



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:13 AM UTC

PDB ID : 4O7U / pdb_00004o7u
Title : Etherocomplex of Enterococcus faecalis thymidylate synthase with 5-hydroxy methylene-6-hydrofolic acid and the phtalimidic inhibitor SS7
Authors : Mangani, S.; Pozzi, C.
Deposited on : 2013-12-26
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (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

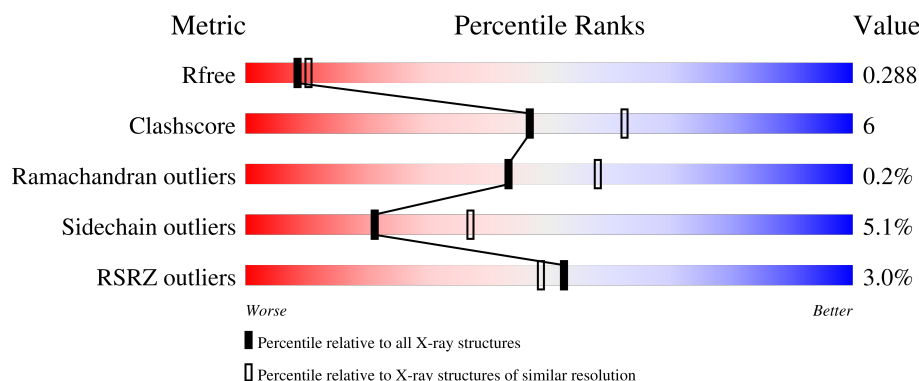
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	315	<div> <div>5%</div> <div>76%</div> <div>16%</div> <div>8%</div> </div>
1	B	315	<div> <div>88%</div> <div>11%</div> <div>.</div> </div>
1	C	315	<div> <div>6%</div> <div>69%</div> <div>16%</div> <div>14%</div> </div>
1	D	315	<div> <div>%</div> <div>84%</div> <div>15%</div> <div>.</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
5	SS7	D	402	-	-	X	-

2 Entry composition [i](#)

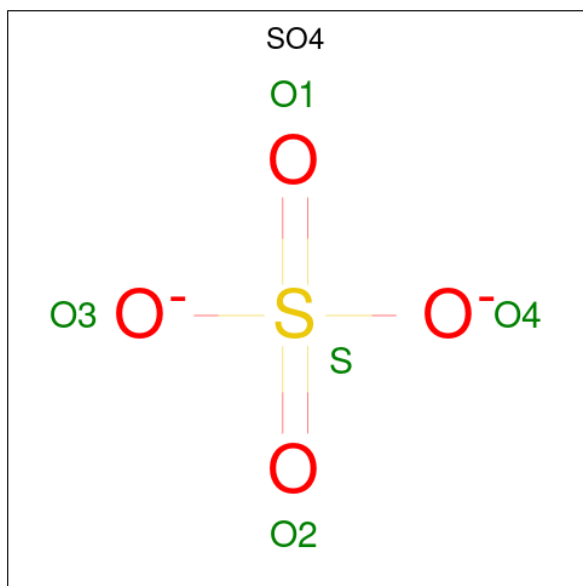
There are 6 unique types of molecules in this entry. The entry contains 10032 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thymidylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2378	1533	394	441	10			
1	B	315	Total	C	N	O	S	0	0	0
			2569	1651	427	480	11			
1	C	272	Total	C	N	O	S	0	0	0
			2222	1436	370	407	9			
1	D	315	Total	C	N	O	S	0	0	0
			2569	1651	427	480	11			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



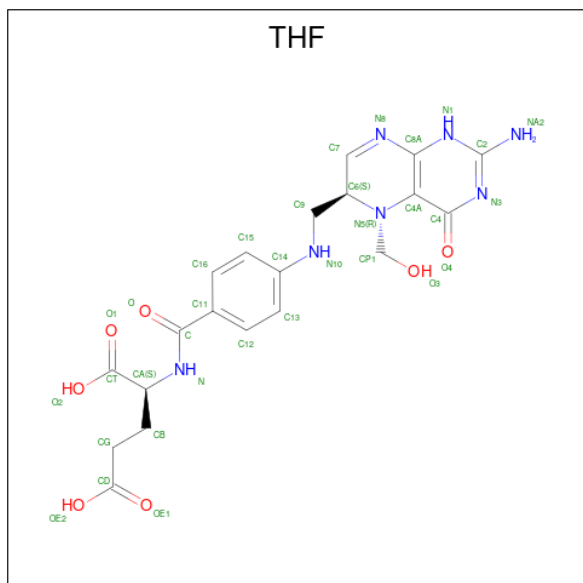
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is 5-HYDROXYMETHYLENE-6-HYDROFOLIC ACID (CCD ID: THF) (formula: $C_{20}H_{23}N_7O_7$).



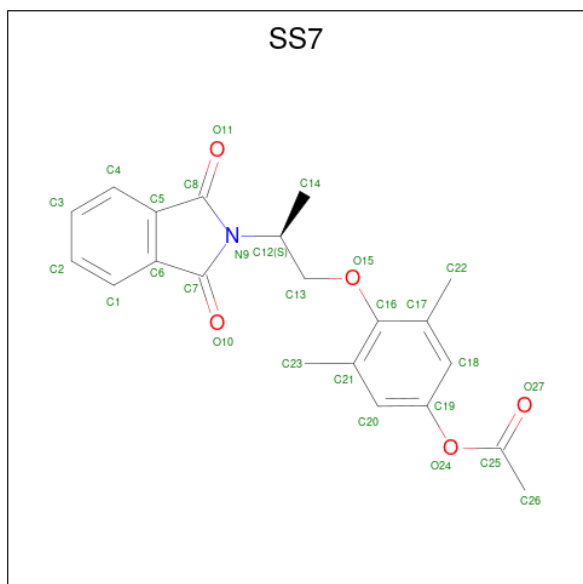
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			34	20	7	7		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 5 is 4-[(2S)-2-(1,3-dioxo-1,3-dihydro-2H-isoindol-2-yl)propyl]oxy}-3,5-dimethylphenyl acetate (CCD ID: SS7) (formula: $C_{21}H_{21}NO_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			27	21	1	5		

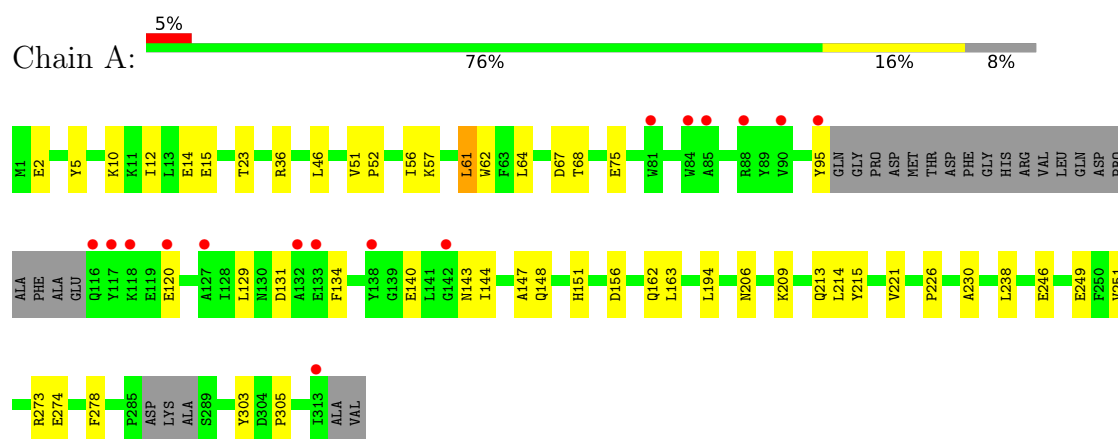
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	38	Total 38	O 38	0	0
6	B	69	Total 69	O 69	0	0
6	C	34	Total 34	O 34	0	0
6	D	64	Total 64	O 64	0	0

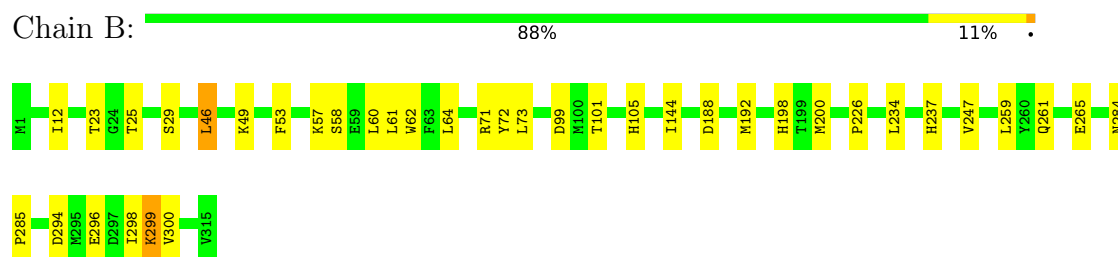
3 Residue-property plots [i](#)

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.

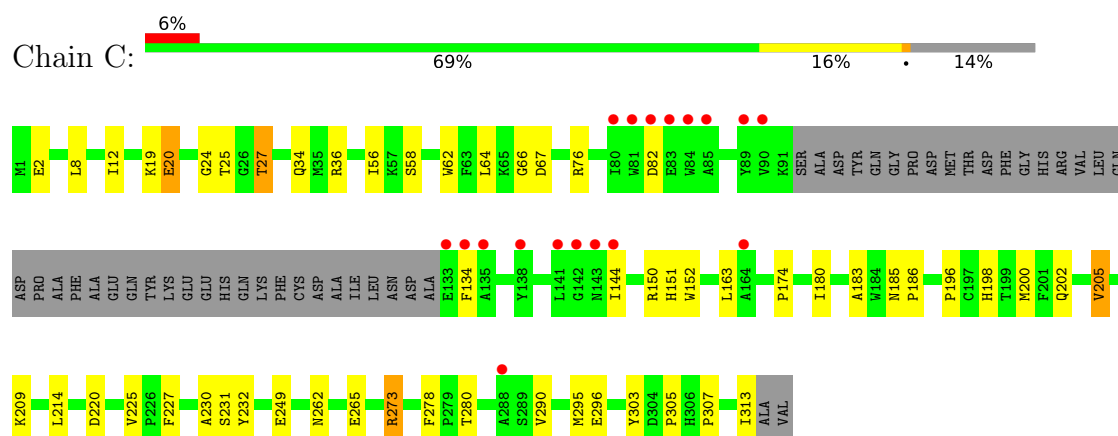
• Molecule 1: Thymidylate synthase



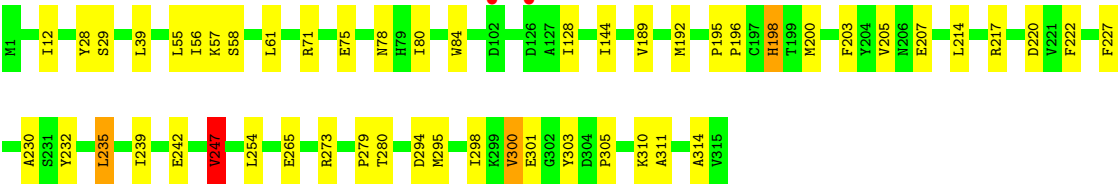
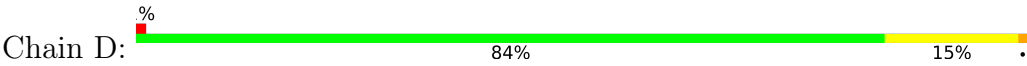
• Molecule 1: Thymidylate synthase



• Molecule 1: Thymidylate synthase



● Molecule 1: Thymidylate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.08Å 94.22Å 96.82Å 90.00° 94.33° 90.00°	Depositor
Resolution (Å)	55.67 – 2.40 55.67 – 2.40	Depositor EDS
% Data completeness (in resolution range)	76.1 (55.67-2.40) 76.1 (55.67-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.212 , 0.287 0.213 , 0.288	Depositor DCC
R_{free} test set	1962 reflections (3.88%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10032	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, SO4, SS7, THF

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	0.55	0/2443	0.86	0/3308
1	B	0.62	0/2641	0.92	0/3579
1	C	0.56	0/2284	0.88	0/3095
1	D	0.60	0/2641	0.91	3/3579 (0.1%)
All	All	0.58	0/10009	0.89	3/13561 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	222	PHE	N-CA-C	5.74	118.40	111.40
1	D	29	SER	N-CA-C	5.59	117.56	109.07
1	D	247	VAL	CB-CA-C	5.58	117.83	110.91

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	2378	0	2287	23	0
1	B	2569	0	2463	20	0
1	C	2222	0	2153	31	0
1	D	2569	0	2463	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	10	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	B	34	0	21	6	0
4	C	4	0	6	1	0
4	D	4	0	6	0	0
5	D	27	0	21	13	0
6	A	38	0	0	0	0
6	B	69	0	0	1	0
6	C	34	0	0	0	0
6	D	64	0	0	0	0
All	All	10032	0	9420	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.

All (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:402:SS7:C22	5:D:402:SS7:H9	1.95	0.95
5:D:402:SS7:H9	5:D:402:SS7:H13	1.47	0.94
1:C:174:PRO:HB2	1:C:205:VAL:HG11	1.56	0.86
1:B:46:LEU:HD22	1:B:226:PRO:HB3	1.57	0.86
3:B:401:THF:HCP1	3:B:401:THF:HN0	1.40	0.83
1:C:200:MET:HE3	1:D:200:MET:HE3	1.60	0.83
1:D:78:ASN:OD1	1:D:80:ILE:HG12	1.82	0.80
1:D:12:ILE:HD11	1:D:28:TYR:O	1.81	0.79
1:D:84:TRP:CZ2	5:D:402:SS7:H1	2.21	0.75
1:C:58:SER:HA	1:C:295:MET:HE1	1.70	0.74
3:B:401:THF:O4	3:B:401:THF:HCP2	1.89	0.71
1:B:25:THR:O	1:B:261:GLN:NE2	2.27	0.68
1:B:296:GLU:OE1	1:B:299:LYS:NZ	2.26	0.67
3:B:401:THF:HCP1	3:B:401:THF:N10	2.09	0.67
1:C:174:PRO:CB	1:C:205:VAL:HG11	2.26	0.65
1:C:180:ILE:HD13	1:C:202:GLN:HB2	1.80	0.64
1:B:12:ILE:HG21	1:B:259:LEU:HB2	1.80	0.64
1:C:19:LYS:NZ	1:D:205:VAL:O	2.31	0.63
1:D:314:ALA:O	5:D:402:SS7:H2	1.99	0.62
5:D:402:SS7:H13	5:D:402:SS7:C13	2.27	0.61
1:A:57:LYS:O	1:A:61:LEU:HB2	2.01	0.61
1:A:64:LEU:CD2	1:A:163:LEU:HD21	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:GLN:O	1:A:162:GLN:NE2	2.36	0.59
1:C:174:PRO:HB2	1:C:205:VAL:CG1	2.30	0.59
1:C:198:HIS:HB3	1:C:214:LEU:HD11	1.84	0.59
5:D:402:SS7:C22	5:D:402:SS7:C13	2.78	0.58
1:A:10:LYS:O	1:A:14:GLU:HG2	2.04	0.57
1:A:75:GLU:HG3	1:A:129:LEU:HD11	1.85	0.57
1:B:23:THR:O	1:B:105:HIS:HE1	1.87	0.57
1:A:95:TYR:HE1	1:A:120:GLU:HG2	1.70	0.56
1:D:220:ASP:OD2	5:D:402:SS7:H4	2.06	0.56
1:C:56:ILE:HD13	1:C:230:ALA:HB3	1.90	0.54
3:B:401:THF:N10	3:B:401:THF:CP1	2.68	0.51
1:D:195:PRO:HG2	1:D:217:ARG:NH2	2.25	0.50
1:B:46:LEU:CD2	1:B:226:PRO:HB3	2.37	0.50
1:D:57:LYS:HD2	1:D:298:ILE:HB	1.93	0.50
1:A:2:GLU:HB2	1:A:278:PHE:CZ	2.47	0.50
1:A:36:ARG:HH21	1:A:249:GLU:CD	2.19	0.50
1:C:8:LEU:O	1:C:12:ILE:HG13	2.12	0.49
1:D:227:PHE:CD2	5:D:402:SS7:H14	2.47	0.49
5:D:402:SS7:C26	5:D:402:SS7:H18	2.42	0.49
1:C:2:GLU:HB2	1:C:278:PHE:CZ	2.48	0.49
1:A:68:THR:HG21	1:A:147:ALA:HB2	1.94	0.48
1:C:58:SER:CA	1:C:295:MET:HE1	2.41	0.48
1:C:62:TRP:CD1	1:C:67:ASP:HB3	2.48	0.47
1:A:64:LEU:HD22	1:A:163:LEU:HD21	1.95	0.47
1:C:220:ASP:O	1:C:225:VAL:HG23	2.14	0.47
1:B:265:GLU:OE1	1:D:265:GLU:OE1	2.32	0.47
1:A:303:TYR:CZ	1:A:305:PRO:HG3	2.49	0.47
1:C:56:ILE:HG23	1:C:231:SER:HB3	1.95	0.47
1:C:152:TRP:CH2	1:C:185:ASN:HB2	2.49	0.47
1:D:220:ASP:CB	5:D:402:SS7:H4	2.45	0.47
1:C:34:GLN:O	4:C:402:EDO:H12	2.15	0.47
1:A:131:ASP:HB3	1:A:134:PHE:HB3	1.97	0.46
1:D:198:HIS:HB2	1:D:214:LEU:HD11	1.96	0.46
1:B:234:LEU:HD11	1:B:298:ILE:HD13	1.98	0.46
1:D:203:PHE:CZ	1:D:239:ILE:HD12	2.51	0.46
1:A:51:VAL:HA	1:A:52:PRO:HD2	1.83	0.45
1:B:188:ASP:O	1:B:192:MET:HG3	2.16	0.45
1:C:313:ILE:H	1:C:313:ILE:HD12	1.81	0.45
1:A:215:TYR:CD2	1:B:200:MET:HE1	2.51	0.45
1:C:20:GLU:OE2	1:C:24:GLY:HA2	2.17	0.45
1:A:62:TRP:CD1	1:A:67:ASP:HB3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:THR:O	1:B:105:HIS:CE1	2.66	0.45
1:C:19:LYS:HB2	1:C:27:THR:HG22	1.99	0.45
1:A:68:THR:OG1	1:A:140:GLU:OE1	2.25	0.45
3:B:401:THF:HN0	3:B:401:THF:CP1	2.16	0.45
5:D:402:SS7:H9	5:D:402:SS7:H12	1.92	0.45
1:A:213:GLN:HA	1:A:251:VAL:O	2.18	0.44
1:B:284:ASN:HA	1:B:285:PRO:HD2	1.88	0.44
1:A:209:LYS:HG2	1:A:246:GLU:HB2	2.00	0.44
1:D:214:LEU:HD22	1:D:232:TYR:CD2	2.53	0.44
1:D:235:LEU:O	1:D:239:ILE:HG13	2.18	0.43
1:A:46:LEU:HD22	1:A:226:PRO:HB3	2.00	0.43
1:C:303:TYR:CZ	1:C:305:PRO:HG3	2.53	0.43
5:D:402:SS7:H18	5:D:402:SS7:H21	1.99	0.43
1:B:60:LEU:O	1:B:64:LEU:HG	2.17	0.43
1:C:36:ARG:NE	1:C:249:GLU:OE2	2.40	0.43
1:D:279:PRO:HB2	1:D:300:VAL:HG13	2.00	0.43
1:B:99:ASP:OD1	1:B:101:THR:HG23	2.18	0.43
1:D:84:TRP:HZ2	5:D:402:SS7:H1	1.75	0.43
1:C:273:ARG:NH2	1:C:307:PRO:O	2.52	0.43
1:D:192:MET:SD	1:D:196:PRO:HD3	2.59	0.43
1:C:64:LEU:HD23	1:C:64:LEU:HA	1.74	0.43
1:C:209:LYS:HB2	1:C:209:LYS:HE3	1.90	0.43
1:D:56:ILE:HD13	1:D:230:ALA:HB3	2.00	0.43
1:A:5:TYR:CE2	1:A:46:LEU:HD21	2.54	0.42
1:A:303:TYR:CE2	1:A:305:PRO:HG3	2.54	0.42
1:B:49:LYS:HD2	6:B:503:HOH:O	2.19	0.42
1:B:62:TRP:CZ3	1:B:73:LEU:HD11	2.54	0.42
1:D:303:TYR:CE2	1:D:305:PRO:HG3	2.54	0.42
1:B:265:GLU:OE1	1:D:265:GLU:CD	2.62	0.42
1:B:72:TYR:CD1	1:B:72:TYR:C	2.97	0.42
1:B:237:HIS:CD2	1:B:247:VAL:HG11	2.55	0.42
1:C:227:PHE:CD1	1:C:227:PHE:C	2.99	0.41
1:C:185:ASN:HA	1:C:186:PRO:HD3	1.90	0.41
1:C:214:LEU:HD13	1:C:232:TYR:CE2	2.55	0.41
1:D:39:LEU:HB2	1:D:247:VAL:HG13	2.02	0.41
1:D:198:HIS:CD2	1:D:198:HIS:H	2.37	0.41
1:A:12:ILE:HD11	1:A:221:VAL:HG21	2.01	0.41
1:C:183:ALA:O	1:C:196:PRO:HG3	2.20	0.41
1:C:58:SER:HA	1:C:295:MET:CE	2.46	0.41
3:B:401:THF:O4	3:B:401:THF:CP1	2.66	0.41
1:D:71:ARG:O	1:D:75:GLU:HG3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:GLY:HA3	1:C:150:ARG:HG3	2.03	0.40
1:D:58:SER:HB3	1:D:295:MET:SD	2.61	0.40
1:B:53:PHE:O	1:B:57:LYS:HG2	2.21	0.40
1:A:56:ILE:HD13	1:A:230:ALA:HB3	2.03	0.40
1:D:310:LYS:HD2	1:D:311:ALA:H	1.87	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	284/315 (90%)	270 (95%)	14 (5%)	0	100	100
1	B	313/315 (99%)	304 (97%)	8 (3%)	1 (0%)	36	50
1	C	268/315 (85%)	253 (94%)	15 (6%)	0	100	100
1	D	313/315 (99%)	303 (97%)	9 (3%)	1 (0%)	36	50
All	All	1178/1260 (94%)	1130 (96%)	46 (4%)	2 (0%)	43	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	144	ILE
1	B	144	ILE

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	254/273 (93%)	241 (95%)	13 (5%)	21	37
1	B	273/273 (100%)	264 (97%)	9 (3%)	33	55
1	C	237/273 (87%)	221 (93%)	16 (7%)	14	25
1	D	273/273 (100%)	258 (94%)	15 (6%)	19	34
All	All	1037/1092 (95%)	984 (95%)	53 (5%)	21	37

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLU
1	A	23	THR
1	A	61	LEU
1	A	143	ASN
1	A	144	ILE
1	A	151	HIS
1	A	156	ASP
1	A	194	LEU
1	A	206	ASN
1	A	214	LEU
1	A	238	LEU
1	A	273	ARG
1	A	274	GLU
1	B	29	SER
1	B	46	LEU
1	B	58	SER
1	B	61	LEU
1	B	71	ARG
1	B	198	HIS
1	B	294	ASP
1	B	299	LYS
1	B	300	VAL
1	C	20	GLU
1	C	25	THR
1	C	27	THR
1	C	76	ARG
1	C	82	ASP
1	C	134	PHE
1	C	144	ILE
1	C	151	HIS
1	C	163	LEU

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Mol	Chain	Res	Type
1	C	205	VAL
1	C	262	ASN
1	C	265	GLU
1	C	273	ARG
1	C	280	THR
1	C	290	VAL
1	C	296	GLU
1	D	55	LEU
1	D	61	LEU
1	D	128	ILE
1	D	189	VAL
1	D	198	HIS
1	D	207	GLU
1	D	235	LEU
1	D	242	GLU
1	D	247	VAL
1	D	254	LEU
1	D	273	ARG
1	D	280	THR
1	D	294	ASP
1	D	300	VAL
1	D	301	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	198	HIS
1	A	206	ASN
1	A	262	ASN
1	A	268	GLN
1	A	270	GLN
1	B	237	HIS
1	B	268	GLN
1	C	34	GLN
1	C	261	GLN
1	D	122	GLN
1	D	261	GLN
1	D	268	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

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$
3	THF	B	401	-	32,36,36	1.40	1 (3%)	37,50,50	1.86	7 (18%)
4	EDO	D	403	-	3,3,3	0.42	0	2,2,2	0.43	0
2	SO4	C	401	-	4,4,4	0.44	0	6,6,6	0.15	0
4	EDO	C	402	-	3,3,3	0.63	0	2,2,2	0.07	0
2	SO4	A	401	-	4,4,4	0.38	0	6,6,6	0.34	0
5	SS7	D	402	-	29,29,29	2.87	6 (20%)	41,42,42	3.10	19 (46%)
2	SO4	A	402	-	4,4,4	0.45	0	6,6,6	0.17	0
2	SO4	D	401	-	4,4,4	0.49	0	6,6,6	0.39	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	THF	B	401	-	-	5/22/37/37	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	D	403	-	-	0/1/1/1	-
5	SS7	D	402	-	-	11/13/29/29	0/3/3/3
4	EDO	C	402	-	-	1/1/1/1	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	402	SS7	O10-C7	12.42	1.47	1.22
5	D	402	SS7	C5-C8	-6.01	1.39	1.48
3	B	401	THF	C7-N8	5.26	1.34	1.28
5	D	402	SS7	C8-N9	-3.46	1.33	1.40
5	D	402	SS7	O24-C25	2.90	1.46	1.36
5	D	402	SS7	O11-C8	2.54	1.27	1.22
5	D	402	SS7	C7-N9	-2.41	1.35	1.40

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	402	SS7	C5-C8-N9	8.83	115.72	105.96
5	D	402	SS7	C7-N9-C8	-8.00	102.47	111.55
3	B	401	THF	C7-N8-C8A	6.96	121.69	115.47
5	D	402	SS7	C6-C7-N9	6.18	112.79	105.96
5	D	402	SS7	C5-C6-C7	-5.67	103.29	108.26
5	D	402	SS7	O24-C25-C26	5.29	122.43	110.97
5	D	402	SS7	O10-C7-C6	-4.96	119.06	128.66
5	D	402	SS7	O24-C25-O27	-3.73	114.37	122.43
3	B	401	THF	C7-C6-N5	3.53	114.80	109.81
3	B	401	THF	C9-N10-C14	3.37	130.58	122.00
5	D	402	SS7	C19-O24-C25	3.27	130.55	118.27
5	D	402	SS7	C20-C21-C16	3.27	122.67	117.94
3	B	401	THF	C11-C-N	3.22	123.00	117.04
5	D	402	SS7	O15-C13-C12	2.99	115.17	108.41
5	D	402	SS7	O11-C8-C5	-2.99	122.87	128.66
5	D	402	SS7	C6-C5-C8	-2.88	105.73	108.26
5	D	402	SS7	C12-N9-C7	2.79	132.07	123.99
3	B	401	THF	CT-CA-N	2.78	117.02	110.57
5	D	402	SS7	C4-C5-C8	2.63	134.00	129.59
5	D	402	SS7	O11-C8-N9	-2.55	121.41	124.99
5	D	402	SS7	O15-C16-C17	2.39	123.33	118.90
3	B	401	THF	O-C-C11	-2.39	116.16	120.90
5	D	402	SS7	C1-C6-C7	2.34	133.52	129.59
5	D	402	SS7	O10-C7-N9	2.25	128.16	124.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	THF	O3-CP1-N5	-2.19	106.62	112.31
5	D	402	SS7	C13-O15-C16	2.19	121.98	114.50

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	401	THF	N5-C6-C9-N10
3	B	401	THF	C7-C6-C9-N10
5	D	402	SS7	C14-C12-N9-C8
5	D	402	SS7	C14-C12-N9-C7
5	D	402	SS7	N9-C12-C13-O15
5	D	402	SS7	C14-C12-C13-O15
5	D	402	SS7	C12-C13-O15-C16
5	D	402	SS7	C17-C16-O15-C13
5	D	402	SS7	C21-C16-O15-C13
5	D	402	SS7	C26-C25-O24-C19
5	D	402	SS7	O27-C25-O24-C19
4	C	402	EDO	O1-C1-C2-O2
5	D	402	SS7	C20-C19-O24-C25
5	D	402	SS7	C18-C19-O24-C25
3	B	401	THF	C6-C9-N10-C14
3	B	401	THF	OE1-CD-CG-CB
3	B	401	THF	OE2-CD-CG-CB

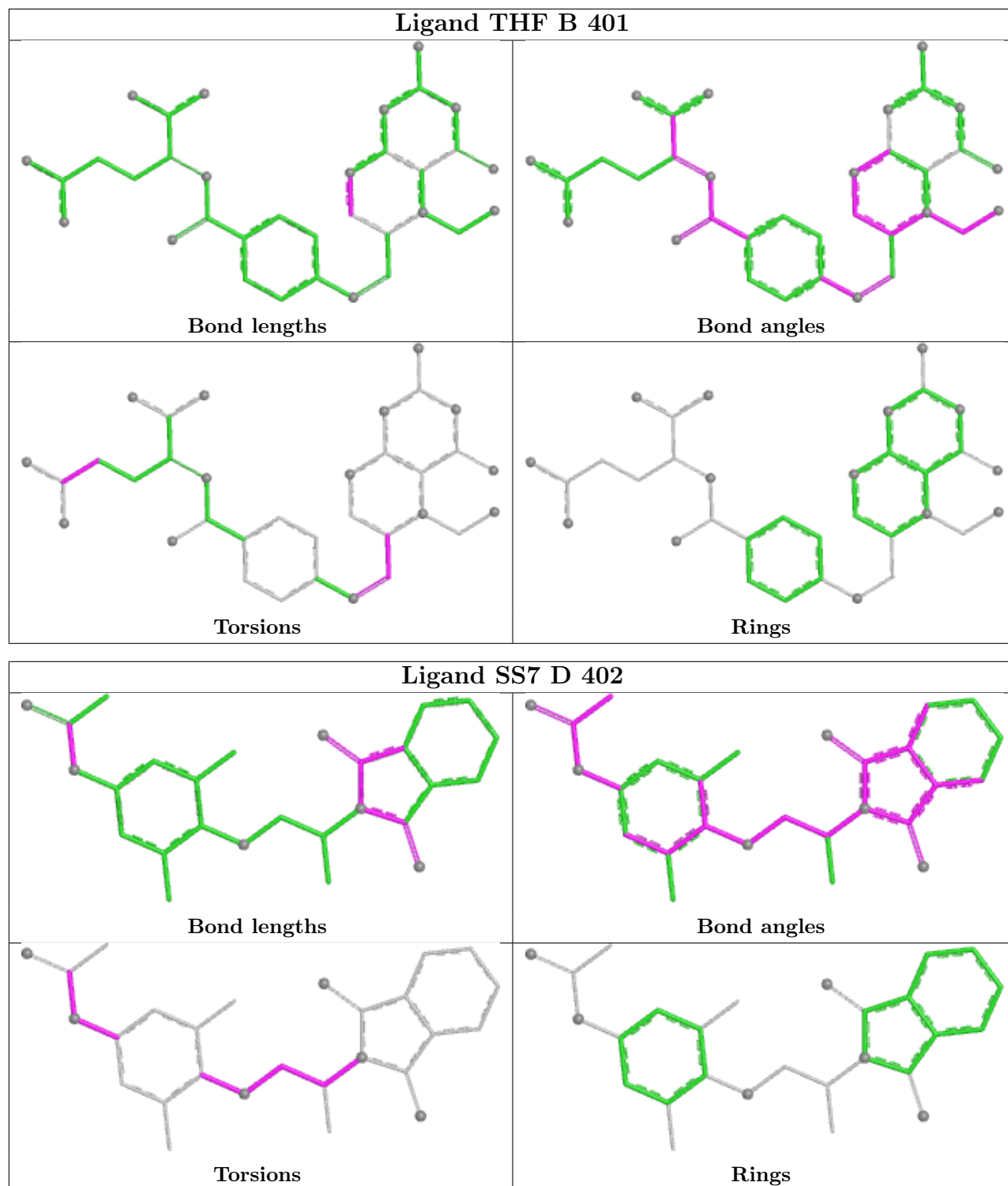
There are no ring outliers.

3 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	THF	6	0
4	C	402	EDO	1	0
5	D	402	SS7	13	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.



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, 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	290/315 (92%)	0.37	16 (5%) 30 27	7, 15, 56, 77	0
1	B	315/315 (100%)	0.04	0 100 100	5, 9, 21, 29	0
1	C	272/315 (86%)	0.35	18 (6%) 24 20	6, 14, 49, 67	0
1	D	315/315 (100%)	0.00	2 (0%) 85 83	4, 9, 28, 39	0
All	All	1192/1260 (94%)	0.18	36 (3%) 52 48	4, 11, 44, 77	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	90	VAL	4.1
1	A	132	ALA	4.1
1	A	117	TYR	4.1
1	A	84	TRP	3.9
1	C	142	GLY	3.8
1	C	89	TYR	3.6
1	C	81	TRP	3.6
1	C	135	ALA	3.4
1	A	85	ALA	3.3
1	A	118	LYS	3.3
1	C	80	ILE	3.2
1	A	95	TYR	3.2
1	A	120	GLU	2.7
1	C	84	TRP	2.6
1	A	313	ILE	2.6
1	C	82	ASP	2.5
1	C	133	GLU	2.5
1	C	83	GLU	2.5
1	C	288	ALA	2.5
1	C	144	ILE	2.4
1	C	143	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	141	LEU	2.3
1	A	138	TYR	2.2
1	D	126	ASP	2.2
1	C	164	ALA	2.2
1	C	134	PHE	2.2
1	C	85	ALA	2.1
1	A	116	GLN	2.1
1	C	138	TYR	2.1
1	D	102	ASP	2.1
1	A	81	TRP	2.1
1	A	88	ARG	2.1
1	A	133	GLU	2.1
1	C	90	VAL	2.0
1	A	127	ALA	2.0
1	A	142	GLY	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](#)

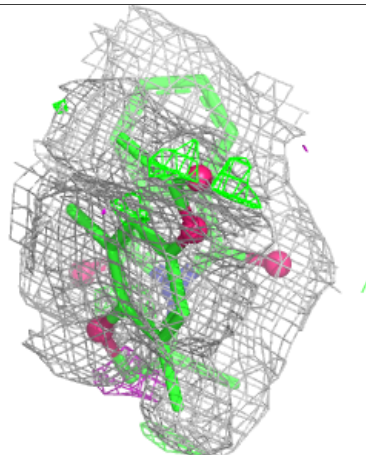
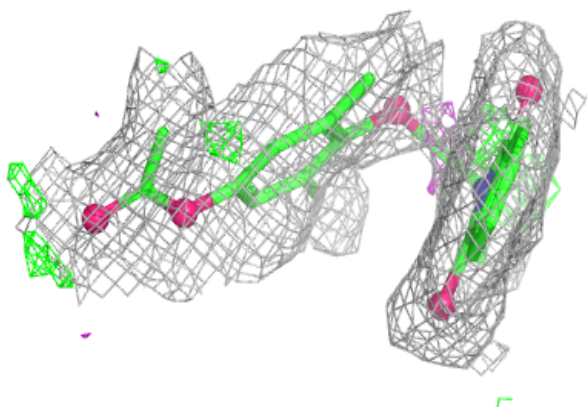
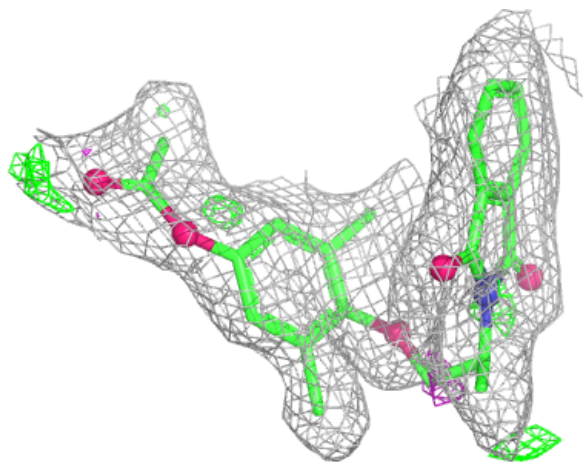
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

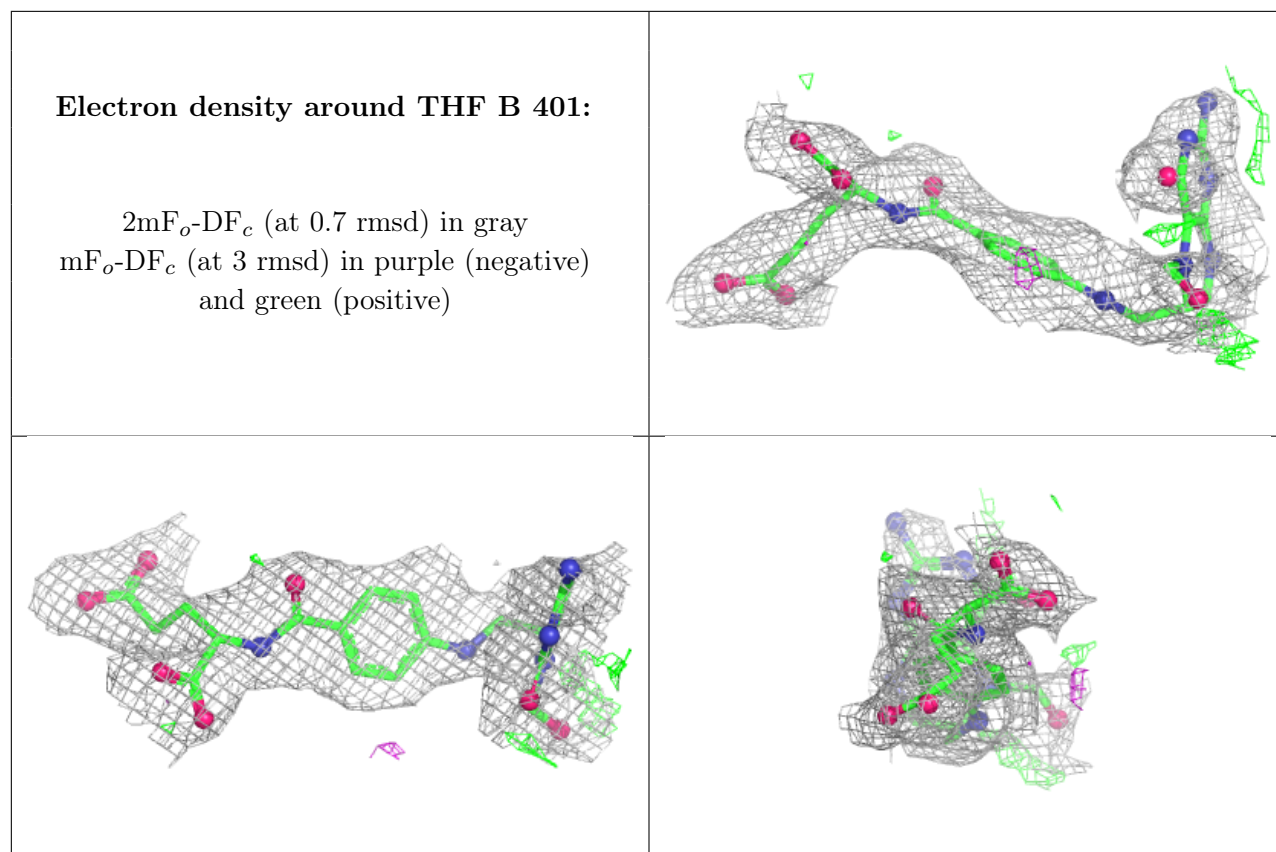
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	SS7	D	402	27/27	0.86	0.13	10,11,13,13	0
4	EDO	C	402	4/4	0.92	0.10	7,7,7,7	0
2	SO4	A	402	5/5	0.92	0.13	34,35,36,36	0
3	THF	B	401	34/34	0.94	0.09	10,10,11,12	0
4	EDO	D	403	4/4	0.95	0.08	17,17,17,18	0
2	SO4	C	401	5/5	0.95	0.10	30,30,31,32	0
2	SO4	A	401	5/5	0.98	0.08	13,13,14,14	0
2	SO4	D	401	5/5	0.99	0.07	10,10,10,10	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 SS7 D 402:

$2mF_o-DF_c$ (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.