

wwPDB X-ray Structure Validation Summary Report (i)

Jan 2, 2024 – 11:46 pm GMT

PDB ID	:	5LLO
Title	:	Crystal structure of human carbonic anhydrase isozyme XII with 3-[(1S)-2,3-
		$\label{eq:linear} Dihydro-1H-inden-1-ylamino]-2,5,6-trifluoro-4-[(2-hy-droxyethyl)sulfonyl] benefative and the second s$
		zenesulfonamide
Authors	:	Smirnov, A.; Manakova, E.; Grazulis, S.
Deposited on	:	2016-07-28
Resolution	:	1.60 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:

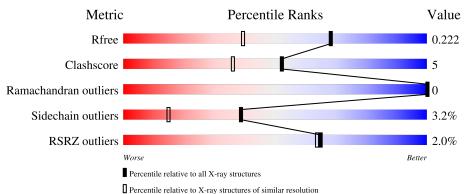
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as 541 be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
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: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665(1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	263	% 91%	7%	
1	В	263	88%	10%	
1	С	263	<u>6%</u> 85%	14%	•
1	D	263	90%	8%	••



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 9754 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	Atoms				ZeroOcc	AltConf	Trace	
1	А	260	Total	С	Ν	Ο	S	0	8	0
1	Л	200	2148	1363	365	413	7	0	8	0
1	В	261	Total	С	Ν	Ο	S	0	6	0
	I D	201	2153	1364	372	410	$\overline{7}$	0		
1	С	261	Total	С	Ν	0	S	0	F	0
		201	2138	1357	367	407	$\overline{7}$	0	5	0
1	Л	961	Total	С	Ν	0	S	0	4	0
		261	2130	1352	363	408	$\overline{7}$		4	0

• Molecule 1 is a protein called Carbonic anhydrase 12.

There are 4 discrepancies between the modelled and reference sequences:

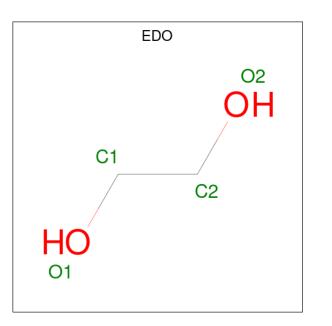
Chain	Residue	Modelled	Actual	Comment	Reference
А	1	MET	-	initiating methionine	UNP O43570
В	1	MET	-	initiating methionine	UNP O43570
С	1	MET	-	initiating methionine	UNP O43570
D	1	MET	-	initiating methionine	UNP O43570

• Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Zn 1 1	0	0
2	В	1	Total Zn 1 1	0	0
2	С	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

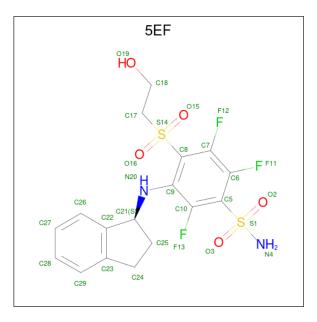
• Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
3	D	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0

• Molecule 4 is 3-[(1S)-2,3-dihydro-1H-inden-1-ylamino]-2,5,6-trifluoro-4-[(2-hydroxyethyl)sul fonyl]benzenesulfonamide (three-letter code: 5EF) (formula: $C_{17}H_{17}F_3N_2O_5S_2$).





Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	А	1	Total		F	Ν	0	\mathbf{S}	0	0
1	11	Ĩ	29	17	3	2	5	2	0	Ū
4	В	1	Total	С	\mathbf{F}	Ν	Ο	\mathbf{S}	0	0
4	D	1	29	17	3	2	5	2	0	0
4	С	1	Total	С	F	Ν	Ο	S	0	0
4	U	1	29	17	3	2	5	2	0	0
4	Л	1	Total	С	F	Ν	Ο	S	0	0
4	D		29	17	3	2	5	2		U

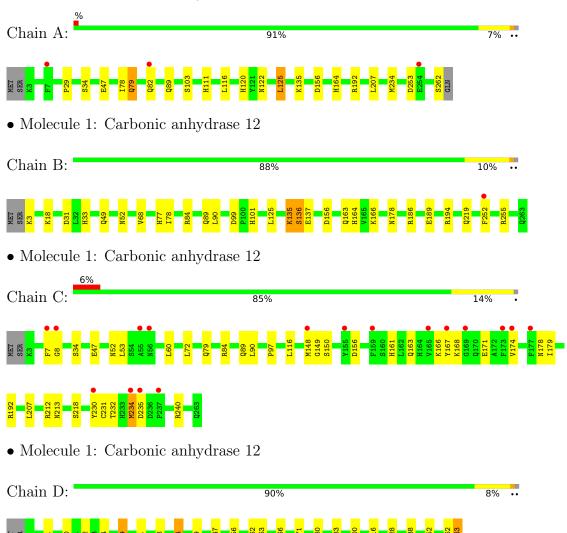
• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	268	Total O 268 268	0	0
5	В	286	Total O 286 286	0	0
5	С	223	Total O 223 223	0	0
5	D	276	Total O 276 276	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: Carbonic anhydrase 12



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	77.29Å 74.11Å 91.58Å	Depositor
a, b, c, α , β , γ	90.00° 108.95° 90.00°	Depositor
Resolution (Å)	73.10 - 1.60	Depositor
Resolution (A)	56.31 - 1.60	EDS
% Data completeness	98.8 (73.10-1.60)	Depositor
(in resolution range)	98.8 (56.31-1.60)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	0.05	Depositor
$< I/\sigma(I) > 1$	$3.34 (at 1.60 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
D D.	0.185 , 0.225	Depositor
R, R_{free}	0.182 , 0.222	DCC
R_{free} test set	12532 reflections (9.83%)	wwPDB-VP
Wilson B-factor $(Å^2)$	16.4	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.34 , 42.5	EDS
L-test for twinning ²	$ < L >=0.47, < L^2>=0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9754	wwPDB-VP
Average B, all atoms $(Å^2)$	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 52.05 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.1414e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

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: EDO, 5EF, 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		
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	1.12	0/2211	1.06	0/3011	
1	В	1.16	0/2216	1.08	0/3014	
1	С	1.05	0/2201	1.03	0/2994	
1	D	1.15	0/2193	1.06	0/2985	
All	All	1.12	0/8821	1.06	0/12004	

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 (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	2148	0	2028	14	0
1	В	2153	0	2042	25	0
1	С	2138	0	2031	32	0
1	D	2130	0	2015	22	0
2	А	1	0	0	0	0
2	В	1	0	0	0	0
2	С	1	0	0	0	0
2	D	1	0	0	0	0
3	А	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	В	4	0	6	0	0
3	D	4	0	6	1	0
4	А	29	0	0	0	0
4	В	29	0	0	0	0
4	С	29	0	0	0	0
4	D	29	0	0	0	0
5	А	268	0	0	3	0
5	В	286	0	0	9	0
5	С	223	0	0	5	0
5	D	276	0	0	4	0
All	All	9754	0	8134	92	0

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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.

The worst 5 of 92 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:D:163[A]:GLN:OE1	1:D:166[A]:LYS:CE	1.63	1.43
1:D:163[A]:GLN:OE1	1:D:166[A]:LYS:HE2	1.02	1.18
1:A:82:GLN:HG2	5:C:517:HOH:O	1.49	1.11
1:B:84[A]:ARG:HD2	5:B:441:HOH:O	1.52	1.08
1:C:148:MET:HB3	5:C:571:HOH:O	1.59	1.03

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	А	266/263~(101%)	262~(98%)	4 (2%)	0	100	100
1	В	265/263~(101%)	257 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured Allowed		Outliers	Percentiles	
1	\mathbf{C}	264/263~(100%)	256~(97%)	8 (3%)	0	100	100
1	D	263/263~(100%)	258~(98%)	5 (2%)	0	100	100
All	All	1058/1052~(101%)	1033 (98%)	25 (2%)	0	100	100

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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	А	240/235~(102%)	232~(97%)	8~(3%)	38	14	
1	В	239/235~(102%)	232~(97%)	7 (3%)	42	18	
1	С	238/235~(101%)	229~(96%)	9 (4%)	33	10	
1	D	237/235~(101%)	230~(97%)	7 (3%)	41	16	
All	All	954/940~(102%)	923~(97%)	31 (3%)	39	15	

5 of 31 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	В	219	GLN
1	D	156	ASP
1	С	84[B]	ARG
1	D	238	SER
1	D	49	GLN

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such side chains are listed below:

Mol	Chain	Res	Type
1	В	101	HIS
1	С	49	GLN
1	D	164	HIS
1	С	233	HIS
1	D	49	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 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	Bo	ond leng	ths	B	ond ang	gles
	ion Type Chain	Ullaili	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	D	302	-	3, 3, 3	0.39	0	2,2,2	1.04	0
4	$5\mathrm{EF}$	А	303	2	29,31,31	2.60	8 (27%)	39,48,48	3.10	11 (28%)
3	EDO	В	302	-	3,3,3	0.79	0	2,2,2	0.07	0
4	$5\mathrm{EF}$	В	303	2	29,31,31	2.55	9 (31%)	39,48,48	2.79	10 (25%)
4	$5\mathrm{EF}$	С	302	2	29,31,31	3.10	9 (31%)	39,48,48	2.85	13 (33%)
4	5EF	D	303	2	29,31,31	2.56	8 (27%)	39,48,48	2.77	15 (38%)
3	EDO	А	302	-	3,3,3	0.65	0	2,2,2	0.06	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
3	EDO	D	302	-	-	0/1/1/1	-
4	5EF	А	303	2	-	7/20/29/29	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	В	302	-	-	1/1/1/1	-
4	$5\mathrm{EF}$	В	303	2	-	14/20/29/29	0/3/3/3
4	$5\mathrm{EF}$	С	302	2	-	9/20/29/29	0/3/3/3
4	$5\mathrm{EF}$	D	303	2	-	10/20/29/29	0/3/3/3
3	EDO	А	302	-	-	0/1/1/1	-

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The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
4	D	303	$5\mathrm{EF}$	O16-S14	10.10	1.59	1.44
4	С	302	5EF	O16-S14	9.40	1.58	1.44
4	А	303	5EF	O16-S14	8.70	1.57	1.44
4	С	302	5EF	O15-S14	8.52	1.56	1.44
4	В	303	$5\mathrm{EF}$	O16-S14	7.64	1.55	1.44

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
4	А	303	5EF	O15-S14-O16	-14.39	101.61	118.44
4	В	303	5EF	O15-S14-O16	-12.25	104.12	118.44
4	D	303	5EF	O15-S14-O16	-10.98	105.60	118.44
4	С	302	5EF	O15-S14-O16	-8.46	108.54	118.44
4	С	302	5EF	O15-S14-C17	7.99	122.54	108.20

There are no chirality outliers.

5 of 41 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	А	303	5EF	C6-C5-S1-N4
4	А	303	5EF	C9-C8-S14-O15
4	В	303	5EF	C18-C17-S14-O16
4	В	303	5EF	C18-C17-S14-O15
4	В	303	5EF	C18-C17-S14-C8

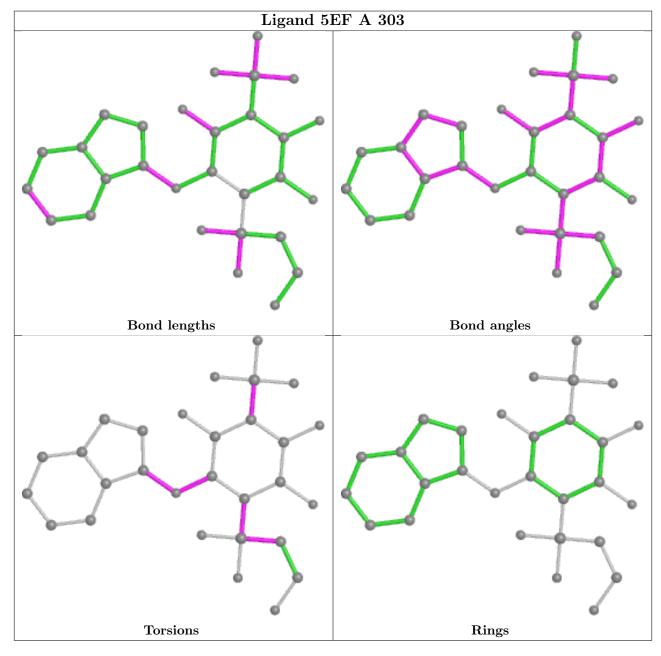
There are no ring outliers.

1 monomer is involved in 1 short contact:

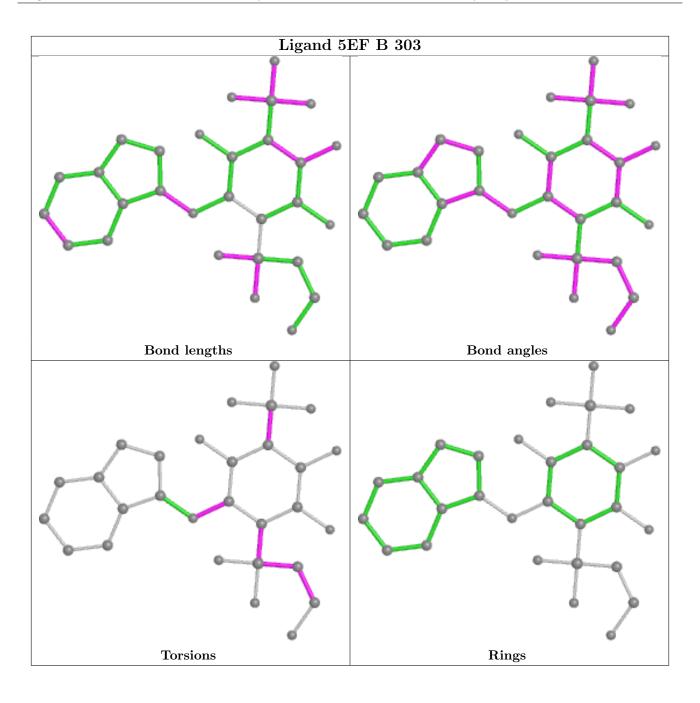
M	ol	Chain	Res	Type	Clashes	Symm-Clashes
3		D	302	EDO	1	0



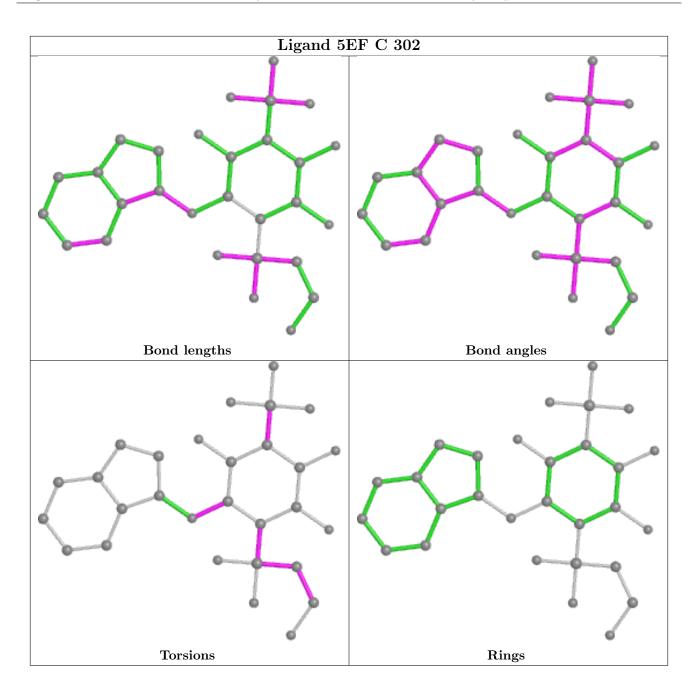
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 sufficient the outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



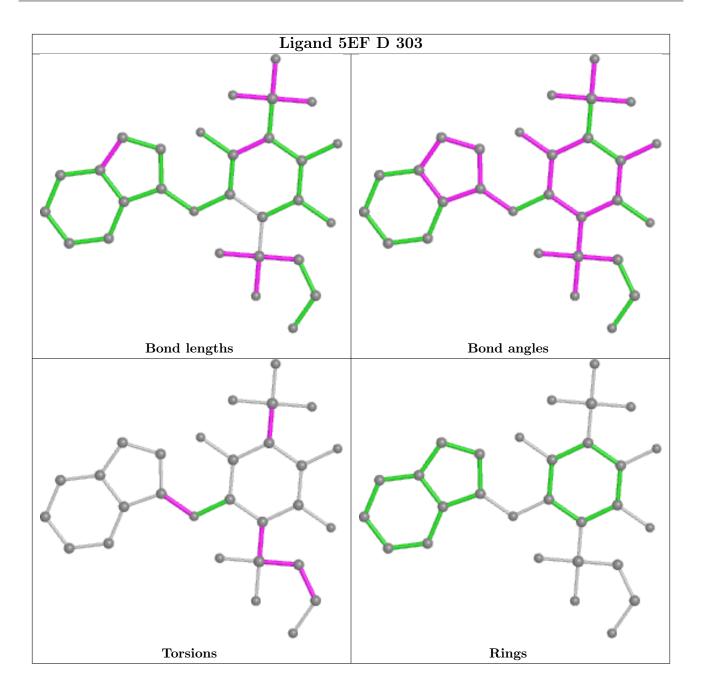












5.7 Other polymers (i)

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	$OWAB(Å^2)$	$\mathbf{Q} \! < \! 0.9$
1	А	260/263~(98%)	-0.10	3 (1%) 79 78	9, 17, 32, 41	0
1	В	261/263~(99%)	-0.14	1 (0%) 92 92	8, 14, 27, 38	0
1	С	261/263~(99%)	0.29	17 (6%) 18 17	9, 21, 41, 58	0
1	D	261/263~(99%)	-0.19	0 100 100	9, 15, 25, 53	0
All	All	1043/1052~(99%)	-0.03	21 (2%) 65 64	8, 16, 33, 58	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	\mathbf{Res}	Type	RSRZ
1	С	55	ALA	6.5
1	С	8	GLY	4.9
1	С	173	PHE	4.1
1	В	252	PHE	3.7
1	А	7	PHE	3.6

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)

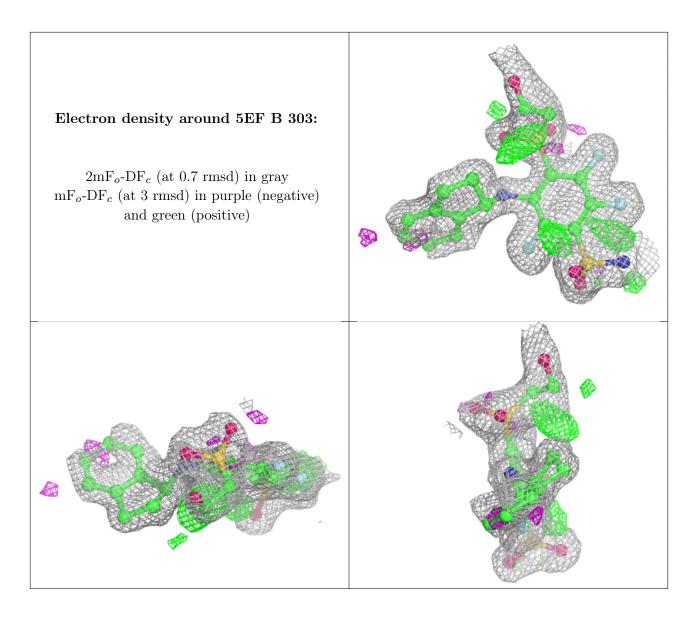
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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



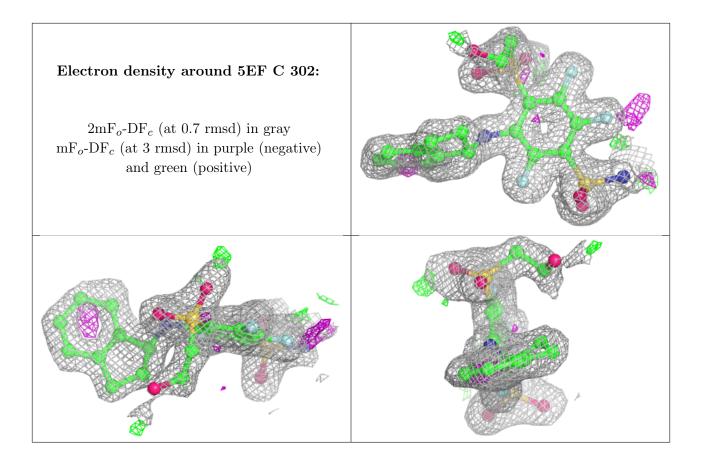
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	EDO	В	302	4/4	0.86	0.11	$29,\!30,\!31,\!32$	0
3	EDO	D	302	4/4	0.89	0.19	27,29,30,35	0
3	EDO	А	302	4/4	0.93	0.10	24,24,27,28	0
4	$5\mathrm{EF}$	В	303	29/29	0.95	0.11	$12,\!23,\!35,\!43$	0
4	5EF	С	302	29/29	0.95	0.13	16,22,28,39	0
4	5EF	D	303	29/29	0.96	0.11	13,21,33,45	0
4	5EF	А	303	29/29	0.97	0.09	12,24,34,40	0
2	ZN	С	301	1/1	0.99	0.07	$14,\!14,\!14,\!14$	0
2	ZN	А	301	1/1	1.00	0.07	10,10,10,10	0
2	ZN	D	301	1/1	1.00	0.08	10,10,10,10	0
2	ZN	В	301	1/1	1.00	0.08	9,9,9,9	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.

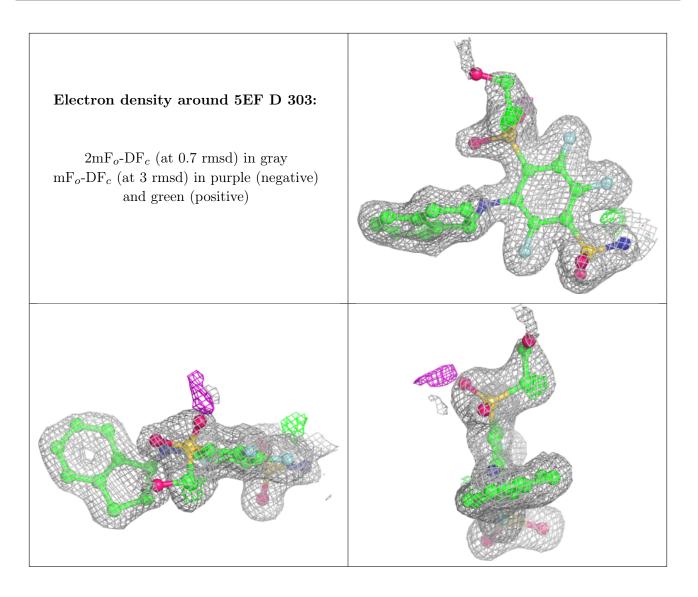




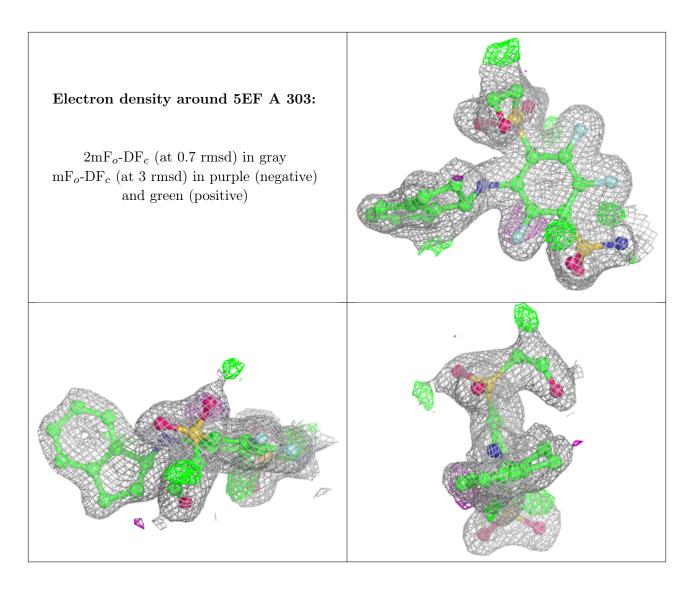












6.5 Other polymers (i)

There are no such residues in this entry.

