wwPDB X-ray Structure Validation Summary Report (i)

Oct 12, 2026 – 10:19 am BST

PDB ID	:	30FA / pdb_000030fa
Title	:	Crystal structure of carbonic anhydrase XII with methyl 4-benzylsulfanyl-3-sulfamoyl-benzoate
Authors	:	Paketuryte-Latve, V.; Smirnov, A.; Manakova, E.N.; Grazulis, S.
Deposited on	:	2026-04-22
Resolution	:	1.13 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 (i)) were used in the production of this report:

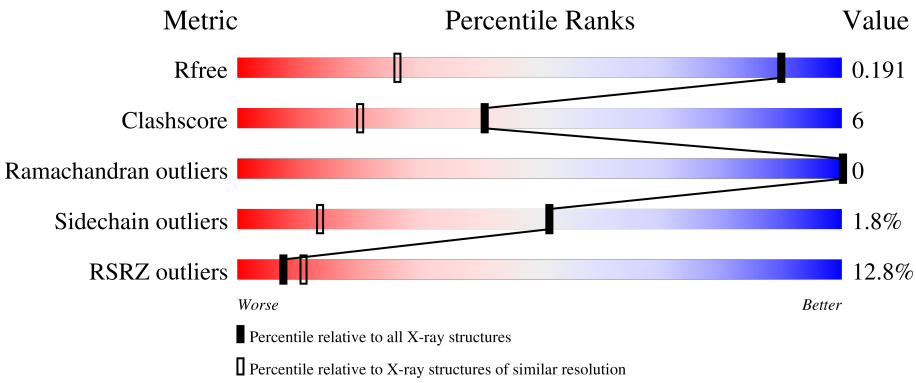
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 1.13 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	1985 (1.14-1.10)
Clashscore	190562	2021 (1.14-1.10)
Ramachandran outliers	187476	1978 (1.14-1.10)
Sidechain outliers	187428	1974 (1.14-1.10)
RSRZ outliers	180081	1984 (1.14-1.10)

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	263	9% 88% 11% •
1	B	263	8% 90% 9% ••
1	C	263	29% 87% 10% ••
1	D	263	5% 93% 6% •

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
4	PEG	A	303	-	-	X	-
4	PEG	B	303	-	-	X	-
5	EDO	A	304	-	-	X	-
5	EDO	D	304	-	X	-	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 9816 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 12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	261	2153	1368	366	412	7	0	8	0
1	B	261	2193	1391	372	422	8	0	12	0
1	C	261	2104	1341	355	400	8	0	4	0
1	D	260	2133	1353	363	410	7	0	5	0

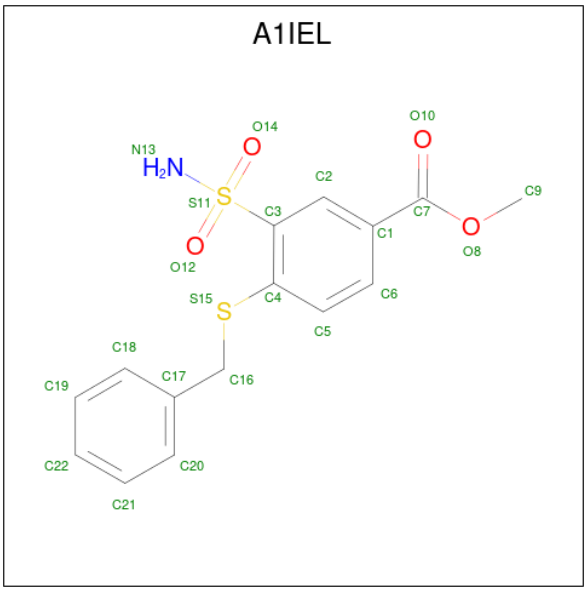
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP O43570
B	1	MET	-	initiating methionine	UNP O43570
C	1	MET	-	initiating methionine	UNP O43570
D	1	MET	-	initiating methionine	UNP O43570

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

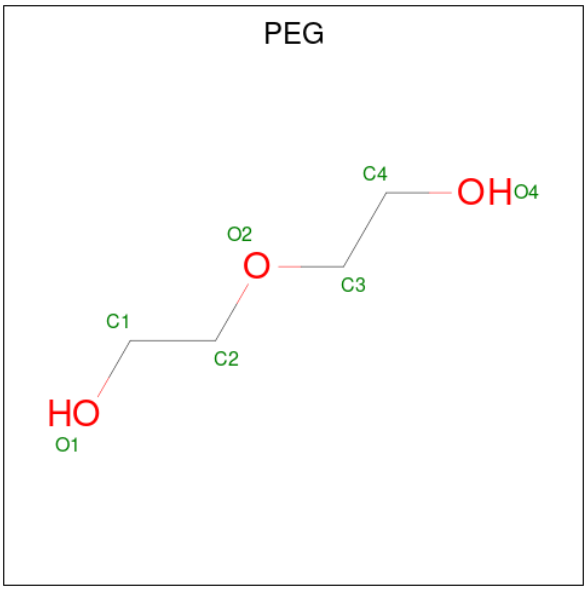
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0
2	D	1	1	1	0	0

- Molecule 3 is methyl 4-(phenylmethylsulfanyl)-3-sulfamoyl-benzoate (CCD ID: A1IEL) (formula: $C_{15}H_{15}NO_4S_2$) (labeled as "Ligand of Interest" by depositor).



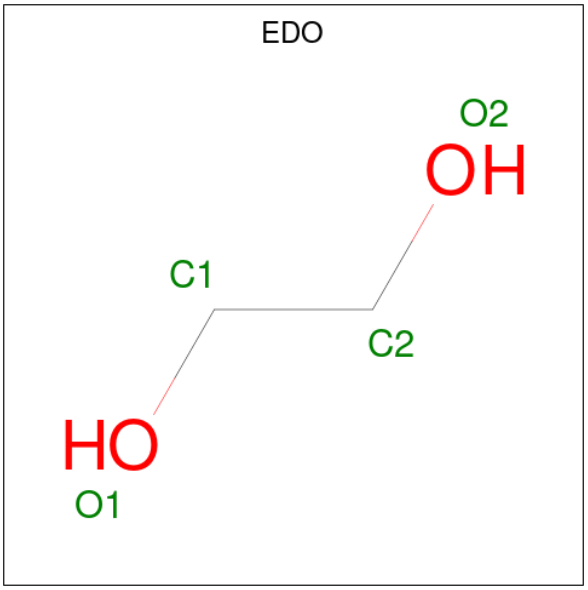
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	22	15	1	4	2	0	0
			Total	C	N	O	S		
3	B	1	22	15	1	4	2	0	0
			Total	C	N	O	S		
3	C	1	22	15	1	4	2	0	0
			Total	C	N	O	S		
3	D	1	22	15	1	4	2	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	7	4	3	0	0
			Total	C	O		
4	B	1	7	4	3	0	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	4	2	2	0	0
			Total	C	O		
5	A	1	4	2	2	0	0
			Total	C	O		
5	A	1	4	2	2	0	0
			Total	C	O		
5	A	1	4	2	2	0	0
			Total	C	O		
5	B	1	4	2	2	0	0
			Total	C	O		
5	D	1	4	2	2	0	0
			Total	C	O		
5	D	1	4	2	2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	281	281	281	0	0

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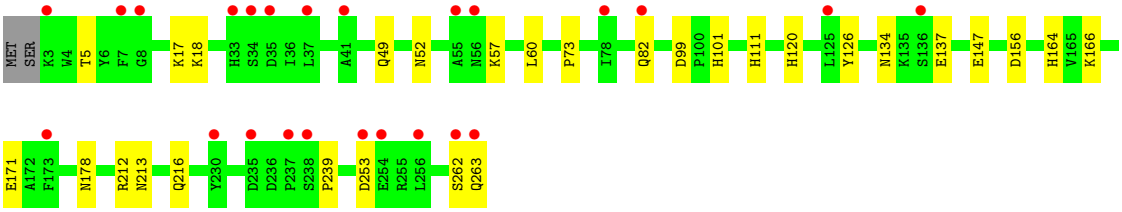
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	B	277	277	277	0	0
			Total	O		
6	C	229	229	229	0	0
			Total	O		
6	D	312	312	312	0	0

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.

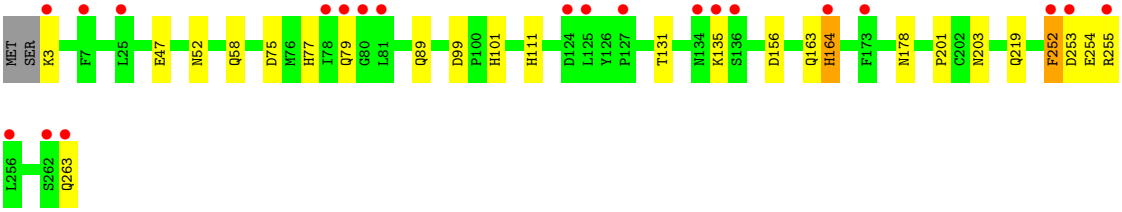
• Molecule 1: Carbonic anhydrase 12

Chain A:



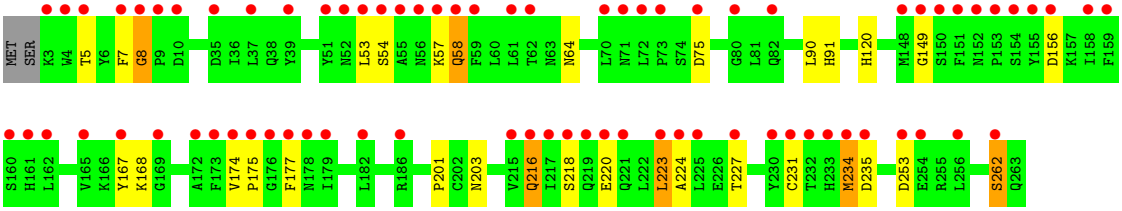
• Molecule 1: Carbonic anhydrase 12

Chain B:



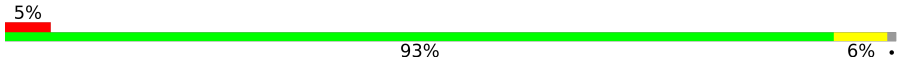
• Molecule 1: Carbonic anhydrase 12

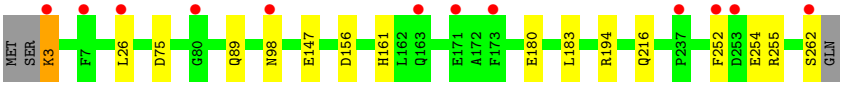
Chain C:



• Molecule 1: Carbonic anhydrase 12

Chain D:




MET
SER
K3
F7
L26
D75
G80
Q89
N98
E147
D156
H161
L162
Q163
E171
A172
F173
E180
L183
R194
Q216
P237
F252
D253
E254
R255
S262
GLN

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	77.22Å 74.07Å 91.65Å 90.00° 108.68° 90.00°	Depositor
Resolution (Å)	86.82 – 1.13 86.82 – 1.13	Depositor EDS
% Data completeness (in resolution range)	97.3 (86.82-1.13) 97.3 (86.82-1.13)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.50 (at 1.13Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.157 , 0.186 0.166 , 0.191	Depositor DCC
R_{free} test set	35287 reflections (9.68%)	wwPDB-VP
Wilson B-factor (Å^2)	11.7	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 47.1	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	9816	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 37.39 % 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 4.3185e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical 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.

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: ZN, PEG, EDO, A1IEL

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	0.88	0/2217	1.13	7/3019 (0.2%)
1	B	0.89	1/2258 (0.0%)	1.12	5/3074 (0.2%)
1	C	1.01	5/2168 (0.2%)	1.34	13/2954 (0.4%)
1	D	0.91	1/2197 (0.0%)	1.10	5/2992 (0.2%)
All	All	0.93	7/8840 (0.1%)	1.18	30/12039 (0.2%)

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	B	0	1
1	C	0	1
All	All	0	2

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	223	LEU	C-N	11.97	1.50	1.34
1	D	161	HIS	CE1-NE2	9.48	1.42	1.32
1	B	77	HIS	CE1-NE2	5.91	1.38	1.32
1	C	227	THR	CB-OG1	5.83	1.53	1.43
1	C	262	SER	CA-CB	-5.75	1.44	1.53

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	75	ASP	CA-CB-CG	11.69	124.29	112.60
1	C	220	GLU	CB-CA-C	-11.13	92.09	110.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	253	ASP	CA-CB-CG	8.63	121.23	112.60
1	D	252	PHE	CA-CB-CG	-7.50	106.30	113.80
1	C	177	PHE	CA-CB-CG	7.38	121.18	113.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	252[B]	PHE	Mainchain
1	C	223	LEU	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	A	2153	0	2033	25	0
1	B	2193	0	2061	45	0
1	C	2104	0	1983	16	0
1	D	2133	0	2005	9	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	22	0	0	0	0
3	B	22	0	0	0	0
3	C	22	0	0	0	0
3	D	22	0	0	0	0
4	A	7	0	10	5	0
4	B	7	0	10	12	0
5	A	16	0	23	9	0
5	B	4	0	6	0	0
5	D	8	0	12	3	0
6	A	281	0	0	10	0
6	B	277	0	0	14	0
6	C	229	0	0	2	0
6	D	312	0	0	6	0
All	All	9816	0	8143	109	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 109 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)
5:D:304:EDO:C2	5:D:304:EDO:O2	1.64	1.40
5:A:307:EDO:H11	6:A:419:HOH:O	1.42	1.18
1:B:252[B]:PHE:CZ	1:B:255:ARG:HB3	1.83	1.12
1:B:252[B]:PHE:CZ	1:B:255:ARG:CB	2.48	0.97
1:B:252[A]:PHE:CZ	1:B:255:ARG:HB3	2.02	0.94

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	267/263 (102%)	262 (98%)	5 (2%)	0	100	100
1	B	271/263 (103%)	263 (97%)	8 (3%)	0	100	100
1	C	263/263 (100%)	255 (97%)	8 (3%)	0	100	100
1	D	263/263 (100%)	256 (97%)	7 (3%)	0	100	100
All	All	1064/1052 (101%)	1036 (97%)	28 (3%)	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	A	239/235 (102%)	234 (98%)	5 (2%)	47	10
1	B	244/235 (104%)	240 (98%)	4 (2%)	55	18
1	C	232/235 (99%)	228 (98%)	4 (2%)	53	16
1	D	237/235 (101%)	231 (98%)	6 (2%)	42	7
All	All	952/940 (101%)	933 (98%)	19 (2%)	51	10

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	89[A]	GLN
1	D	254	GLU
1	D	262	SER
1	D	156	ASP
1	B	156	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	13	ASN
1	C	58	GLN
1	C	56	ASN
1	C	79	GLN
1	A	213	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 17 ligands modelled in this entry, 4 are monoatomic - leaving 13 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	PEG	B	303	-	6,6,6	0.60	0	5,5,5	0.95	0
5	EDO	A	304	-	3,3,3	0.23	0	2,2,2	2.29	1 (50%)
3	A1IEL	D	302	2	23,23,23	2.69	13 (56%)	31,32,32	2.77	11 (35%)
3	A1IEL	A	302	2	23,23,23	1.36	4 (17%)	31,32,32	1.43	4 (12%)
5	EDO	A	307	-	3,3,3	1.80	2 (66%)	2,2,2	0.06	0
5	EDO	A	305	-	3,3,3	0.22	0	2,2,2	0.23	0
3	A1IEL	C	302	2	23,23,23	1.35	4 (17%)	31,32,32	1.54	7 (22%)
3	A1IEL	B	302	2	23,23,23	1.83	4 (17%)	31,32,32	1.98	8 (25%)
5	EDO	D	304	-	3,3,3	3.19	2 (66%)	2,2,2	1.55	0
5	EDO	D	303	-	3,3,3	0.82	0	2,2,2	0.69	0
5	EDO	B	304	-	3,3,3	0.89	0	2,2,2	0.01	0
5	EDO	A	306	-	3,3,3	0.32	0	2,2,2	0.58	0
4	PEG	A	303	-	6,6,6	0.30	0	5,5,5	0.38	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
4	PEG	B	303	-	-	4/4/4/4	-
5	EDO	A	304	-	-	0/1/1/1	-
3	A1IEL	D	302	2	-	3/17/17/17	0/2/2/2
3	A1IEL	A	302	2	-	2/17/17/17	0/2/2/2
5	EDO	A	307	-	-	0/1/1/1	-
5	EDO	A	305	-	-	1/1/1/1	-
3	A1IEL	C	302	2	-	3/17/17/17	0/2/2/2
3	A1IEL	B	302	2	-	2/17/17/17	0/2/2/2
5	EDO	D	304	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	D	303	-	-	0/1/1/1	-
5	EDO	B	304	-	-	1/1/1/1	-
5	EDO	A	306	-	-	0/1/1/1	-
4	PEG	A	303	-	-	4/4/4/4	-

The worst 5 of 29 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	302	A1IEL	C3-S11	6.89	1.86	1.77
3	B	302	A1IEL	C4-C3	5.34	1.49	1.40
3	D	302	A1IEL	C4-S15	-4.38	1.70	1.77
5	D	304	EDO	O2-C2	4.34	1.64	1.42
3	D	302	A1IEL	C1-C7	-4.25	1.40	1.50

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	302	A1IEL	O8-C7-C1	8.08	124.59	112.34
3	D	302	A1IEL	C4-C3-S11	-6.60	117.22	122.42
3	D	302	A1IEL	C9-O8-C7	6.26	127.89	115.83
3	B	302	A1IEL	C4-C3-S11	-5.62	117.99	122.42
3	C	302	A1IEL	C9-O8-C7	4.79	125.07	115.83

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	302	A1IEL	C4-C3-S11-O14
3	B	302	A1IEL	C4-C3-S11-O14
3	C	302	A1IEL	C4-C3-S11-O14
3	D	302	A1IEL	C4-C3-S11-O14
4	B	303	PEG	C4-C3-O2-C2

There are no ring outliers.

5 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	303	PEG	12	0
5	A	304	EDO	7	0
5	A	307	EDO	2	0
5	D	304	EDO	3	0

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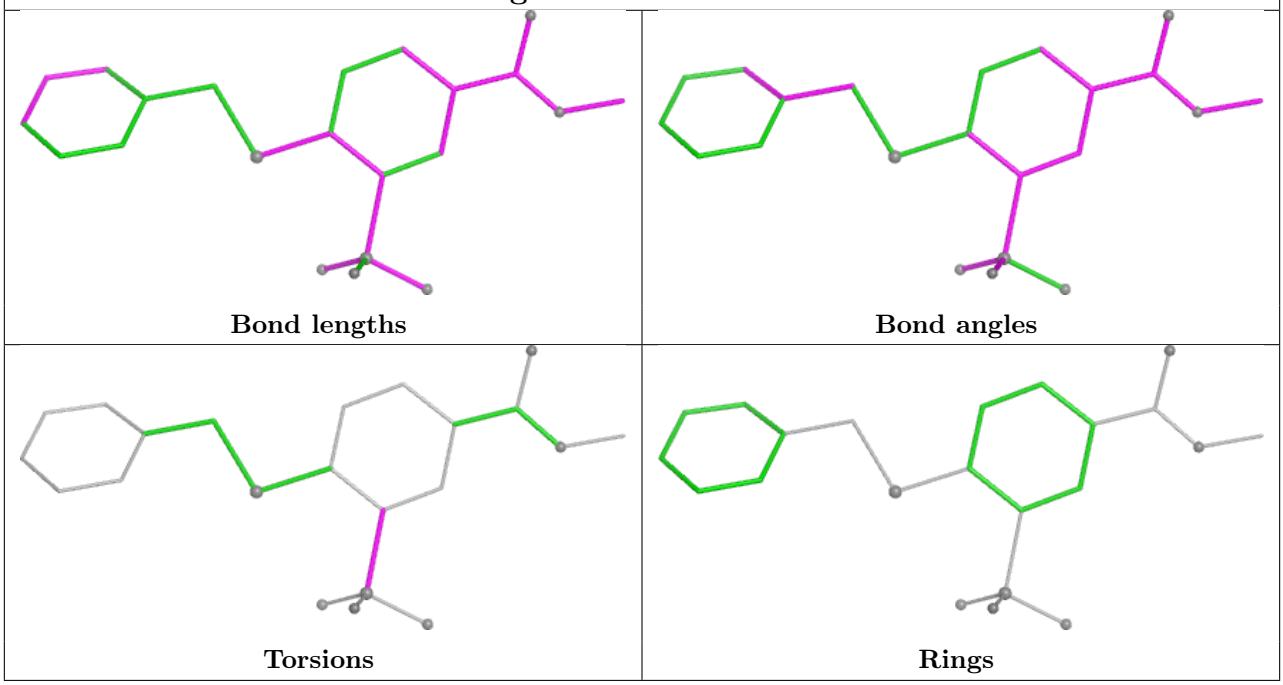
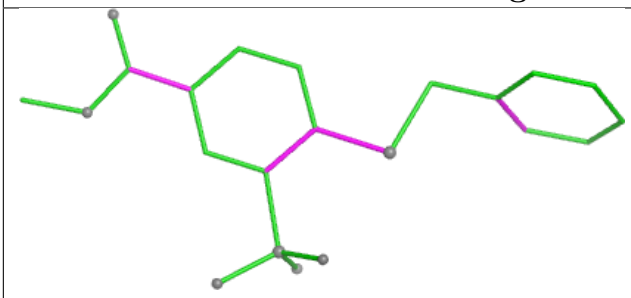
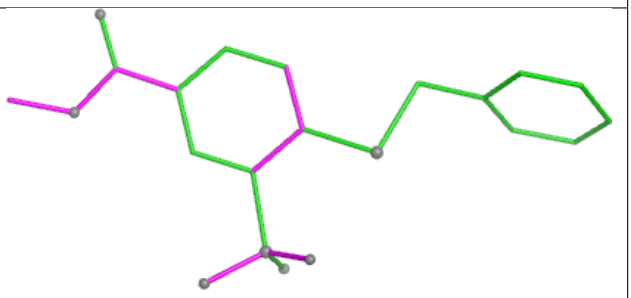
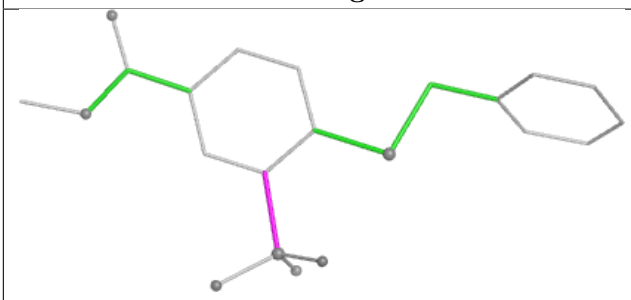
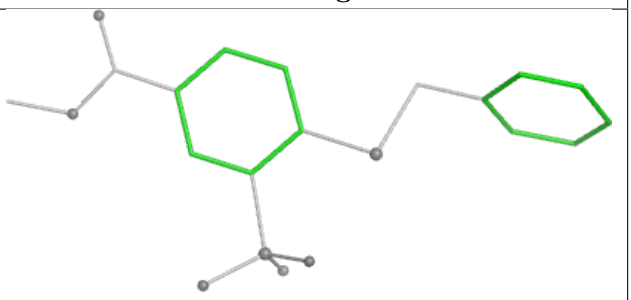
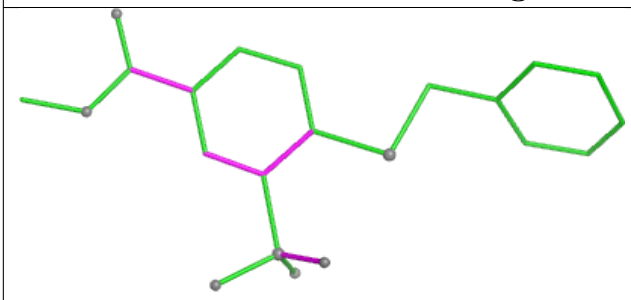
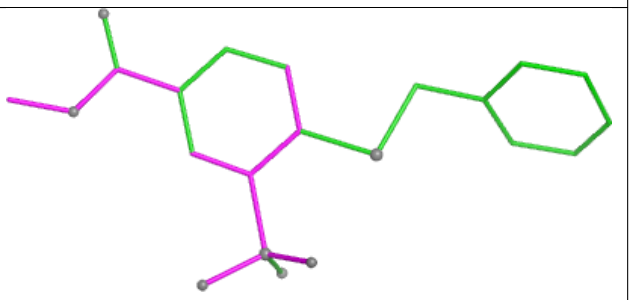
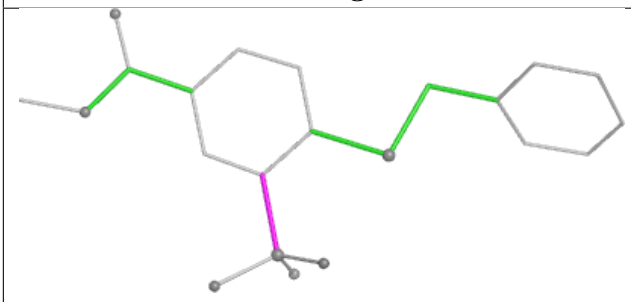
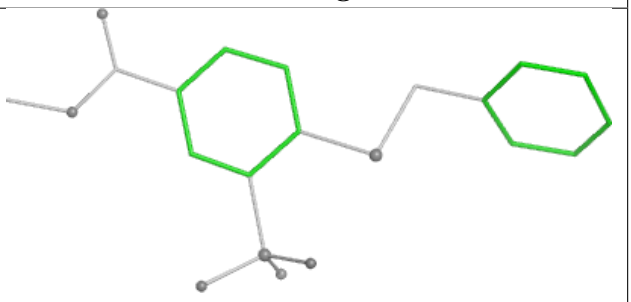
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	PEG	5	0

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.

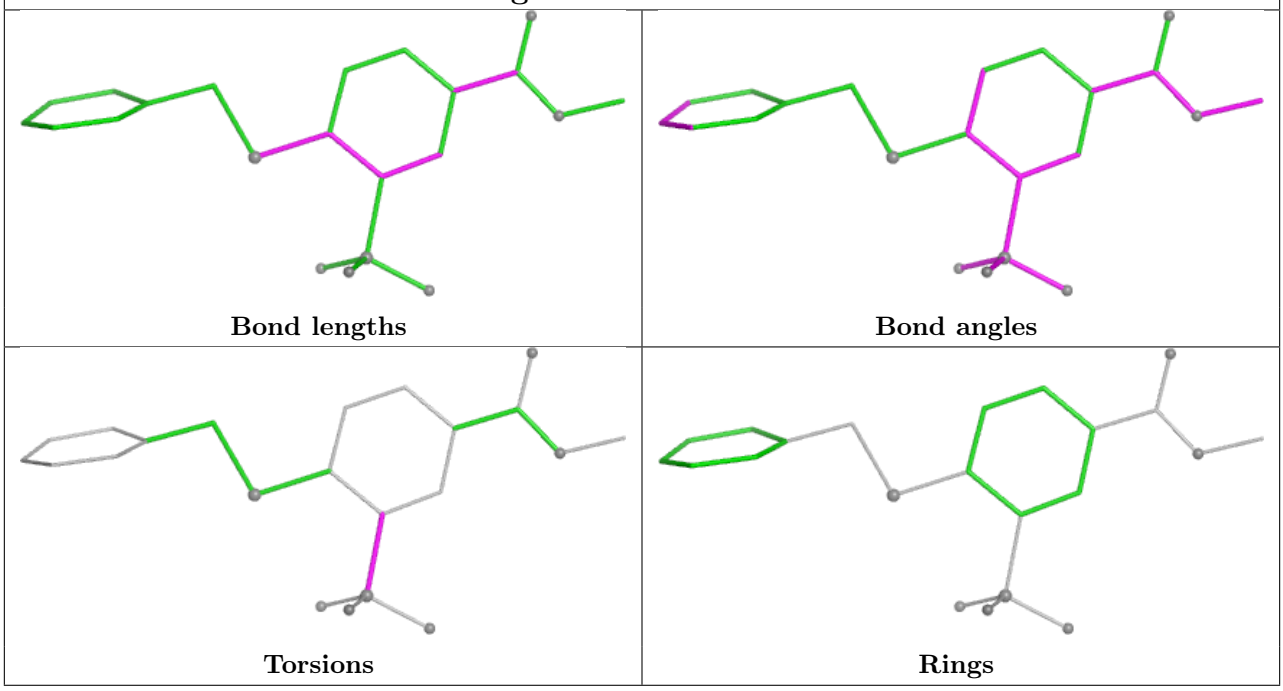
Ligand A1IEL D 302



Ligand A1IEL A 302	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand A1IEL C 302	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand A1IEL B 302



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	261/263 (99%)	0.60	24 (9%)	14	20	6, 16, 34, 60	8 (3%)
1	B	261/263 (99%)	0.41	21 (8%)	18	25	5, 13, 26, 66	12 (4%)
1	C	261/263 (99%)	1.53	76 (29%)	1	3	7, 21, 51, 178	4 (1%)
1	D	260/263 (98%)	0.42	12 (4%)	37	43	6, 14, 29, 53	5 (1%)
All	All	1043/1052 (99%)	0.74	133 (12%)	7	11	5, 15, 37, 178	29 (2%)

The worst 5 of 133 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	7	PHE	12.0
1	C	151	PHE	11.0
1	C	224	ALA	8.3
1	C	8	GLY	7.5
1	C	223	LEU	7.4

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)

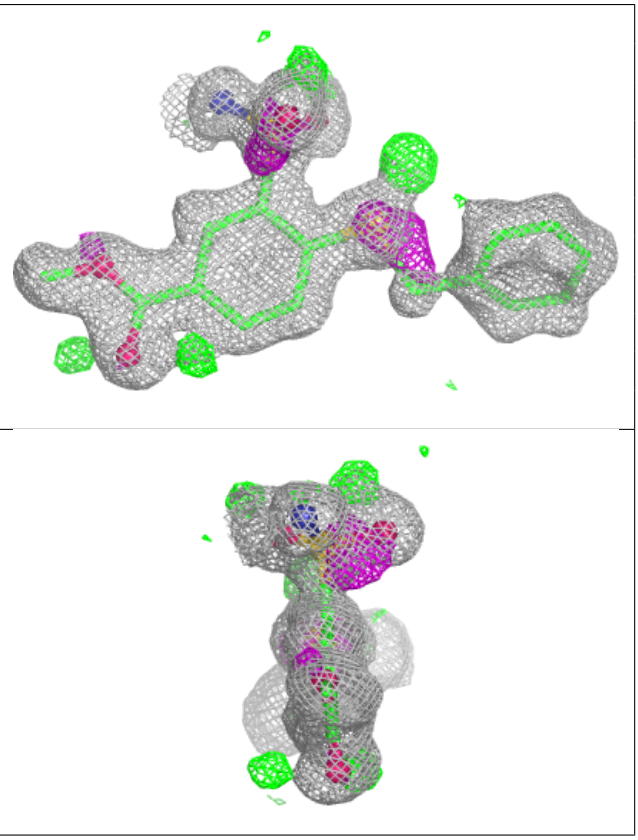
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	PEG	B	303	7/7	0.84	0.13	20,25,35,36	0
5	EDO	A	306	4/4	0.87	0.14	21,22,24,24	0
4	PEG	A	303	7/7	0.90	0.11	22,31,33,38	0
5	EDO	D	303	4/4	0.91	0.11	19,21,22,26	0
5	EDO	B	304	4/4	0.92	0.11	19,22,26,27	0
5	EDO	D	304	4/4	0.92	0.11	17,19,19,26	0
5	EDO	A	305	4/4	0.93	0.10	22,24,24,25	0
5	EDO	A	304	4/4	0.94	0.10	24,27,29,30	0
5	EDO	A	307	4/4	0.95	0.11	13,15,19,19	0
3	A1IEL	D	302	22/22	0.98	0.08	10,16,23,25	0
3	A1IEL	B	302	22/22	0.99	0.06	9,14,25,27	0
3	A1IEL	C	302	22/22	0.99	0.06	11,17,26,28	0
3	A1IEL	A	302	22/22	0.99	0.05	10,14,26,28	0
2	ZN	B	301	1/1	1.00	0.01	7,7,7,7	0
2	ZN	C	301	1/1	1.00	0.01	11,11,11,11	0
2	ZN	D	301	1/1	1.00	0.01	8,8,8,8	0
2	ZN	A	301	1/1	1.00	0.01	8,8,8,8	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.

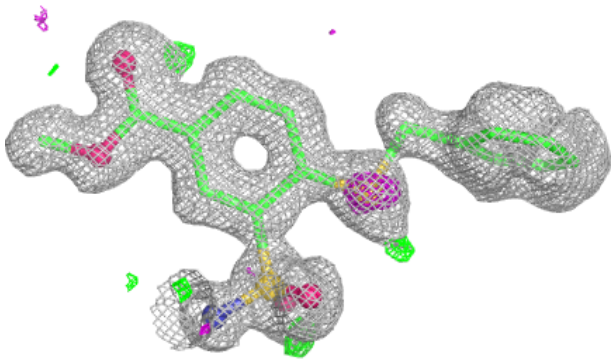
Electron density around A1IEL D 302:

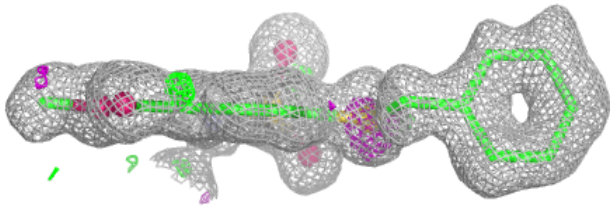
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

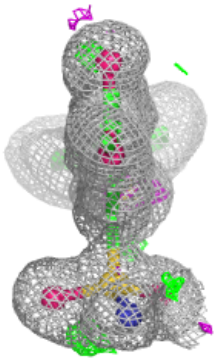


Electron density around A1IEL B 302:

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







Electron density around A1IEL C 302:

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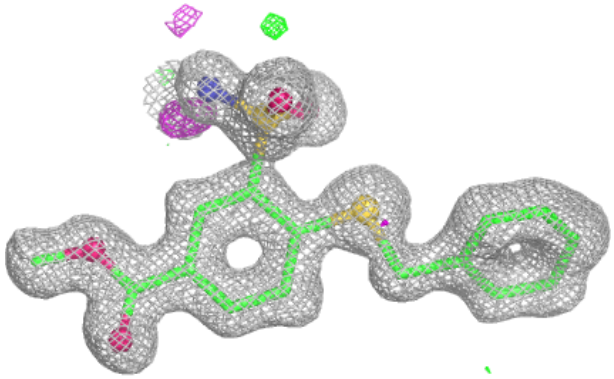
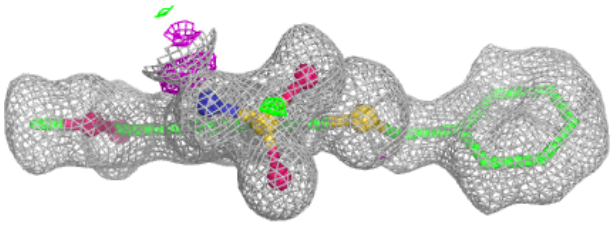
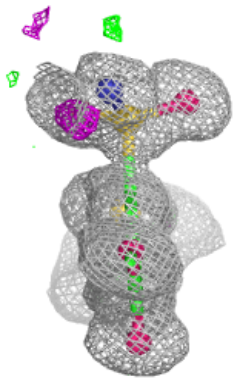
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

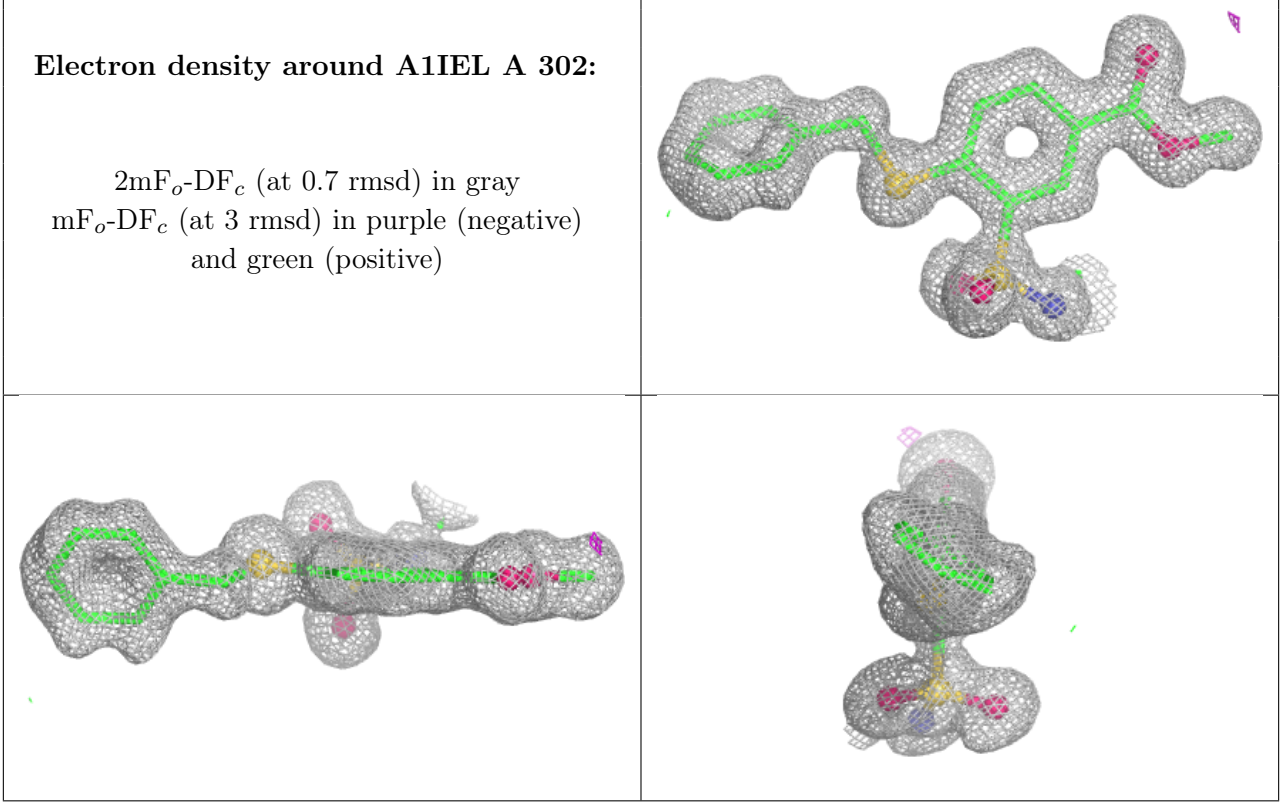






Electron density around A1IEL A 302:

$2mF_o$-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.