



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2021 – 06:05 PM EST

PDB ID : 6XE3  
Title : Salmonella typhimurium Tryptophan Synthase beta-S377A mutant in complex with inhibitor F9 at the enzyme alpha-site, cesium ion at the metal coordination site and carbanion III E(C3) at the enzyme beta-site.  
Authors : Hilario, E.; Dunn, M.F.; Mueller, L.J.; Fan, L.  
Deposited on : 2020-06-11  
Resolution : 1.55 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](mailto: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.

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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 : 2.17.1  
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) : 2.17.1

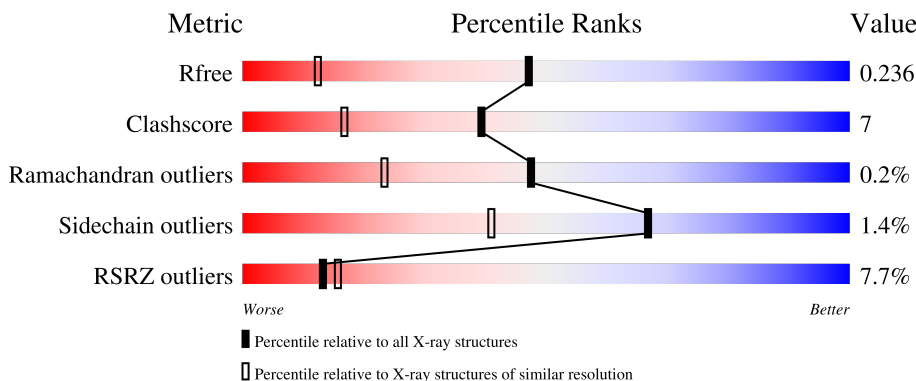
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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  $\geq 3$ , 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	268	<div> <div>10%</div> <div>87%</div> <div>12%</div> </div>
2	B	397	<div> <div>6%</div> <div>89%</div> <div>11%</div> </div>

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	EDO	A	305	-	-	X	-

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 6081 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 Tryptophan synthase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	4	0
			2051	1298	357	388	8			

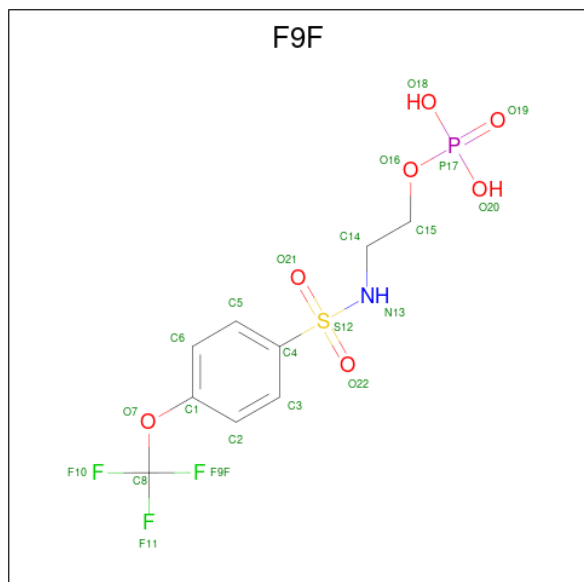
- Molecule 2 is a protein called Tryptophan synthase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	396	Total	C	N	O	S	0	11	0
			3101	1946	545	591	19			

There is a discrepancy between the modelled and reference sequences:

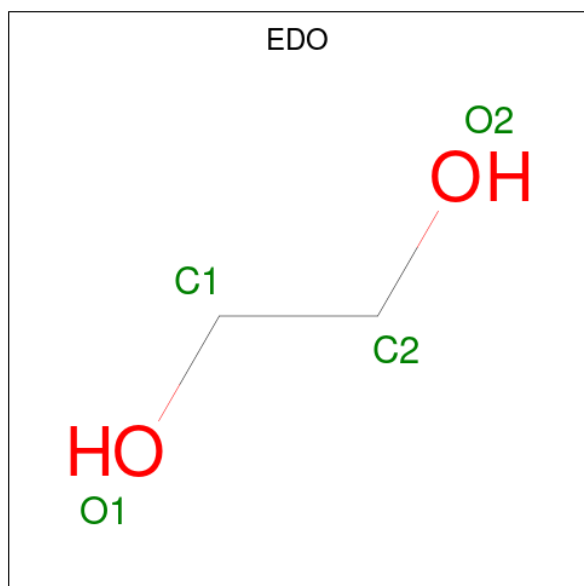
Chain	Residue	Modelled	Actual	Comment	Reference
B	377	ALA	SER	engineered mutation	UNP P0A2K1

- Molecule 3 is 2-([4-(TRIFLUOROMETHOXY)PHENYL]SULFONYL)AMINO)ETHYL DIHYDROGEN PHOSPHATE (three-letter code: F9F) (formula: C<sub>9</sub>H<sub>11</sub>F<sub>3</sub>NO<sub>7</sub>PS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	P	S	0	0
			22	9	3	1	7	1	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C<sub>2</sub>H<sub>6</sub>OS).



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

- | Mol | Chain | Residues | Atoms           | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 6   | A     | 1        | Total Cs<br>1 1 | 0       | 0       |
| 6   | B     | 3        | Total Cs<br>4 4 | 0       | 1       |

- | Mol | Chain | Residues | Atoms             | ZeroOcc | AltConf |
|-----|-------|----------|-------------------|---------|---------|
| 7   | A     | 1        | Total Cl<br>1 1   | 0       | 0       |
| 7   | B     | 14       | Total Cl<br>14 14 | 0       | 0       |

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	1	Total	C	N	O	P	0	0
			29	17	3	8	1		

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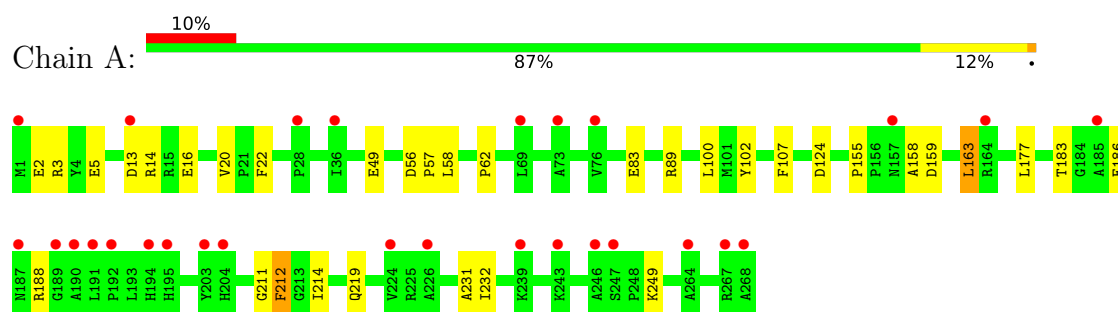
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	273	Total 284	O 284	0	12
9	B	469	Total 486	O 486	0	18



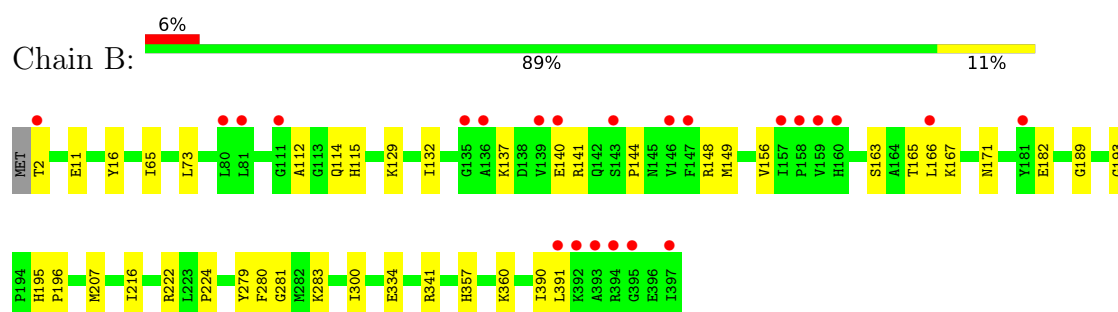
### 3 Residue-property plots [i](#)

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: Tryptophan synthase alpha chain



- Molecule 2: Tryptophan synthase beta chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	181.52Å 59.43Å 67.06Å 90.00° 94.64° 90.00°	Depositor
Resolution (Å)	38.95 – 1.55 38.95 – 1.55	Depositor EDS
% Data completeness (in resolution range)	94.7 (38.95-1.55) 94.7 (38.95-1.55)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.15 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.174 , 0.236 0.174 , 0.236	Depositor DCC
$R_{free}$ test set	4983 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	15.1	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 58.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6081	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 1D0, EDO, CS, F9F, DMS, CL

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	0.34	0/2091	0.64	0/2839
2	B	0.38	0/3159	0.72	0/4263
All	All	0.36	0/5250	0.69	0/7102

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	A	2051	0	2047	30	0
2	B	3101	0	3071	42	0
3	A	22	0	9	2	0
4	A	8	0	12	6	0
4	B	24	0	35	0	0
5	A	24	0	36	3	0
5	B	32	0	48	4	0
6	A	1	0	0	0	0
6	B	4	0	0	0	0
7	A	1	0	0	0	0
7	B	14	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	29	0	17	0	0
9	A	284	0	0	5	0
9	B	486	0	0	30	0
All	All	6081	0	5275	78	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 7.

All (78) 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)
2:B:148:ARG:NH1	9:B:501:HOH:O	1.90	1.02
2:B:166[B]:LEU:HD12	2:B:166[B]:LEU:H	1.40	0.86
2:B:114:GLN:HB2	9:B:775:HOH:O	1.76	0.85
4:A:305:EDO:H12	9:A:406:HOH:O	1.76	0.84
2:B:222:ARG:NH2	9:B:503:HOH:O	2.10	0.83
1:A:56:ASP:OD2	9:A:401:HOH:O	1.99	0.80
1:A:107:PHE:CE2	4:A:305:EDO:H11	2.17	0.79
7:B:416:CL:CL	9:B:570:HOH:O	2.39	0.77
2:B:132:ILE:O	9:B:502:HOH:O	2.05	0.73
2:B:166[B]:LEU:HD11	9:B:678:HOH:O	1.89	0.71
1:A:58:LEU:CD2	9:B:608:HOH:O	2.39	0.69
2:B:144:PRO:C	9:B:501:HOH:O	2.32	0.67
2:B:182[B]:GLU:OE2	9:B:504:HOH:O	2.12	0.66
1:A:107:PHE:CZ	4:A:305:EDO:H11	2.30	0.66
7:B:415:CL:CL	9:B:821:HOH:O	2.51	0.65
1:A:107:PHE:CE2	4:A:305:EDO:H22	2.32	0.65
1:A:159:ASP:O	1:A:163:LEU:HD23	1.97	0.65
2:B:132:ILE:HG22	9:B:502:HOH:O	1.97	0.65
2:B:341:ARG:HG3	9:B:660:HOH:O	1.98	0.64
9:A:401:HOH:O	2:B:171:ASN:ND2	2.30	0.64
2:B:166[B]:LEU:HD12	2:B:166[B]:LEU:N	2.13	0.63
2:B:166[B]:LEU:H	2:B:166[B]:LEU:CD1	2.11	0.62
2:B:156:VAL:HG12	9:B:574:HOH:O	2.02	0.60
2:B:216[A]:ILE:HG21	2:B:224:PRO:HD3	1.84	0.59
1:A:58:LEU:HD21	9:B:608:HOH:O	1.99	0.59
1:A:188:ARG:HG2	1:A:219:GLN:NE2	2.17	0.59
1:A:214:ILE:HD13	1:A:231:ALA:HB1	1.84	0.57
1:A:83[B]:GLU:HA	1:A:83[B]:GLU:OE1	2.05	0.56
1:A:3:ARG:NH1	5:A:308:DMS:H22	2.22	0.54
2:B:357:HIS:HE1	5:B:408:DMS:H21	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:112:ALA:CB	9:B:598:HOH:O	2.56	0.52
2:B:165:THR:HG21	9:B:841:HOH:O	2.09	0.52
3:A:301:F9F:H5	9:A:465:HOH:O	2.09	0.51
1:A:3:ARG:HH12	5:A:308:DMS:H22	1.76	0.50
2:B:300:ILE:HD11	2:B:390:ILE:HD13	1.93	0.50
1:A:107:PHE:CE2	4:A:305:EDO:C2	2.95	0.50
2:B:165:THR:OG1	2:B:166[B]:LEU:HD12	2.11	0.50
1:A:14:ARG:NH2	9:A:407:HOH:O	2.45	0.49
1:A:100:LEU:C	1:A:100:LEU:HD13	2.32	0.49
5:B:408:DMS:H22	9:B:920:HOH:O	2.13	0.49
2:B:115:HIS:CE1	2:B:189:GLY:HA2	2.47	0.49
2:B:283:LYS:NZ	9:B:506[A]:HOH:O	2.32	0.49
2:B:357:HIS:HE1	5:B:408:DMS:C2	2.25	0.49
1:A:3:ARG:HH12	5:A:308:DMS:C2	2.26	0.48
2:B:163:SER:OG	2:B:165:THR:HG22	2.13	0.48
2:B:300:ILE:HD11	2:B:390:ILE:CD1	2.44	0.48
1:A:14:ARG:HD2	1:A:16:GLU:OE1	2.13	0.48
2:B:137:LYS:HB2	9:B:528:HOH:O	2.15	0.47
5:B:430:DMS:H23	9:B:846:HOH:O	2.15	0.47
2:B:129:LYS:HE3	9:B:552:HOH:O	2.15	0.47
2:B:193:GLY:HA2	2:B:280:PHE:O	2.15	0.47
2:B:132:ILE:C	9:B:502:HOH:O	2.51	0.46
1:A:155:PRO:HD2	1:A:158:ALA:HB2	1.97	0.45
1:A:183:THR:HG22	3:A:301:F9F:H142	1.98	0.45
2:B:156:VAL:HG13	9:B:502:HOH:O	2.17	0.45
1:A:58:LEU:HD23	9:B:608:HOH:O	2.12	0.44
1:A:211:GLY:O	1:A:212:PHE:HB2	2.18	0.44
1:A:155:PRO:HA	1:A:177:LEU:HB3	1.99	0.43
2:B:341:ARG:NH1	9:B:511:HOH:O	2.39	0.43
1:A:107:PHE:CE2	4:A:305:EDO:C1	2.95	0.43
2:B:132:ILE:HD13	2:B:149:MET:SD	2.59	0.43
2:B:195:HIS:CG	2:B:196:PRO:HA	2.54	0.43
1:A:83[B]:GLU:OE1	1:A:83[B]:GLU:CA	2.67	0.42
2:B:11[A]:GLU:O	2:B:11[A]:GLU:HG2	2.19	0.42
2:B:16:TYR:O	2:B:281:GLY:HA2	2.20	0.42
1:A:62:PRO:HG3	9:B:902:HOH:O	2.20	0.42
1:A:2:GLU:HG2	1:A:5:GLU:OE1	2.20	0.41
1:A:22:PHE:HA	1:A:49:GLU:O	2.21	0.41
2:B:360[A]:LYS:HG3	9:B:633:HOH:O	2.21	0.41
1:A:57:PRO:HA	1:A:102:TYR:CZ	2.56	0.41
1:A:89:ARG:NH2	1:A:124:ASP:OD2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:65:ILE:HD11	2:B:73:LEU:HD23	2.03	0.41
2:B:141:ARG:NH1	9:B:528:HOH:O	2.53	0.41
1:A:20:VAL:O	1:A:232:ILE:HA	2.20	0.41
2:B:334[B]:GLU:HG2	9:B:564:HOH:O	2.21	0.41
2:B:195:HIS:ND1	2:B:196:PRO:HA	2.36	0.41
2:B:2:THR:HG22	9:B:842:HOH:O	2.21	0.40
2:B:167:LYS:HB2	2:B:279:TYR:OH	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	A	270/268 (101%)	264 (98%)	5 (2%)	1 (0%)	34	14
2	B	406/397 (102%)	399 (98%)	7 (2%)	0	100	100
All	All	676/665 (102%)	663 (98%)	12 (2%)	1 (0%)	47	26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	212	PHE

### 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	212/208 (102%)	208 (98%)	4 (2%)	57	28
2	B	321/310 (104%)	318 (99%)	3 (1%)	78	61
All	All	533/518 (103%)	526 (99%)	7 (1%)	67	44

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASP
1	A	163	LEU
1	A	186	GLU
1	A	249	LYS
2	B	140	GLU
2	B	207	MET
2	B	391	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	342	HIS

### 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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 44 ligands modelled in this entry, 20 are monoatomic - leaving 24 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	DMS	B	402	-	3,3,3	0.25	0	3,3,3	0.11	0
5	DMS	B	403	-	3,3,3	0.57	0	3,3,3	0.27	0
5	DMS	B	431	-	3,3,3	0.23	0	3,3,3	0.12	0
5	DMS	B	427	-	3,3,3	0.78	0	3,3,3	0.68	0
5	DMS	A	308	-	3,3,3	0.17	0	3,3,3	0.46	0
5	DMS	B	408	-	3,3,3	0.20	0	3,3,3	0.26	0
4	EDO	B	407	-	3,3,3	0.13	0	2,2,2	0.25	0
4	EDO	A	305	-	3,3,3	0.68	0	2,2,2	1.28	0
5	DMS	B	430	-	3,3,3	0.31	0	3,3,3	0.22	0
4	EDO	B	409	-	3,3,3	0.06	0	2,2,2	0.18	0
5	DMS	A	309	-	3,3,3	0.25	0	3,3,3	0.31	0
4	EDO	B	432	-	3,3,3	0.07	0	2,2,2	0.56	0
5	DMS	A	304	-	3,3,3	0.20	0	3,3,3	0.10	0
5	DMS	A	303	-	3,3,3	0.21	0	3,3,3	0.10	0
4	EDO	A	302	-	3,3,3	0.28	0	2,2,2	0.60	0
5	DMS	A	311	-	3,3,3	0.25	0	3,3,3	0.18	0
4	EDO	B	406	-	3,3,3	0.08	0	2,2,2	0.14	0
8	1D0	B	401	-	27,30,30	2.19	2 (7%)	33,42,42	0.91	1 (3%)
5	DMS	B	429	-	3,3,3	0.34	0	3,3,3	0.04	0
3	F9F	A	301	-	22,22,22	0.62	1 (4%)	32,33,33	0.75	1 (3%)
5	DMS	B	428	-	3,3,3	0.24	0	3,3,3	0.21	0
4	EDO	B	405	6	3,3,3	0.08	0	2,2,2	0.38	0
4	EDO	B	404	-	3,3,3	0.06	0	2,2,2	0.22	0
5	DMS	A	310	-	3,3,3	0.22	0	3,3,3	0.12	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	EDO	A	302	-	-	1/1/1/1	-
4	EDO	B	407	-	-	1/1/1/1	-
4	EDO	B	406	-	-	1/1/1/1	-
4	EDO	A	305	-	-	0/1/1/1	-
8	1D0	B	401	-	-	4/13/20/20	0/2/2/2
4	EDO	B	432	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	405	6	-	1/1/1/1	-
3	F9F	A	301	-	-	5/20/20/20	0/1/1/1
4	EDO	B	404	-	-	1/1/1/1	-
4	EDO	B	409	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	401	1D0	C-CA	-10.33	1.36	1.52
8	B	401	1D0	CB-CA	-3.41	1.47	1.51
3	A	301	F9F	P17-O19	2.36	1.58	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	401	1D0	CB-N-C1	-3.05	119.61	123.98
3	A	301	F9F	O22-S12-O21	2.21	122.26	119.55

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	401	1D0	C51-C41-C4A-N2
8	B	401	1D0	C31-C41-C4A-N2
8	B	401	1D0	C6-C1-N-CB
8	B	401	1D0	C2-C1-N-CB
4	B	405	EDO	O1-C1-C2-O2
4	B	409	EDO	O1-C1-C2-O2
4	A	302	EDO	O1-C1-C2-O2
4	B	406	EDO	O1-C1-C2-O2
4	B	407	EDO	O1-C1-C2-O2
3	A	301	F9F	N13-C14-C15-O16
3	A	301	F9F	C5-C4-S12-O21
3	A	301	F9F	C2-C1-O7-C8
3	A	301	F9F	C6-C1-O7-C8
3	A	301	F9F	C3-C4-S12-O21
4	B	404	EDO	O1-C1-C2-O2
4	B	432	EDO	O1-C1-C2-O2

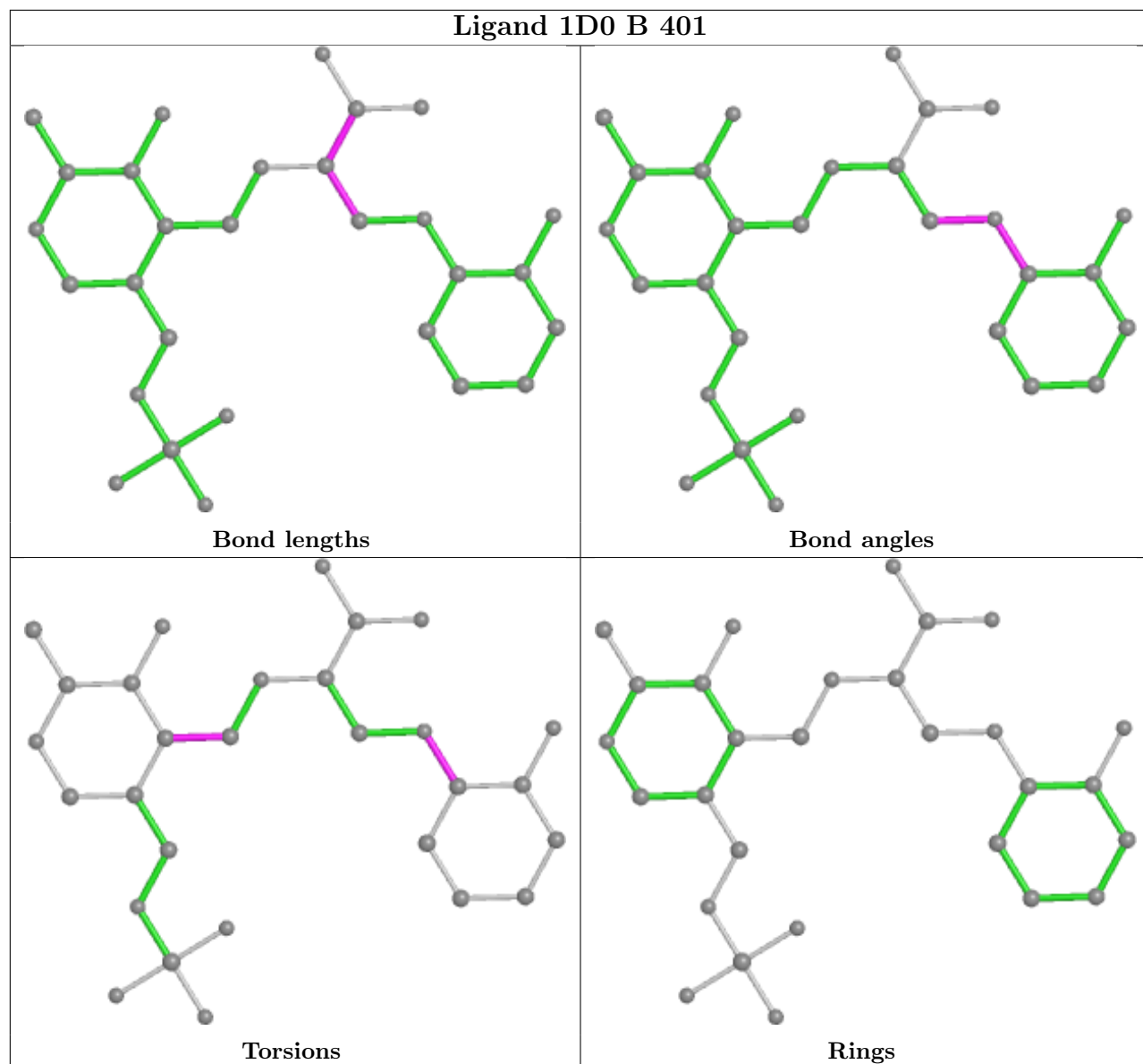
There are no ring outliers.

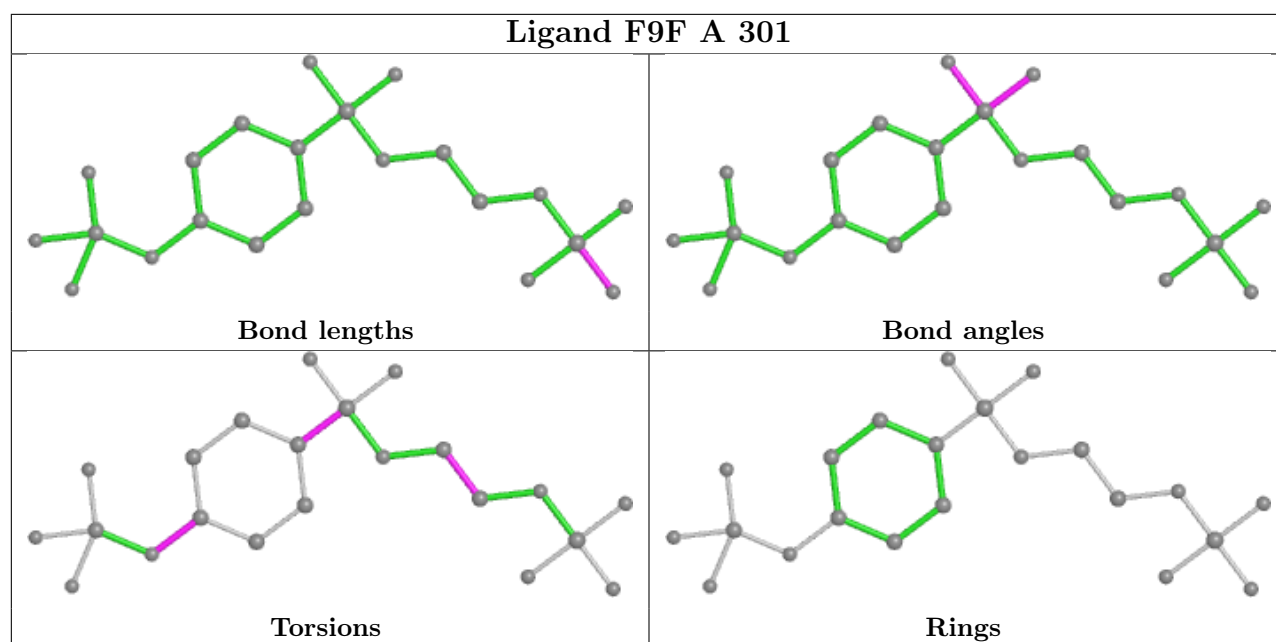
5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	308	DMS	3	0
5	B	408	DMS	3	0
4	A	305	EDO	6	0
5	B	430	DMS	1	0
3	A	301	F9F	2	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.

## Ligand 1D0 B 401





## 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

### 6.1 Protein, DNA and RNA chains

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	268/268 (100%)	0.67	28 (10%) <b>6</b> <b>6</b>	12, 22, 42, 60	0
2	B	396/397 (99%)	0.26	23 (5%) <b>23</b> <b>26</b>	9, 15, 30, 64	0
All	All	664/665 (99%)	0.43	51 (7%) <b>13</b> <b>16</b>	9, 19, 36, 64	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	397	ILE	9.9
1	A	190	ALA	5.3
2	B	139	VAL	5.0
1	A	203	TYR	4.6
1	A	185	ALA	4.5
1	A	246	ALA	4.4
1	A	69	LEU	4.3
2	B	146	VAL	4.3
1	A	73	ALA	4.3
2	B	393	ALA	4.2
1	A	1	MET	4.2
1	A	224	VAL	4.1
1	A	192	PRO	4.1
2	B	395	GLY	4.1
2	B	135	GLY	3.7
1	A	195	HIS	3.6
2	B	166[A]	LEU	3.6
1	A	247	SER	3.6
2	B	147	PHE	3.5
2	B	159	VAL	3.5
1	A	239	LYS	3.4
2	B	158	PRO	3.4
1	A	76	VAL	3.2
1	A	28	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	189	GLY	3.2
2	B	181	TYR	3.1
2	B	136	ALA	3.1
1	A	13	ASP	2.9
1	A	243	LYS	2.9
2	B	111	GLY	2.9
1	A	194	HIS	2.8
2	B	394	ARG	2.8
1	A	204	HIS	2.8
2	B	140	GLU	2.5
1	A	191	LEU	2.4
1	A	226	ALA	2.4
1	A	187	ASN	2.4
2	B	160	HIS	2.4
2	B	392	LYS	2.4
1	A	267	ARG	2.3
1	A	157	ASN	2.3
2	B	80	LEU	2.3
2	B	143	SER	2.2
1	A	264	ALA	2.2
2	B	391	LEU	2.2
1	A	268	ALA	2.1
2	B	81	LEU	2.1
2	B	2	THR	2.1
1	A	36	ILE	2.1
2	B	157	ILE	2.0
1	A	164	ARG	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 [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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
7	CL	B	420	1/1	0.64	0.22	73,73,73,73	0
7	CL	A	307	1/1	0.81	0.08	82,82,82,82	0
4	EDO	B	405	4/4	0.84	0.14	39,45,48,50	0
4	EDO	B	406	4/4	0.85	0.20	35,36,37,46	0
5	DMS	A	303	4/4	0.87	0.12	45,50,52,65	0
5	DMS	B	403	4/4	0.87	0.15	26,37,41,46	0
5	DMS	A	308	4/4	0.89	0.19	22,36,39,41	0
4	EDO	B	404	4/4	0.92	0.10	36,36,41,41	0
4	EDO	B	407	4/4	0.92	0.09	31,31,34,43	0
4	EDO	B	432	4/4	0.92	0.12	25,28,35,37	0
7	CL	B	418	1/1	0.92	0.08	66,66,66,66	0
4	EDO	A	302	4/4	0.92	0.12	25,29,33,34	0
7	CL	B	425	1/1	0.92	0.11	61,61,61,61	0
4	EDO	B	409	4/4	0.93	0.09	37,38,46,48	0
7	CL	B	419	1/1	0.93	0.04	73,73,73,73	0
3	F9F	A	301	22/22	0.93	0.11	17,20,22,24	0
7	CL	B	424	1/1	0.93	0.12	70,70,70,70	0
7	CL	B	413	1/1	0.93	0.04	62,62,62,62	0
7	CL	B	426	1/1	0.93	0.04	71,71,71,71	0
4	EDO	A	305	4/4	0.94	0.11	12,14,17,25	0
5	DMS	A	304	4/4	0.94	0.11	39,43,52,52	0
7	CL	B	417	1/1	0.94	0.04	58,58,58,58	0
5	DMS	B	408	4/4	0.95	0.12	29,39,48,49	0
5	DMS	B	402	4/4	0.96	0.14	47,48,59,59	0
7	CL	B	422	1/1	0.96	0.06	44,44,44,44	0
5	DMS	B	429	4/4	0.96	0.15	29,32,36,38	0
5	DMS	B	431	4/4	0.96	0.10	33,43,45,47	0
5	DMS	A	311	4/4	0.96	0.12	37,38,41,51	0
8	1D0	B	401	29/29	0.96	0.12	11,20,27,27	0
5	DMS	B	430	4/4	0.97	0.10	28,33,35,39	0
5	DMS	A	309	4/4	0.98	0.10	23,24,26,32	0
7	CL	B	421	1/1	0.98	0.08	53,53,53,53	0
7	CL	B	414	1/1	0.98	0.12	56,56,56,56	0
7	CL	B	416	1/1	0.98	0.05	41,41,41,41	0
5	DMS	A	310	4/4	0.98	0.08	21,22,23,24	0
5	DMS	B	427	4/4	0.98	0.07	10,15,16,22	0
5	DMS	B	428	4/4	0.98	0.13	29,34,35,39	0
7	CL	B	415	1/1	0.99	0.07	40,40,40,40	0
7	CL	B	423	1/1	0.99	0.02	34,34,34,34	0
6	CS	B	410[A]	1/1	1.00	0.06	17,17,17,17	1
6	CS	B	410[B]	1/1	1.00	0.06	22,22,22,22	1

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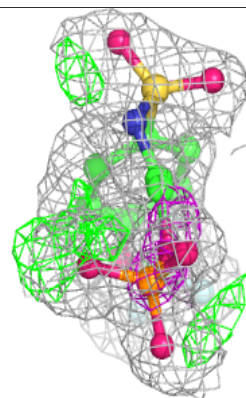
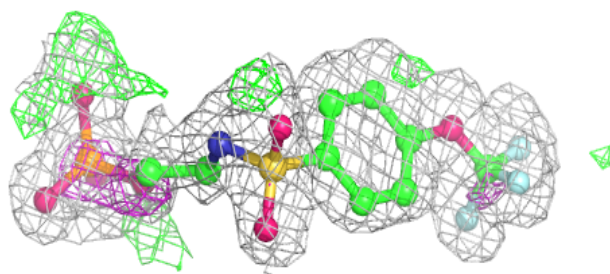
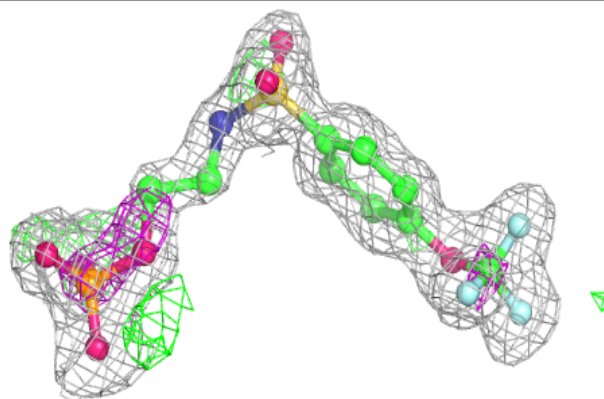
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	CS	B	411	1/1	1.00	0.04	17,17,17,17	0
6	CS	B	412	1/1	1.00	0.02	26,26,26,26	1
6	CS	A	306	1/1	1.00	0.05	34,34,34,34	1

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 F9F A 301:**

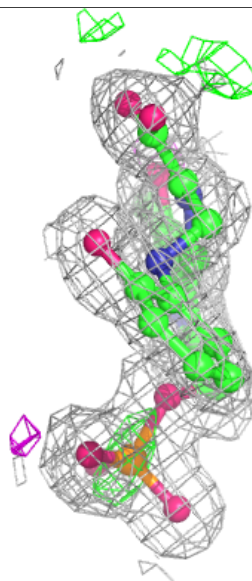
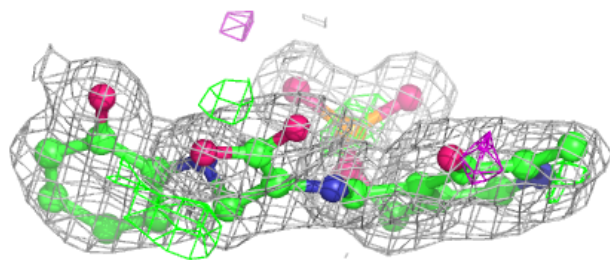
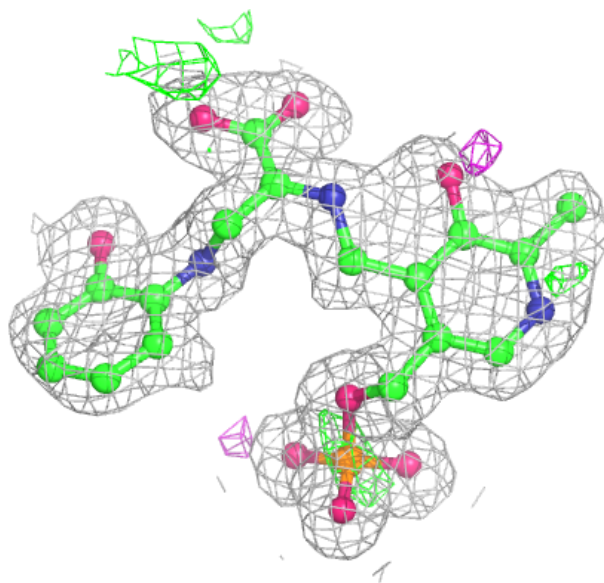
2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)





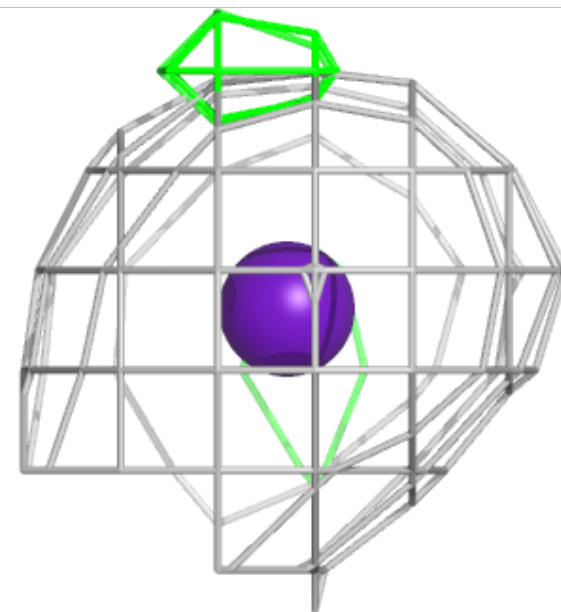
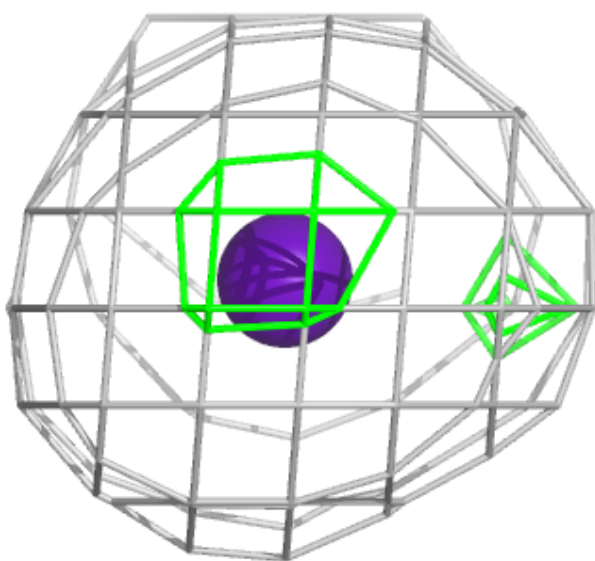
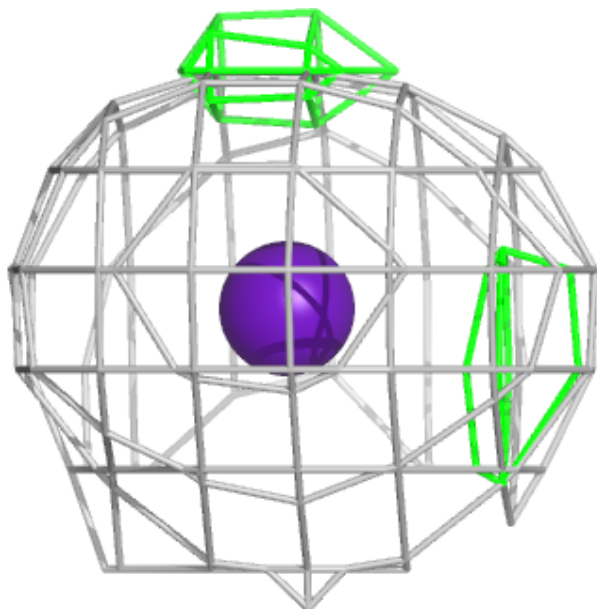
**Electron density around 1D0 B 401:**

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



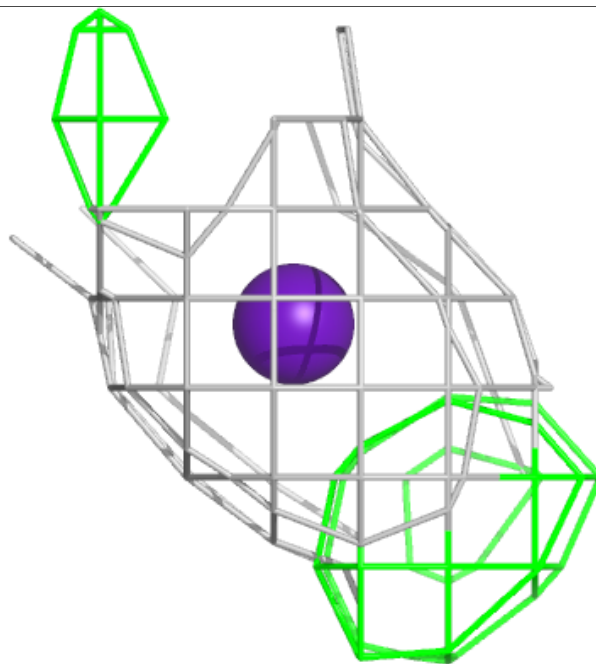
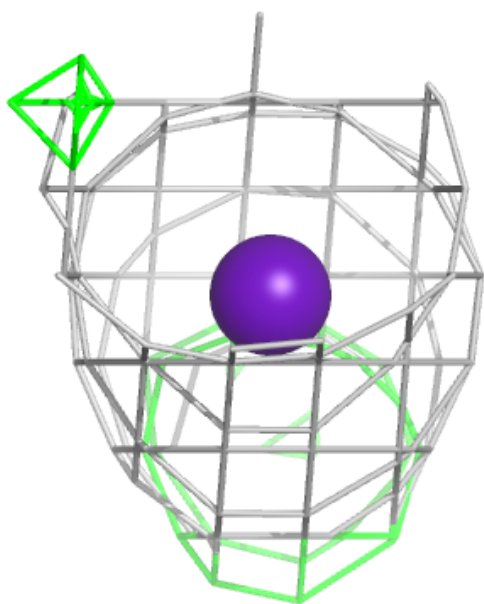
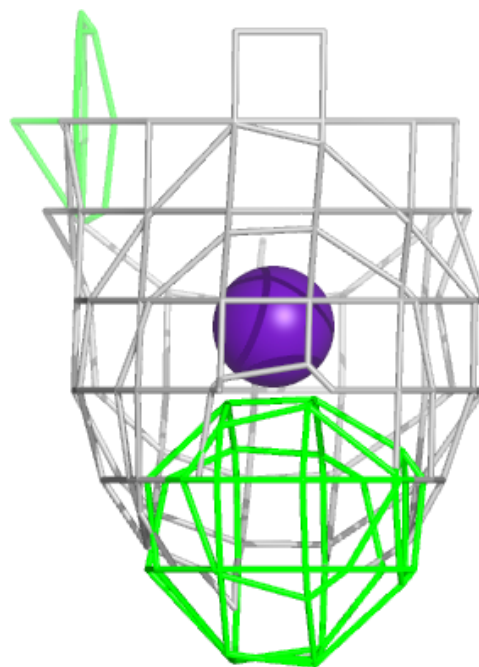
**Electron density around CS B 410 (A):**

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



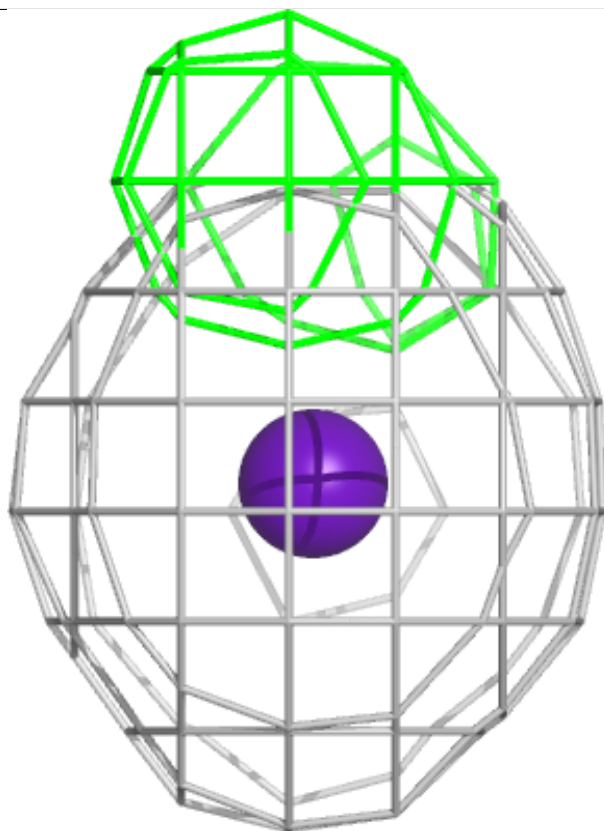
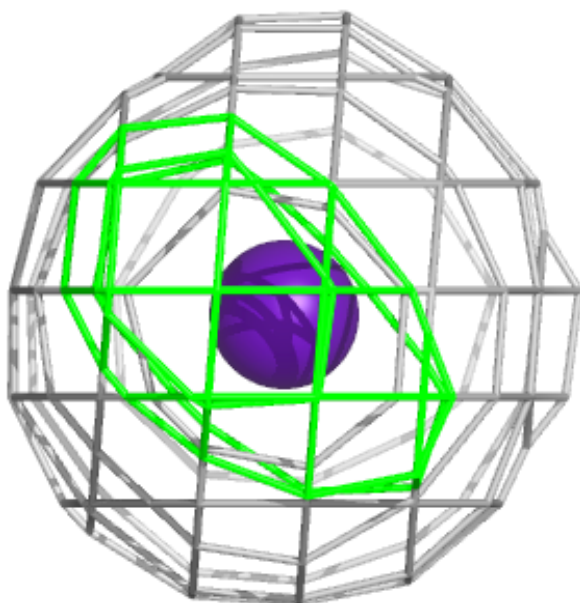
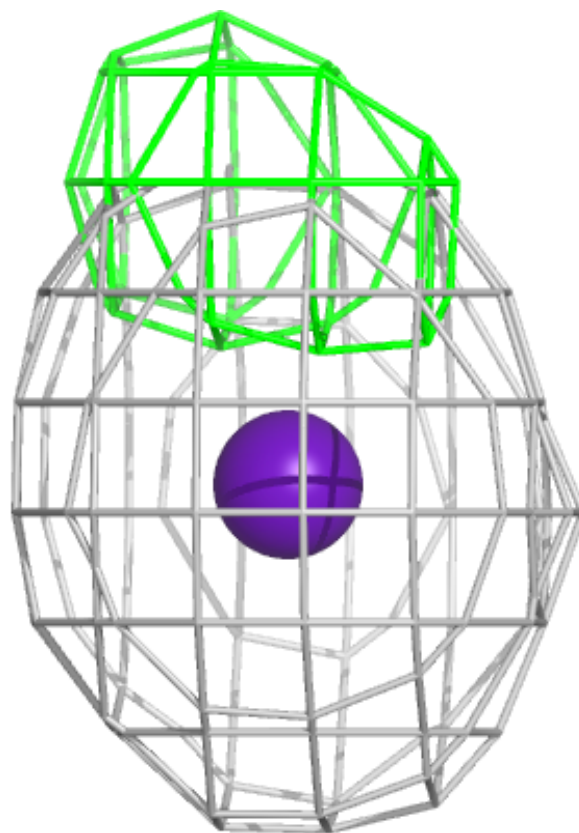
**Electron density around CS B 410 (B):**

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



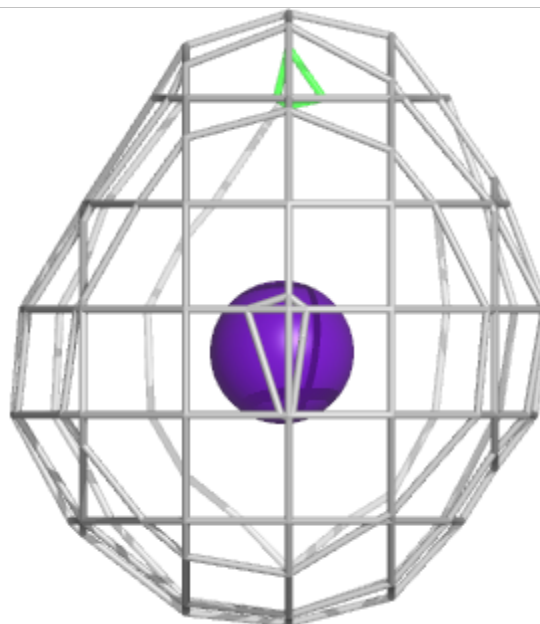
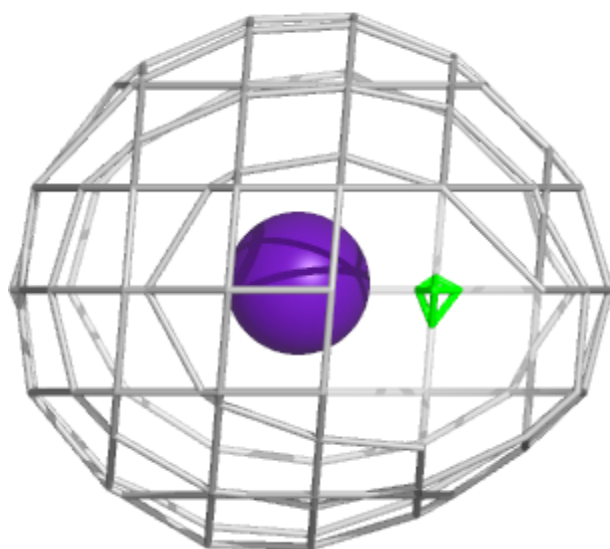
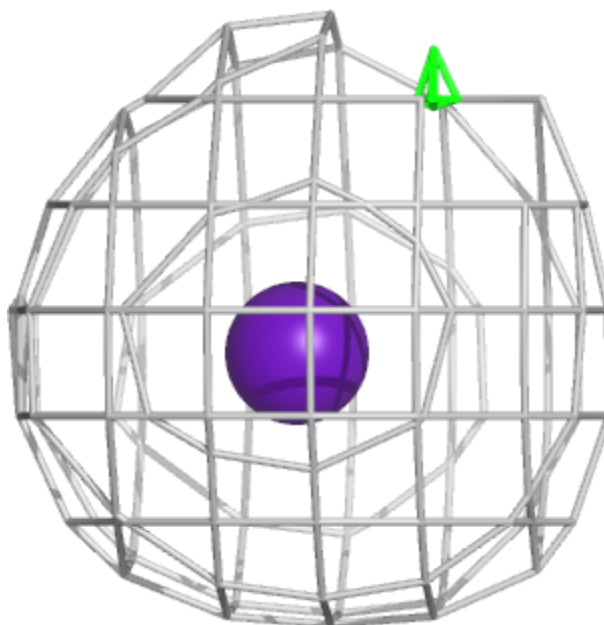
**Electron density around CS B 411:**

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



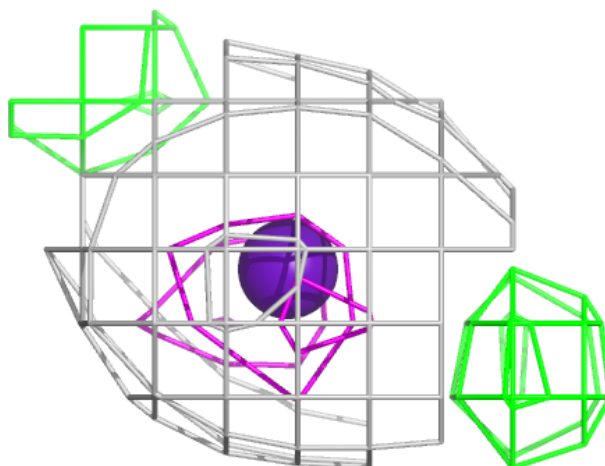
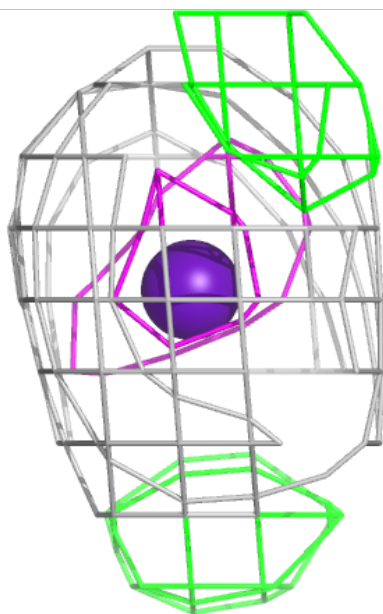
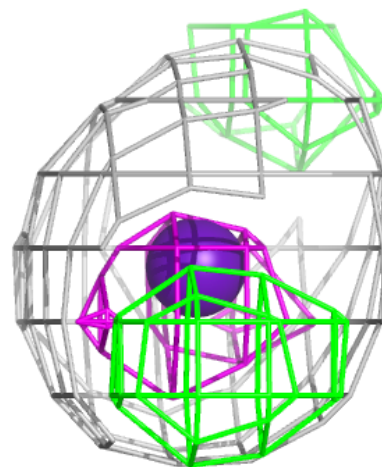
**Electron density around CS B 412:**

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



**Electron density around CS A 306:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.