



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 12:15 PM UTC

PDB ID : 7OPI / pdb_00007opi
Title : Structure of a minimal SF3B core in complex with the inactive modulator spliceostatin E (form I)
Authors : Cretu, C.; Pena, V.
Deposited on : 2021-05-31
Resolution : 3.10 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 ⓘ 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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

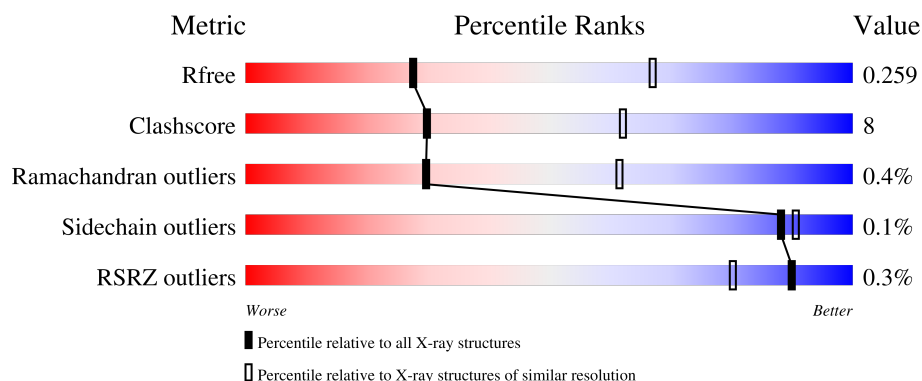
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

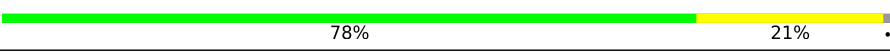
The reported resolution of this entry is 3.10 Å.

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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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 $\leq 5\%$. 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	899	 74% 22% .
2	B	86	 63% 13% 24%
3	C	852	 78% 21% .
4	D	108	 73% 13% 14%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14801 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 Splicing factor 3B subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	863	Total	C	N	O	S	0	0	0
			6823	4338	1166	1288	31			

There are 43 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	GLY	-	expression tag	UNP Q921M3
A	-8	ALA	-	expression tag	UNP Q921M3
A	-7	GLU	-	expression tag	UNP Q921M3
A	-6	PHE	-	expression tag	UNP Q921M3
A	-5	LYS	-	expression tag	UNP Q921M3
A	-4	GLY	-	expression tag	UNP Q921M3
A	-3	LEU	-	expression tag	UNP Q921M3
A	-2	ARG	-	expression tag	UNP Q921M3
A	-1	ARG	-	expression tag	UNP Q921M3
A	0	HIS	-	expression tag	UNP Q921M3
A	761	GLY	-	linker	UNP Q921M3
A	762	GLY	-	linker	UNP Q921M3
A	763	ASN	-	linker	UNP Q921M3
A	764	GLY	-	linker	UNP Q921M3
A	765	ASN	-	linker	UNP Q921M3
A	766	SER	-	linker	UNP Q921M3
A	767	GLY	-	linker	UNP Q921M3
A	?	-	GLU	deletion	UNP Q15393
A	?	-	ASP	deletion	UNP Q15393
A	?	-	PRO	deletion	UNP Q15393
A	?	-	THR	deletion	UNP Q15393
A	?	-	GLY	deletion	UNP Q15393
A	?	-	ASN	deletion	UNP Q15393
A	?	-	LYS	deletion	UNP Q15393
A	?	-	ALA	deletion	UNP Q15393
A	?	-	LEU	deletion	UNP Q15393
A	?	-	TRP	deletion	UNP Q15393

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ASP	deletion	UNP Q15393
A	?	-	ARG	deletion	UNP Q15393
A	?	-	GLY	deletion	UNP Q15393
A	?	-	LEU	deletion	UNP Q15393
A	?	-	LEU	deletion	UNP Q15393
A	?	-	ASN	deletion	UNP Q15393
A	?	-	GLY	deletion	UNP Q15393
A	?	-	ALA	deletion	UNP Q15393
A	1200	ASP	-	expression tag	UNP Q15393
A	1201	TYR	-	expression tag	UNP Q15393
A	1202	LYS	-	expression tag	UNP Q15393
A	1203	ASP	-	expression tag	UNP Q15393
A	1204	ASP	-	expression tag	UNP Q15393
A	1205	ASP	-	expression tag	UNP Q15393
A	1206	ASP	-	expression tag	UNP Q15393
A	1207	LYS	-	expression tag	UNP Q15393

- Molecule 2 is a protein called Splicing factor 3B subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	65	Total	C	N	O	S	0	0	0
			535	340	93	97	5			

- Molecule 3 is a protein called Splicing factor 3B subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	842	Total	C	N	O	S	0	0	0
			6707	4307	1152	1209	39			

- Molecule 4 is a protein called PHD finger-like domain-containing protein 5A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	93	Total	C	N	O	S	0	0	0
			700	429	123	135	13			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-9	GLY	-	expression tag	UNP Q7RTV0
D	-8	PRO	-	expression tag	UNP Q7RTV0
D	-7	LEU	-	expression tag	UNP Q7RTV0
D	-6	GLY	-	expression tag	UNP Q7RTV0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	SER	-	expression tag	UNP Q7RTV0
D	-4	PRO	-	expression tag	UNP Q7RTV0
D	-3	GLY	-	expression tag	UNP Q7RTV0
D	-2	SER	-	expression tag	UNP Q7RTV0
D	-1	ARG	-	expression tag	UNP Q7RTV0
D	0	ALA	-	expression tag	UNP Q7RTV0

- 071
-
- Chemical structure diagram showing a complex molecule with multiple stereocenters and functional groups. The structure includes a central chain with a nitrogen atom (N) and a carbonyl group (C19, O). The molecule features two substituted benzene rings. The top ring is a 1,3-diene derivative with a carboxylate group (O1, O2) and a methoxy group (C13, O3). The bottom ring is a 1,3-diene derivative with a carboxylate group (O5, O4) and a methoxy group (C23, O2). Stereochemistry is indicated by wedges and dashes at various chiral centers.

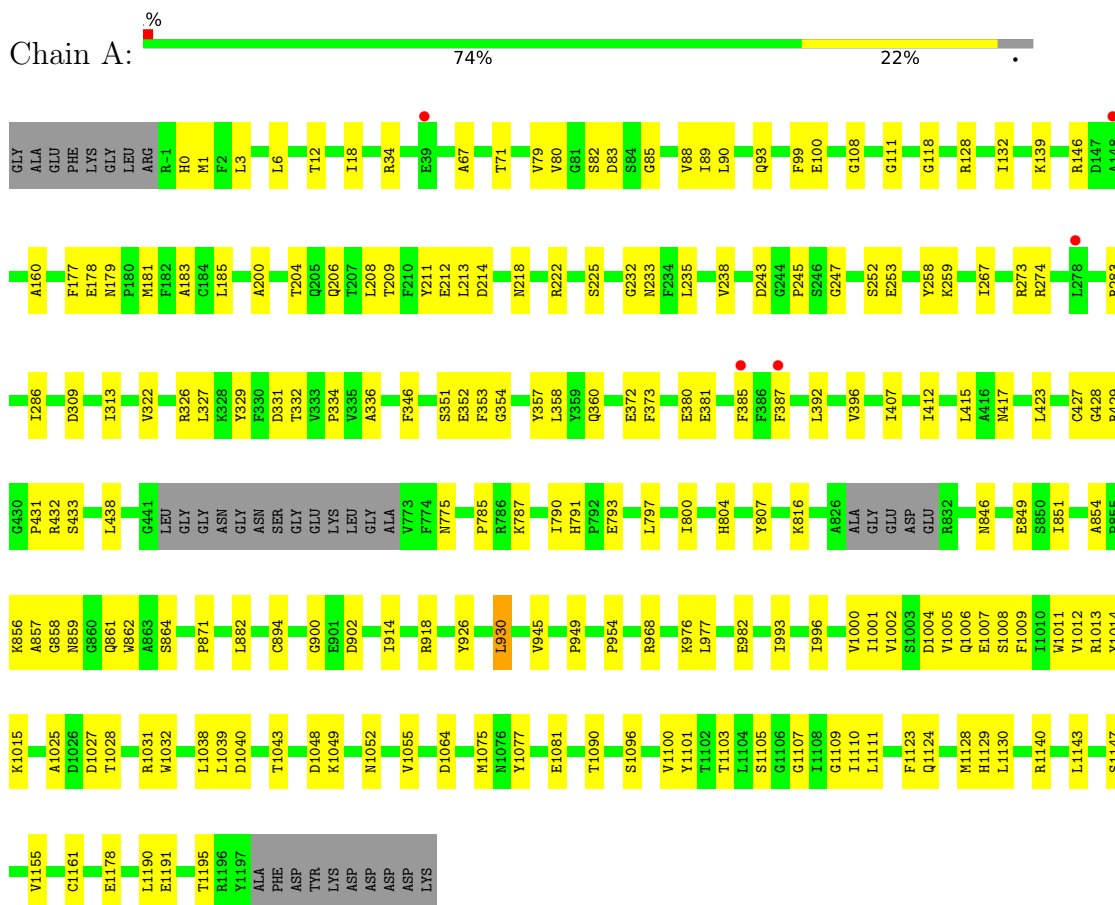
- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	3	Total Zn 3 3	0	0

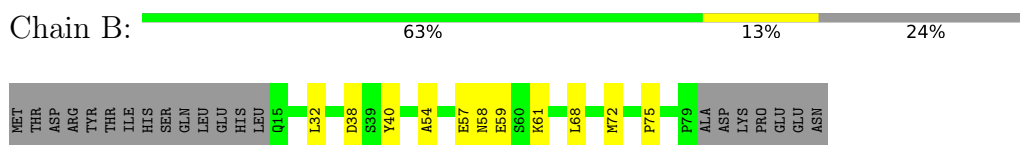
3 Residue-property plots

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.

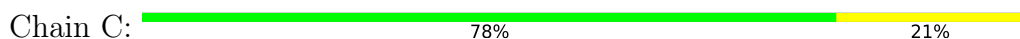
- Molecule 1: Splicing factor 3B subunit 3

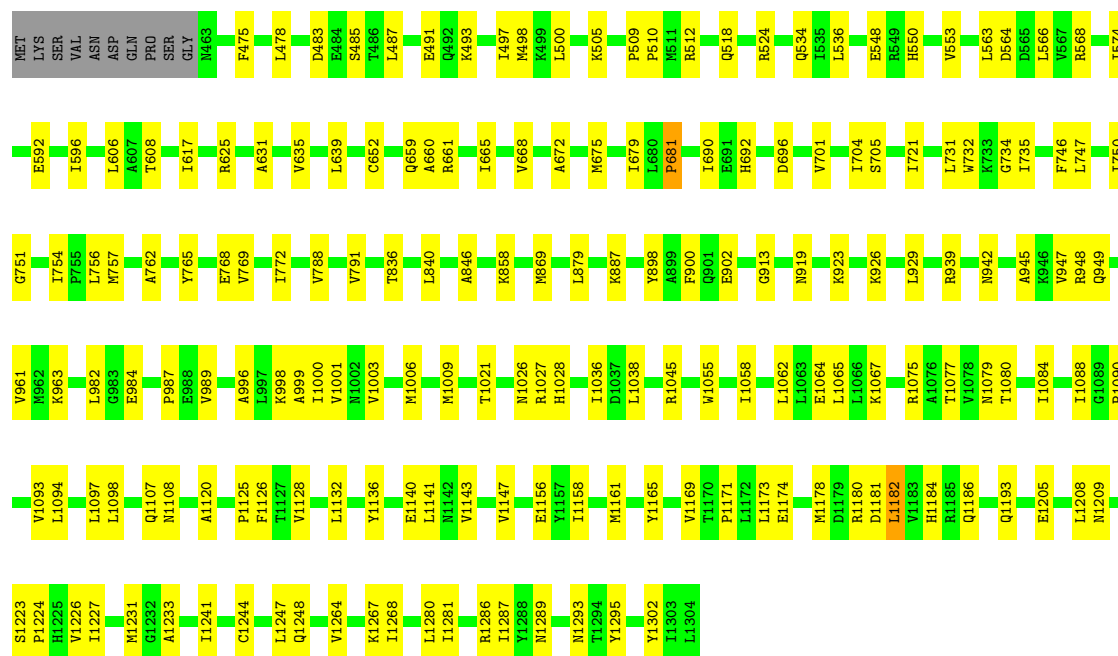


- Molecule 2: Splicing factor 3B subunit 5



- Molecule 3: Splicing factor 3B subunit 1





● Molecule 4: PHD finger-like domain-containing protein 5A

Chain D: 73% 13% 14%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.53Å 108.28Å 253.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.43 – 3.10 48.43 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.43-3.10) 100.0 (48.43-3.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.228 , 0.269 (Not available) , 0.259	Depositor DCC
R_{free} test set	3041 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	114.4	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 61.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14801	wwPDB-VP
Average B, all atoms (Å ²)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 07I, 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	0.12	0/6971	0.35	0/9452
2	B	0.14	0/551	0.32	0/744
3	C	0.13	0/6835	0.31	0/9256
4	D	0.11	0/709	0.31	0/952
All	All	0.13	0/15066	0.33	0/20404

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	A	6823	0	6744	118	0
2	B	535	0	504	9	0
3	C	6707	0	6915	125	0
4	D	700	0	682	9	0
5	C	33	0	0	0	0
6	D	3	0	0	0	0
All	All	14801	0	14845	252	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 8.

The worst 5 of 252 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)
3:C:487:LEU:HD12	3:C:491:GLU:HG2	1.50	0.93
3:C:553:VAL:HG23	3:C:596:ILE:HG23	1.59	0.84
3:C:1289:ASN:HB3	3:C:1295:TYR:H	1.42	0.83
3:C:1094:LEU:HD11	3:C:1128:VAL:HA	1.62	0.80
1:A:1011:TRP:HB2	1:A:1025:ALA:HB3	1.68	0.75

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	857/899 (95%)	813 (95%)	40 (5%)	4 (0%)	24	57
2	B	63/86 (73%)	59 (94%)	4 (6%)	0	100	100
3	C	840/852 (99%)	814 (97%)	23 (3%)	3 (0%)	30	61
4	D	91/108 (84%)	86 (94%)	5 (6%)	0	100	100
All	All	1851/1945 (95%)	1772 (96%)	72 (4%)	7 (0%)	30	61

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	1182	LEU
1	A	417	ASN
1	A	1009	PHE
1	A	900	GLY
3	C	564	ASP

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	747/774 (96%)	747 (100%)	0	100	100
2	B	57/77 (74%)	57 (100%)	0	100	100
3	C	728/737 (99%)	727 (100%)	1 (0%)	88	90
4	D	80/90 (89%)	80 (100%)	0	100	100
All	All	1612/1678 (96%)	1611 (100%)	1 (0%)	88	90

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	982	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	518	GLN
3	C	820	GLN
4	D	55	GLN
4	D	14	GLN
4	D	50	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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, 3 are monoatomic - leaving 1 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	$\# Z > 2$
5	07I	C	1401	-	34,34,34	0.18	0	40,46,46	1.51	6 (15%)

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
5	07I	C	1401	-	-	3/23/51/51	0/2/2/2

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	1401	07I	C6-C7-C8	-4.69	117.22	127.50
5	C	1401	07I	C2-C3-C4	4.38	117.92	109.53
5	C	1401	07I	O-C4-C3	4.22	118.45	109.31
5	C	1401	07I	C11-C10-C8	-3.41	118.99	126.32
5	C	1401	07I	O-C5-C6	2.12	117.48	109.98

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	1401	07I	C1-C5-C6-C7

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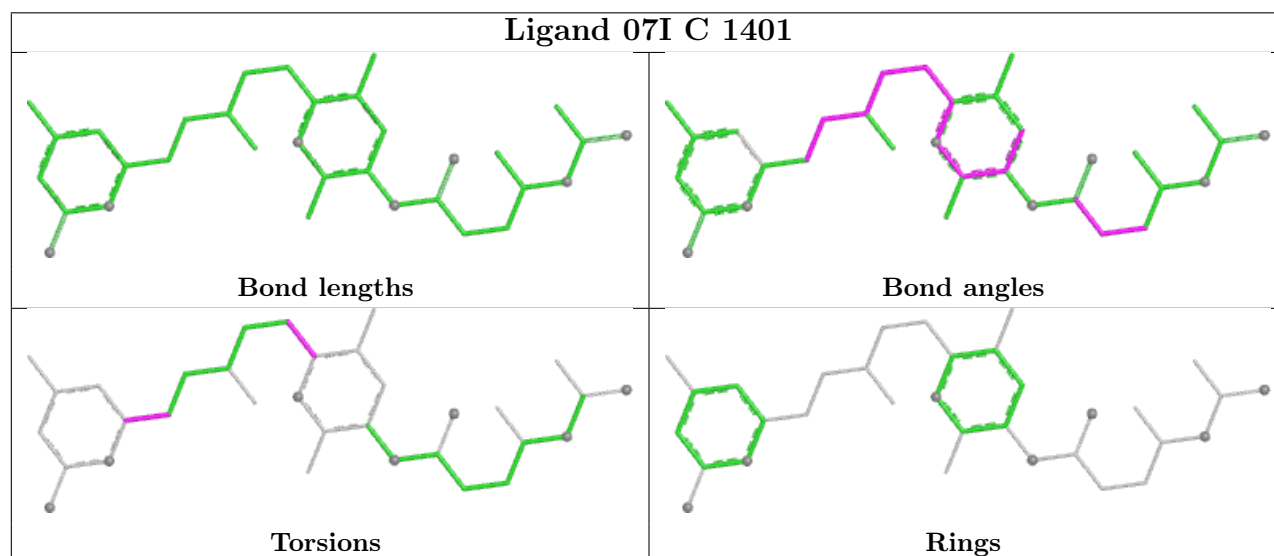
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Mol	Chain	Res	Type	Atoms
5	C	1401	07I	O-C5-C6-C7
5	C	1401	07I	C10-C11-C12-O2

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	863/899 (95%)	-0.18	5 (0%) 85 70	77, 115, 154, 183	0
2	B	65/86 (75%)	-0.03	0 100 100	86, 104, 118, 159	0
3	C	842/852 (98%)	-0.25	0 100 100	81, 121, 151, 192	0
4	D	93/108 (86%)	-0.35	0 100 100	80, 97, 146, 151	0
All	All	1863/1945 (95%)	-0.21	5 (0%) 90 80	77, 117, 151, 192	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	278	LEU	3.1
1	A	387	PHE	2.7
1	A	385	PHE	2.5
1	A	148	ALA	2.2
1	A	39	GLU	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 oligosaccharides in this entry.

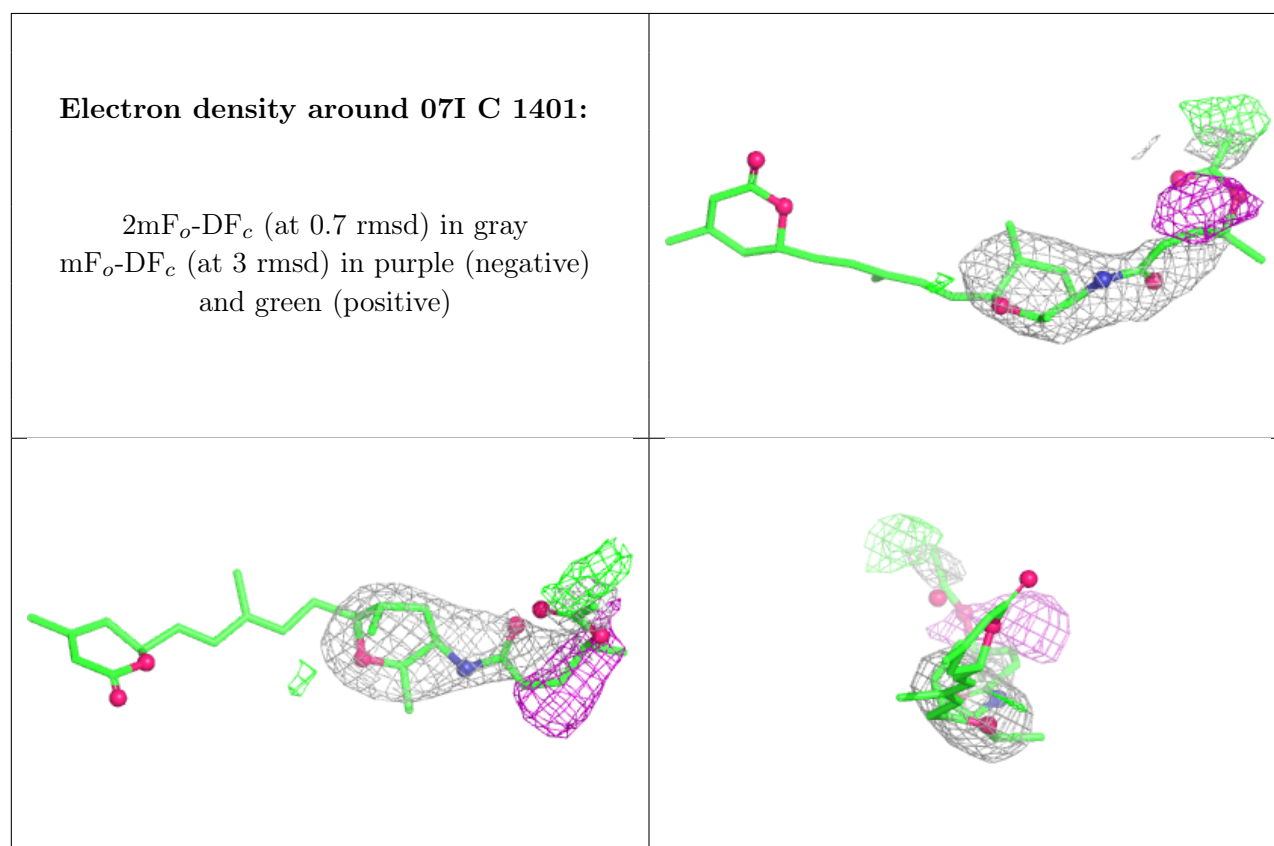
6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
5	07I	C	1401	33/33	0.86	0.20	97,116,133,134	15
6	ZN	D	103	1/1	0.99	0.09	113,113,113,113	0
6	ZN	D	102	1/1	1.00	0.01	96,96,96,96	0
6	ZN	D	101	1/1	1.00	0.03	100,100,100,100	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.