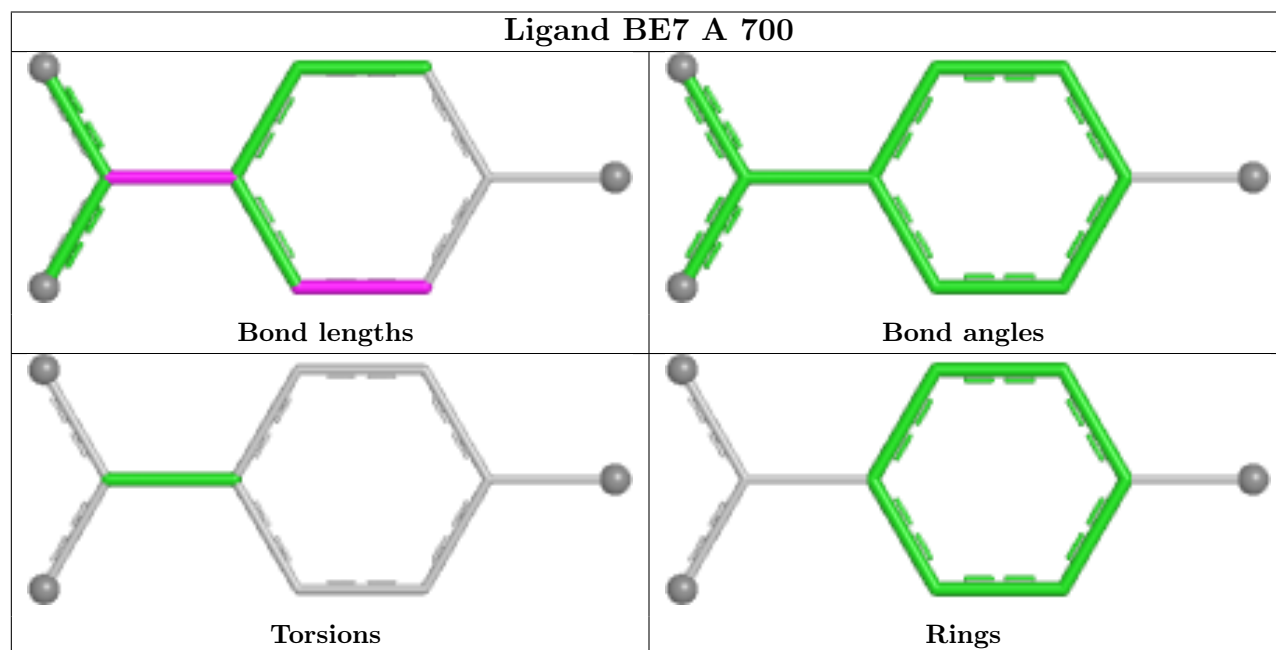
Mol	Chain	Res	Type	Atoms
3	A	500	FUN	C11-C1-S1-N1
3	A	500	FUN	C2-C1-S1-N1

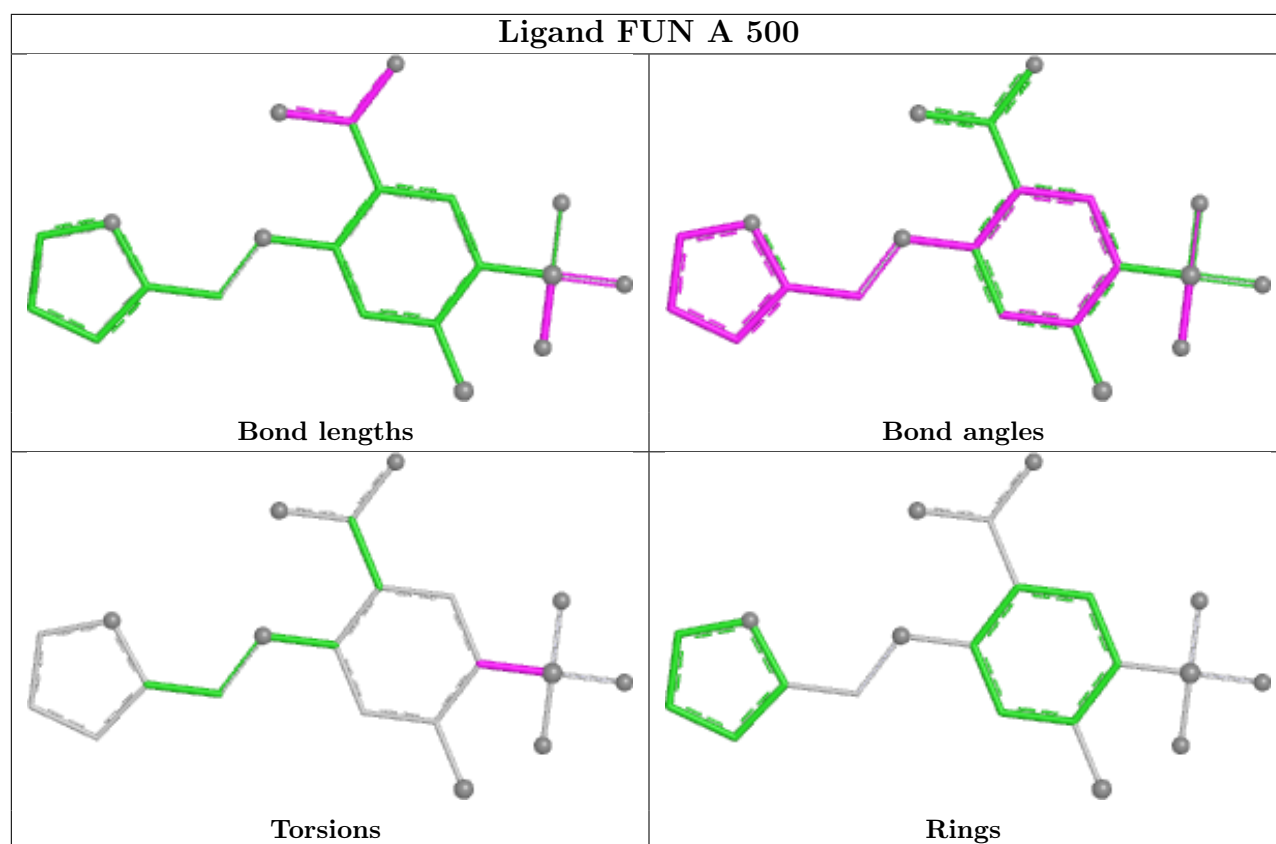
There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [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	A	256/259 (98%)	-1.19	0 100 100	9, 17, 31, 54	0

There are no RSRZ outliers to report.

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](#)

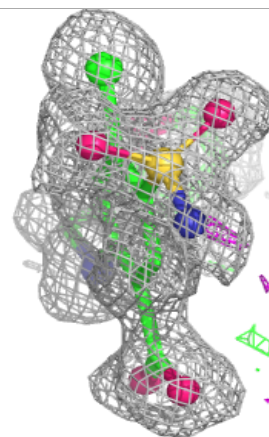
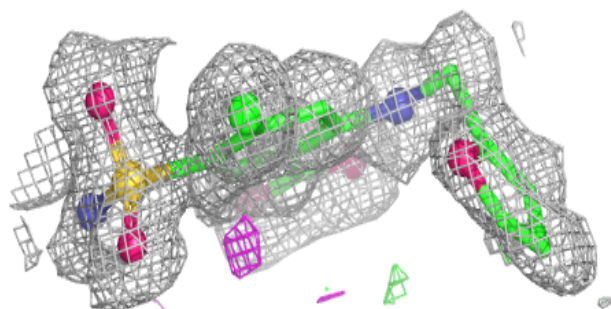
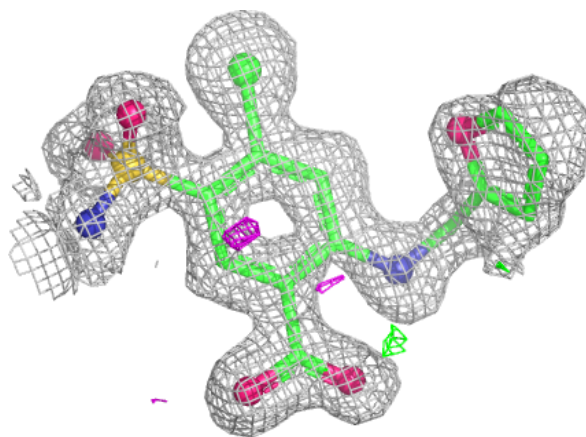
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	A	300	1/1	1.00	0.01	11,11,11,11	0
2	ZN	A	400	1/1	1.00	0.04	10,10,10,10	0
3	FUN	A	500	21/21	1.00	0.03	8,14,29,34	0
4	BE7	A	700	10/11	1.00	0.03	13,27,30,32	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.

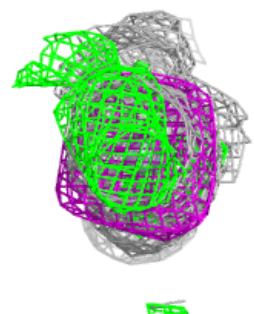
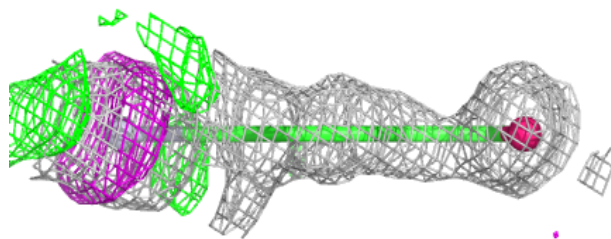
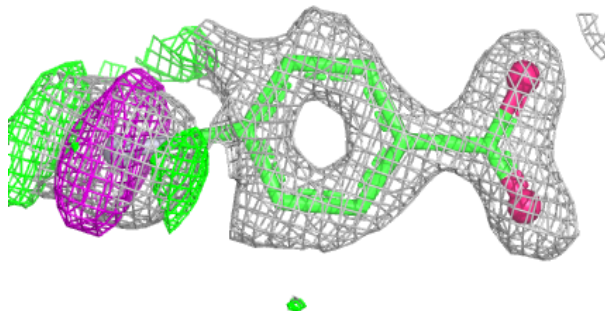
Electron density around FUN A 500:

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



Electron density around BE7 A 700:

$2mF_o-DF_c$ (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.