



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2021 – 04:44 PM EST

PDB ID : 5Q05
Title : Human liver fructose-1,6-bisphosphatase 1 (fructose 1,6-bisphosphate 1-phosphatase, E.C.3.1.3.11) complexed with the allosteric inhibitor 1-(5-bromo-1,3-thiazol-2-yl)-3-[5-(2-methoxyethyl)-4-methylthiophen-2-yl]sulfonyleurea
Authors : Ruf, A.; Joseph, C.; Alker, A.; Banner, D.; Tetaz, T.; Benz, J.; Kuhn, B.; Rudolph, M.G.; Yang, H.; Shao, C.; Burley, S.K.
Deposited on : 2017-04-20
Resolution : 2.20 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 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)
Xtrriage (Phenix) : 1.13
EDS : 2.16
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) : 2.16

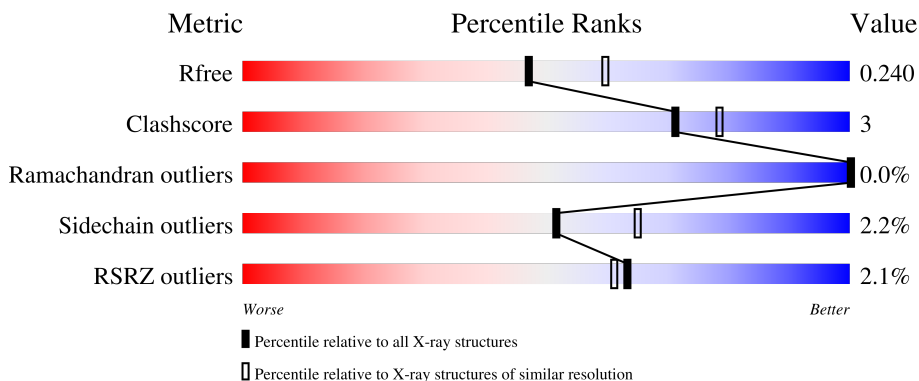
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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.



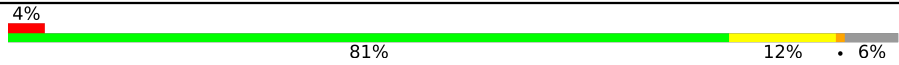


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	338	
1	B	338	
1	C	338	
1	D	338	
1	E	338	

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Mol	Chain	Length	Quality of chain
1	F	338	 4% 81% 12% • 6%
1	G	338	 % 86% 7% • 6%
1	H	338	 % 86% 7% 6%

2 Entry composition

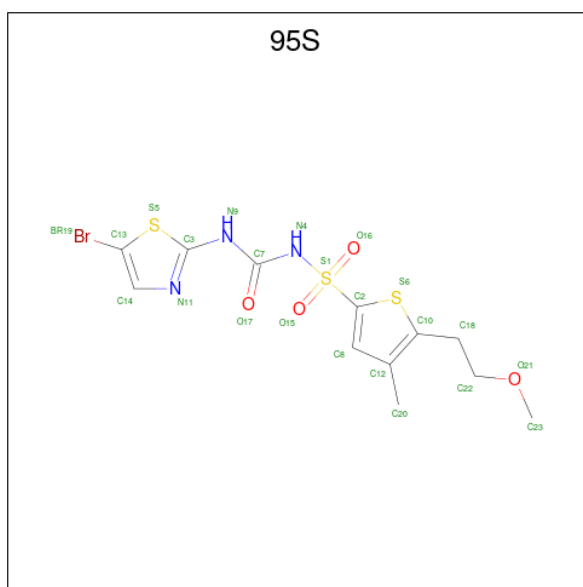
There are 3 unique types of molecules in this entry. The entry contains 20555 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fructose-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	319	2446	1558	412	459	17	0	1	0
1	B	317	2431	1548	409	457	17	0	1	0
1	C	317	2425	1545	406	457	17	0	0	0
1	D	317	2431	1548	409	457	17	0	1	0
1	E	319	2447	1559	412	459	17	0	1	0
1	F	317	2430	1547	409	457	17	0	1	0
1	G	317	2430	1547	409	457	17	0	1	0
1	H	317	2425	1545	406	457	17	0	0	0

- Molecule 2 is N-[(5-bromo-1,3-thiazol-2-yl)carbamoyl]-5-(2-methoxyethyl)-4-methylthiophene-2-sulfonamide (three-letter code: 95S) (formula: C₁₂H₁₄BrN₃O₄S₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	Br	C	N	O			S
2	A	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	B	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	C	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	D	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	E	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	F	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	G	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0
2	H	1	Total 23	Br 1	C 12	N 3	O 4	S 3	0	0

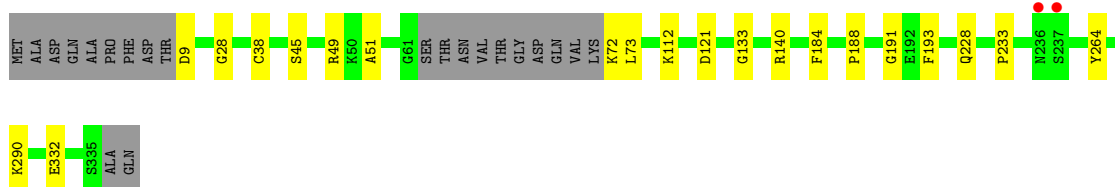
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	103	Total 103	O 103	0	0
3	B	106	Total 106	O 106	0	0
3	C	142	Total 142	O 142	0	0
3	D	139	Total 139	O 139	0	0

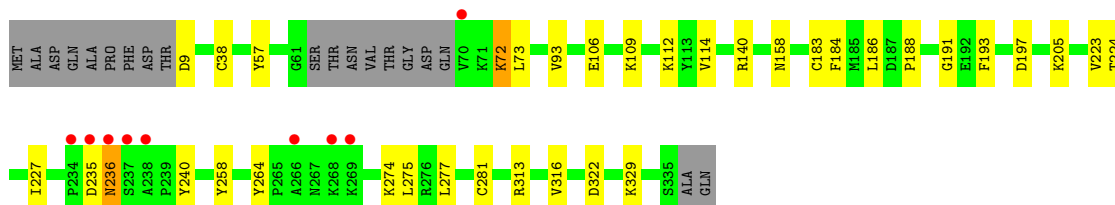
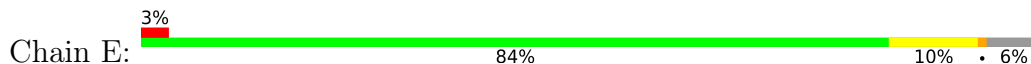
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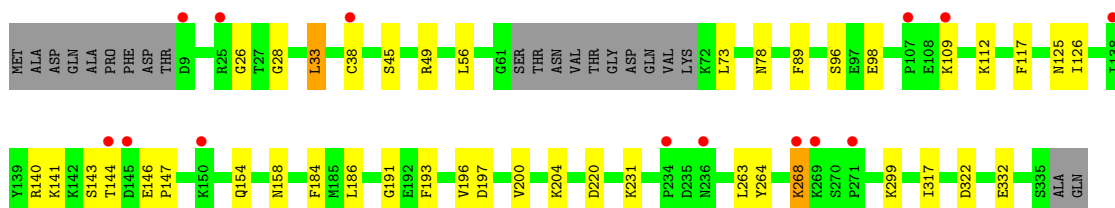
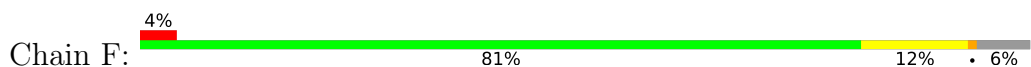
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	75	Total 75	O 75	0	0
3	F	104	Total 104	O 104	0	0
3	G	123	Total 123	O 123	0	0
3	H	114	Total 114	O 114	0	0



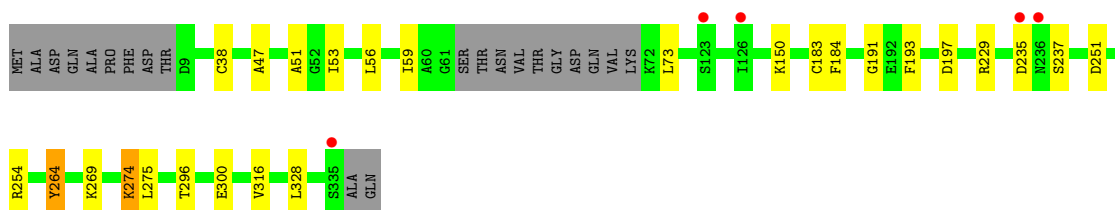
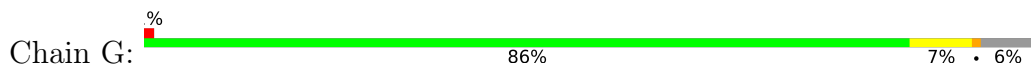
- Molecule 1: Fructose-1,6-bisphosphatase 1



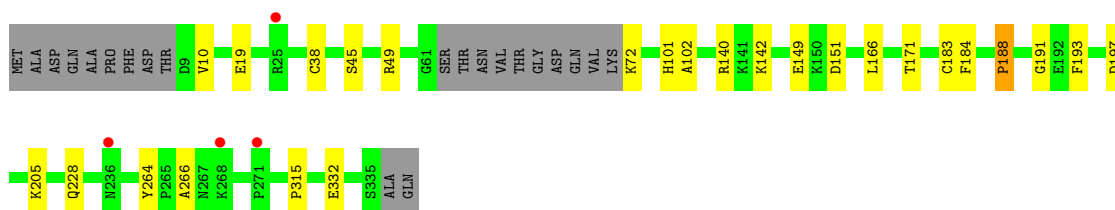
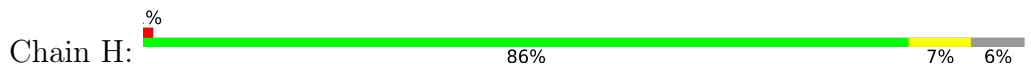
- Molecule 1: Fructose-1,6-bisphosphatase 1



- Molecule 1: Fructose-1,6-bisphosphatase 1



- Molecule 1: Fructose-1,6-bisphosphatase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.73Å 285.43Å 83.91Å 90.00° 97.47° 90.00°	Depositor
Resolution (Å)	29.12 – 2.20 29.12 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.1 (29.12-2.20) 96.1 (29.12-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.187 , 0.240 0.187 , 0.240	Depositor DCC
R_{free} test set	7552 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtrriage
Anisotropy	0.398	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20555	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: 95S

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	0.40	0/2495	0.59	1/3368 (0.0%)
1	B	0.40	0/2479	0.60	1/3347 (0.0%)
1	C	0.41	0/2468	0.60	0/3333
1	D	0.45	0/2479	0.62	0/3347
1	E	0.37	0/2495	0.56	0/3368
1	F	0.37	0/2479	0.57	1/3347 (0.0%)
1	G	0.41	0/2479	0.58	0/3347
1	H	0.43	0/2468	0.60	0/3333
All	All	0.40	0/19842	0.59	3/26790 (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	A	33	LEU	CA-CB-CG	6.92	131.22	115.30
1	B	33	LEU	CA-CB-CG	6.31	129.81	115.30
1	F	33	LEU	CA-CB-CG	5.42	127.77	115.30

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	2446	0	2496	21	1
1	B	2431	0	2478	21	0
1	C	2425	0	2469	16	0
1	D	2431	0	2478	13	0
1	E	2447	0	2500	23	0
1	F	2430	0	2474	23	1
1	G	2430	0	2474	15	0
1	H	2425	0	2469	16	0
2	A	23	0	0	1	0
2	B	23	0	0	1	0
2	C	23	0	0	1	0
2	D	23	0	0	1	0
2	E	23	0	0	1	0
2	F	23	0	0	2	0
2	G	23	0	0	0	0
2	H	23	0	0	0	0
3	A	103	0	0	6	0
3	B	106	0	0	1	0
3	C	142	0	0	4	0
3	D	139	0	0	4	0
3	E	75	0	0	5	0
3	F	104	0	0	4	0
3	G	123	0	0	3	0
3	H	114	0	0	5	0
All	All	20555	0	19838	138	1

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 3.

All (138) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:251:ASP:OD1	1:C:254:ARG:NH2	2.17	0.77
1:B:102:ALA:HB3	1:B:149:GLU:HG2	1.71	0.71
1:F:220:ASP:OD1	1:F:268:LYS:NZ	2.23	0.71
1:E:158:ASN:O	3:E:501:HOH:O	2.10	0.69
1:B:191:GLY:HA3	1:D:191:GLY:HA3	1.76	0.68
1:D:72:LYS:N	3:D:504:HOH:O	2.25	0.68
1:F:197:ASP:OD2	3:F:501:HOH:O	2.11	0.68
1:F:146:GLU:HG3	1:F:147:PRO:HD2	1.76	0.67
1:A:276:ARG:NH1	1:A:311:HIS:O	2.28	0.66
1:F:231:LYS:O	3:F:502:HOH:O	2.13	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:53:ILE:HA	1:G:56:LEU:HD23	1.77	0.66
1:B:109:LYS:NZ	1:D:9:ASP:OD2	2.28	0.65
1:D:51:ALA:O	3:D:501:HOH:O	2.14	0.65
1:G:183:CYS:HB2	1:G:197:ASP:HB2	1.78	0.64
1:G:251:ASP:OD1	1:G:254:ARG:NH2	2.26	0.64
1:A:29:GLU:OE2	3:A:502:HOH:O	2.14	0.63
1:H:140:ARG:NH1	3:H:506:HOH:O	2.31	0.63
1:G:296:THR:HG21	1:G:328:LEU:HD21	1.81	0.63
1:A:79:ASP:OD1	3:A:503:HOH:O	2.16	0.61
1:H:151:ASP:OD1	3:H:502:HOH:O	2.16	0.61
1:E:183:CYS:HB2	1:E:197:ASP:HB2	1.82	0.61
1:A:264:TYR:OH	1:A:274:LYS:HD3	1.99	0.60
1:B:225:GLU:OE1	1:B:334:HIS:NE2	2.31	0.60
1:B:141:LYS:NZ	1:B:143:SER:OG	2.33	0.59
1:A:51:ALA:HA	1:B:188:PRO:HD2	1.85	0.58
1:E:224:THR:OG1	3:E:502:HOH:O	2.17	0.56
1:F:73:LEU:HD23	1:F:126:ILE:HD13	1.87	0.56
1:H:183:CYS:HB2	1:H:197:ASP:HB2	1.88	0.55
1:A:188:PRO:HD2	1:B:51:ALA:HA	1.87	0.55
1:F:96:SER:HB2	1:F:117:PHE:CZ	2.42	0.54
1:E:191:GLY:HA3	1:G:191:GLY:HA3	1.90	0.54
1:C:51:ALA:HA	1:D:188:PRO:HD2	1.90	0.54
1:H:142:LYS:NZ	3:H:513:HOH:O	2.40	0.53
1:A:191:GLY:HA3	1:C:191:GLY:HA3	1.89	0.53
1:C:217:LYS:NZ	3:C:506:HOH:O	2.42	0.53
1:G:235:ASP:O	1:G:237:SER:N	2.37	0.52
1:H:102:ALA:HB3	1:H:149:GLU:HG2	1.91	0.52
1:H:266:ALA:HB2	1:H:315:PRO:HG3	1.92	0.52
1:G:275:LEU:HD12	1:G:316:VAL:HG11	1.90	0.51
1:B:296:THR:HG21	1:B:328:LEU:HD21	1.93	0.51
1:F:78:ASN:HB2	1:F:98:GLU:HG3	1.92	0.51
1:B:112:LYS:HD2	1:B:140[B]:ARG:NE	2.26	0.51
1:B:112:LYS:HD2	1:B:140[B]:ARG:HE	1.76	0.51
1:D:184:PHE:HB3	1:D:193:PHE:HB3	1.93	0.51
1:E:277:LEU:HA	1:E:281:CYS:HB2	1.93	0.51
1:B:22:ARG:HG2	2:B:401:95S:C14	2.41	0.51
1:C:269:LYS:NZ	3:C:508:HOH:O	2.43	0.50
1:D:121:ASP:N	1:D:133:GLY:O	2.42	0.50
1:E:106:GLU:OE2	1:E:109:LYS:NZ	2.32	0.50
1:E:184:PHE:HB3	1:E:193:PHE:HB3	1.93	0.50
1:B:140[B]:ARG:NH2	3:B:511:HOH:O	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:GLY:HA2	2:D:401:95S:C7	2.42	0.49
1:F:140[A]:ARG:NH2	3:F:515:HOH:O	2.45	0.49
1:H:266:ALA:N	1:H:315:PRO:HB3	2.27	0.49
1:D:332:GLU:OE1	3:D:503:HOH:O	2.19	0.49
1:E:274:LYS:O	1:E:313:ARG:NH1	2.45	0.49
1:C:183:CYS:HB2	1:C:197:ASP:HB2	1.95	0.49
1:D:45:SER:O	1:D:49:ARG:HD3	2.13	0.49
1:A:50:LYS:NZ	3:A:501:HOH:O	2.12	0.49
1:C:72:LYS:N	3:C:510:HOH:O	2.45	0.48
1:C:45:SER:O	1:C:49:ARG:HD3	2.14	0.48
1:B:184:PHE:HB3	1:B:193:PHE:HB3	1.95	0.48
1:E:72:LYS:NZ	3:E:516:HOH:O	2.45	0.48
1:G:269:LYS:NZ	3:G:512:HOH:O	2.46	0.48
1:F:141:LYS:NZ	1:F:143:SER:OG	2.46	0.48
1:A:166:LEU:HD13	1:A:249:VAL:HG12	1.96	0.48
1:E:223:VAL:O	1:E:227:ILE:HG12	2.14	0.48
1:A:229:ARG:NH1	1:A:234:PRO:HG3	2.29	0.47
1:F:154:GLN:HB3	1:F:158:ASN:HD22	1.78	0.47
1:D:290:LYS:NZ	3:D:508:HOH:O	2.45	0.47
1:E:322:ASP:OD1	3:E:503:HOH:O	2.20	0.47
1:E:9:ASP:N	3:E:517:HOH:O	2.47	0.47
1:E:274:LYS:HA	1:E:313:ARG:HD3	1.97	0.47
1:F:89:PHE:HA	1:F:109:LYS:O	2.15	0.46
1:F:191:GLY:HA3	1:H:191:GLY:HA3	1.97	0.46
1:F:186:LEU:HB2	1:F:193:PHE:CE1	2.50	0.46
1:G:51:ALA:HA	1:H:188:PRO:HD2	1.97	0.46
1:A:55:HIS:HA	1:A:59:ILE:HG22	1.98	0.46
1:F:28:GLY:HA2	2:F:401:95S:C7	2.45	0.46
1:G:59:ILE:HA	1:H:10:VAL:HG11	1.98	0.45
1:E:57:TYR:CZ	1:F:196:VAL:HG12	2.52	0.45
1:H:184:PHE:HB3	1:H:193:PHE:HB3	1.98	0.45
1:B:206:GLY:HA3	1:B:260:GLY:N	2.31	0.45
1:A:205:LYS:NZ	3:A:513:HOH:O	2.44	0.45
1:B:277:LEU:HA	1:B:281:CYS:HB2	1.97	0.45
1:E:93:VAL:HB	1:E:114:VAL:HG13	1.99	0.45
1:E:258:TYR:OH	1:F:125:ASN:HA	2.17	0.45
1:C:270:SER:HB3	1:C:273:GLY:O	2.17	0.45
1:F:112:LYS:NZ	3:F:520:HOH:O	2.49	0.45
1:B:195:LEU:HD21	1:B:198:LYS:HG2	1.98	0.45
3:C:506:HOH:O	1:D:233:PRO:HD2	2.16	0.45
1:A:112:LYS:HE3	1:A:140[A]:ARG:HH21	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:228:GLN:NE2	3:H:512:HOH:O	2.39	0.44
1:H:166:LEU:O	1:H:171:THR:HA	2.18	0.44
1:F:204:LYS:NZ	1:F:322:ASP:OD2	2.36	0.44
1:A:103:ILE:N	1:A:103:ILE:HD12	2.33	0.44
1:E:112:LYS:HB2	1:E:140[B]:ARG:NH1	2.32	0.44
1:F:299:LYS:HB2	1:F:299:LYS:HE3	1.76	0.44
1:H:45:SER:O	1:H:49:ARG:HD3	2.18	0.44
1:E:275:LEU:HD12	1:E:316:VAL:HG11	2.00	0.44
1:G:47:ALA:HB1	1:G:73:LEU:HD11	2.00	0.44
1:E:235:ASP:O	1:E:236:ASN:HB2	2.18	0.43
1:G:229:ARG:HD2	3:G:501:HOH:O	2.18	0.43
1:A:27:THR:HG23	3:A:579:HOH:O	2.18	0.43
1:B:73:LEU:HD12	1:B:73:LEU:HA	1.72	0.43
1:A:272:ASN:HA	3:A:514:HOH:O	2.19	0.43
1:D:112:LYS:HB2	1:D:140[B]:ARG:NH1	2.33	0.43
1:E:205:LYS:NZ	1:E:240:TYR:OH	2.35	0.43
1:G:264:TYR:CZ	1:G:274:LYS:HG2	2.54	0.43
1:A:140[B]:ARG:HH11	1:A:142:LYS:HD3	1.82	0.43
1:F:184:PHE:HB3	1:F:193:PHE:HB3	1.99	0.43
1:B:141:LYS:HE3	1:B:143:SER:O	2.19	0.42
1:C:120:LEU:HA	1:C:120:LEU:HD12	1.81	0.42
1:B:278:LEU:HD23	1:B:278:LEU:HA	1.75	0.42
1:C:55:HIS:HA	1:C:59:ILE:HG22	2.00	0.42
1:B:276:ARG:NH1	1:B:311:HIS:O	2.50	0.42
2:E:401:95S:O17	2:E:401:95S:S5	2.76	0.42
1:F:26:GLY:HA3	2:F:401:95S:O15	2.19	0.42
1:H:205:LYS:N	3:H:504:HOH:O	2.27	0.42
1:H:72:LYS:HE3	1:H:72:LYS:HB2	1.78	0.42
1:A:112:LYS:HE3	1:A:140[A]:ARG:NH2	2.34	0.42
1:A:37:LEU:HD21	1:A:136:PHE:CD2	2.54	0.42
1:C:53:ILE:HA	1:C:56:LEU:HD23	2.02	0.42
2:A:401:95S:S5	2:A:401:95S:O17	2.77	0.41
1:E:112:LYS:HD3	1:E:140[A]:ARG:NE	2.35	0.41
1:F:45:SER:O	1:F:49:ARG:HD3	2.20	0.41
1:F:263:LEU:HG	1:F:317:ILE:HG23	2.02	0.41
1:A:73:LEU:HD22	1:A:120:LEU:HD22	2.02	0.41
1:A:223:VAL:O	1:A:227:ILE:HG12	2.21	0.41
1:C:28:GLY:HA2	2:C:401:95S:C7	2.51	0.41
1:E:186:LEU:HB2	1:E:193:PHE:CE1	2.56	0.41
1:C:184:PHE:HB3	1:C:193:PHE:HB3	2.01	0.40
1:G:184:PHE:HB3	1:G:193:PHE:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:CYS:HB2	1:B:197:ASP:HB2	2.04	0.40
1:G:229:ARG:NH1	3:G:501:HOH:O	2.21	0.40
1:C:216:ALA:HA	1:C:219:PHE:CD2	2.57	0.40
1:E:329:LYS:HA	1:E:329:LYS:HD3	1.77	0.40
1:C:92:CYS:SG	1:C:141:LYS:HD2	2.62	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:ARG:NH2	1:F:332:GLU:OE1[1_655]	2.16	0.04

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	316/338 (94%)	309 (98%)	7 (2%)	0	100	100
1	B	314/338 (93%)	304 (97%)	10 (3%)	0	100	100
1	C	313/338 (93%)	307 (98%)	6 (2%)	0	100	100
1	D	314/338 (93%)	306 (98%)	8 (2%)	0	100	100
1	E	316/338 (94%)	307 (97%)	8 (2%)	1 (0%)	41	46
1	F	314/338 (93%)	304 (97%)	10 (3%)	0	100	100
1	G	314/338 (93%)	306 (98%)	8 (2%)	0	100	100
1	H	313/338 (93%)	306 (98%)	7 (2%)	0	100	100
All	All	2514/2704 (93%)	2449 (97%)	64 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	236	ASN

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	267/281 (95%)	260 (97%)	7 (3%)	46	58
1	B	265/281 (94%)	255 (96%)	10 (4%)	33	42
1	C	264/281 (94%)	261 (99%)	3 (1%)	73	85
1	D	265/281 (94%)	261 (98%)	4 (2%)	65	78
1	E	267/281 (95%)	262 (98%)	5 (2%)	57	71
1	F	265/281 (94%)	258 (97%)	7 (3%)	46	58
1	G	265/281 (94%)	260 (98%)	5 (2%)	57	71
1	H	264/281 (94%)	258 (98%)	6 (2%)	50	63
All	All	2122/2248 (94%)	2075 (98%)	47 (2%)	52	65

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	38	CYS
1	A	73	LEU
1	A	140[A]	ARG
1	A	140[B]	ARG
1	A	228	GLN
1	A	264	TYR
1	B	33	LEU
1	B	38	CYS
1	B	56	LEU
1	B	144	THR
1	B	200	VAL
1	B	228	GLN
1	B	236	ASN
1	B	264	TYR

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Mol	Chain	Res	Type
1	B	268	LYS
1	B	299	LYS
1	C	38	CYS
1	C	140	ARG
1	C	264	TYR
1	D	38	CYS
1	D	73	LEU
1	D	228	GLN
1	D	264	TYR
1	E	38	CYS
1	E	72	LYS
1	E	73	LEU
1	E	188	PRO
1	E	264	TYR
1	F	33	LEU
1	F	38	CYS
1	F	56	LEU
1	F	144	THR
1	F	200	VAL
1	F	264	TYR
1	F	268	LYS
1	G	38	CYS
1	G	150	LYS
1	G	264	TYR
1	G	274	LYS
1	G	300	GLU
1	H	19	GLU
1	H	38	CYS
1	H	101	HIS
1	H	188	PRO
1	H	264	TYR
1	H	332	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	236	ASN
1	C	55	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
2	95S	B	401	-	19,24,24	1.02	1 (5%)	14,34,34	1.10	2 (14%)
2	95S	H	401	-	19,24,24	1.04	1 (5%)	14,34,34	1.12	2 (14%)
2	95S	D	401	-	19,24,24	0.72	0	14,34,34	1.25	2 (14%)
2	95S	G	401	-	19,24,24	1.54	3 (15%)	14,34,34	1.56	2 (14%)
2	95S	E	401	-	19,24,24	1.29	3 (15%)	14,34,34	1.59	2 (14%)
2	95S	C	401	-	19,24,24	1.16	2 (10%)	14,34,34	1.31	2 (14%)
2	95S	F	401	-	19,24,24	1.29	3 (15%)	14,34,34	1.32	2 (14%)
2	95S	A	401	-	19,24,24	1.12	1 (5%)	14,34,34	1.58	2 (14%)

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
2	95S	B	401	-	-	3/10/19/19	0/2/2/2
2	95S	H	401	-	-	4/10/19/19	0/2/2/2
2	95S	D	401	-	-	2/10/19/19	0/2/2/2
2	95S	G	401	-	-	2/10/19/19	0/2/2/2
2	95S	E	401	-	-	2/10/19/19	0/2/2/2
2	95S	C	401	-	-	2/10/19/19	0/2/2/2
2	95S	F	401	-	-	3/10/19/19	0/2/2/2
2	95S	A	401	-	-	3/10/19/19	0/2/2/2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	401	95S	C2-S1	-4.42	1.69	1.76
2	C	401	95S	C3-N9	2.90	1.41	1.36
2	E	401	95S	C18-C22	2.86	1.60	1.50
2	C	401	95S	C2-S1	-2.85	1.72	1.76
2	B	401	95S	C2-S1	-2.68	1.72	1.76
2	H	401	95S	C2-S1	-2.46	1.72	1.76
2	A	401	95S	C3-N9	2.43	1.41	1.36
2	F	401	95S	C2-S1	-2.41	1.72	1.76
2	F	401	95S	C18-C22	-2.36	1.43	1.50
2	G	401	95S	C2-S6	-2.33	1.68	1.72
2	F	401	95S	C18-C10	-2.30	1.49	1.50
2	E	401	95S	C18-C10	-2.30	1.49	1.50
2	E	401	95S	C8-C12	2.21	1.44	1.38
2	G	401	95S	C3-N9	2.07	1.40	1.36

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	401	95S	C2-S1-N4	5.04	113.16	107.27
2	G	401	95S	C2-S1-N4	4.68	112.74	107.27
2	A	401	95S	C2-S1-N4	4.65	112.70	107.27
2	C	401	95S	C2-S1-N4	3.79	111.70	107.27
2	D	401	95S	C2-S1-N4	3.48	111.33	107.27
2	H	401	95S	C2-S1-N4	2.97	110.74	107.27
2	F	401	95S	C2-S1-N4	2.91	110.67	107.27
2	A	401	95S	BR19-C13-C14	2.60	128.61	124.51
2	F	401	95S	BR19-C13-C14	2.44	128.35	124.51
2	G	401	95S	BR19-C13-C14	2.34	128.19	124.51
2	E	401	95S	BR19-C13-C14	2.33	128.18	124.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	95S	BR19-C13-C14	2.31	128.14	124.51
2	C	401	95S	BR19-C13-C14	2.24	128.03	124.51
2	D	401	95S	BR19-C13-C14	2.21	128.00	124.51
2	B	401	95S	C2-S1-N4	2.20	109.84	107.27
2	H	401	95S	BR19-C13-C14	2.14	127.89	124.51

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	95S	N4-C7-N9-C3
2	H	401	95S	N4-C7-N9-C3
2	H	401	95S	C12-C10-C18-C22
2	H	401	95S	C10-C18-C22-O21
2	D	401	95S	N4-C7-N9-C3
2	G	401	95S	N4-C7-N9-C3
2	E	401	95S	N4-C7-N9-C3
2	C	401	95S	N4-C7-N9-C3
2	C	401	95S	C12-C10-C18-C22
2	F	401	95S	N4-C7-N9-C3
2	F	401	95S	C12-C10-C18-C22
2	A	401	95S	N4-C7-N9-C3
2	B	401	95S	C10-C18-C22-O21
2	A	401	95S	C18-C22-O21-C23
2	B	401	95S	C18-C22-O21-C23
2	H	401	95S	C18-C22-O21-C23
2	E	401	95S	C10-C18-C22-O21
2	G	401	95S	C18-C22-O21-C23
2	A	401	95S	C12-C10-C18-C22
2	D	401	95S	C10-C18-C22-O21
2	F	401	95S	C18-C22-O21-C23

There are no ring outliers.

6 monomers are involved in 7 short contacts:

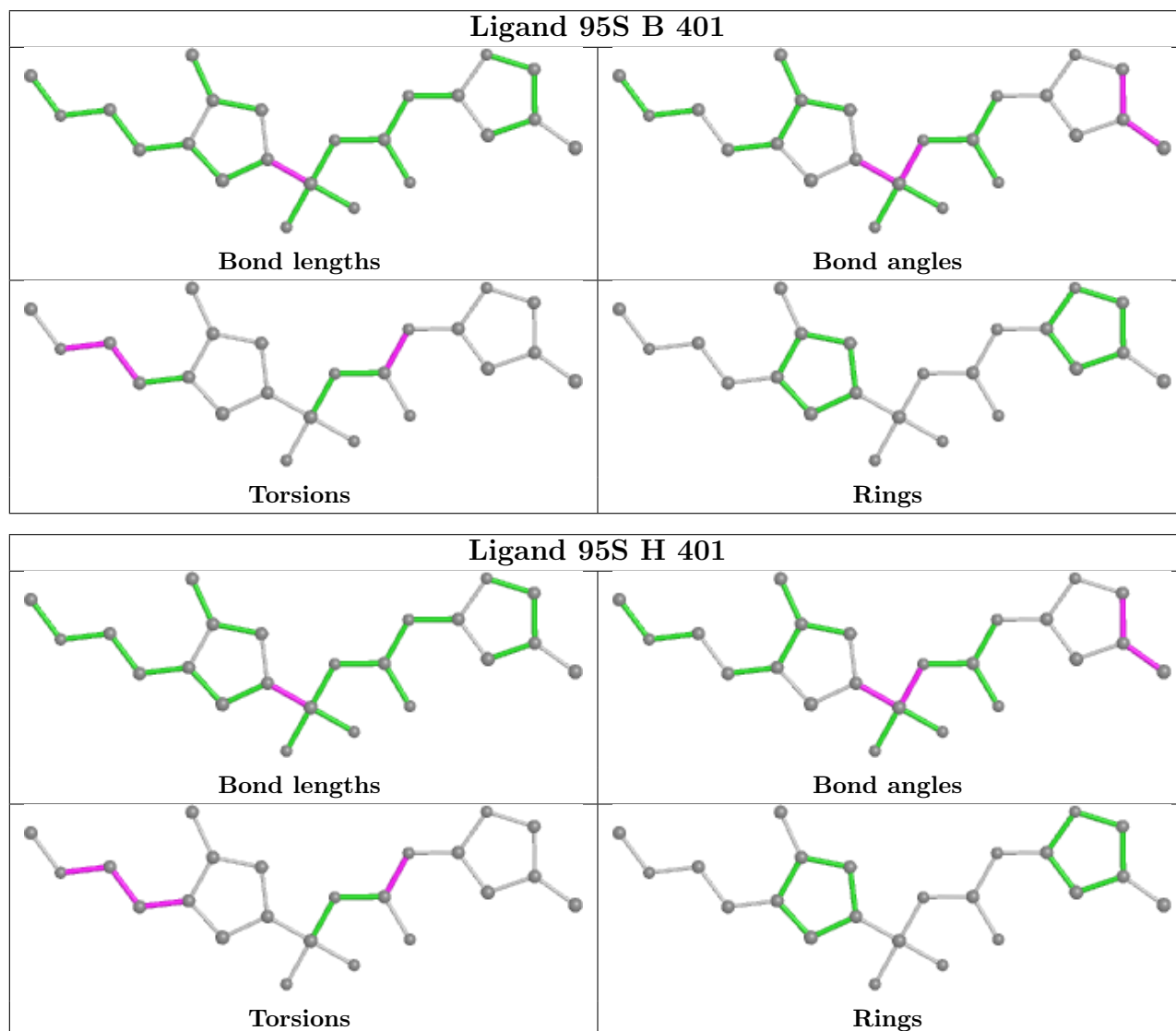
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	95S	1	0
2	D	401	95S	1	0
2	E	401	95S	1	0
2	C	401	95S	1	0
2	F	401	95S	2	0

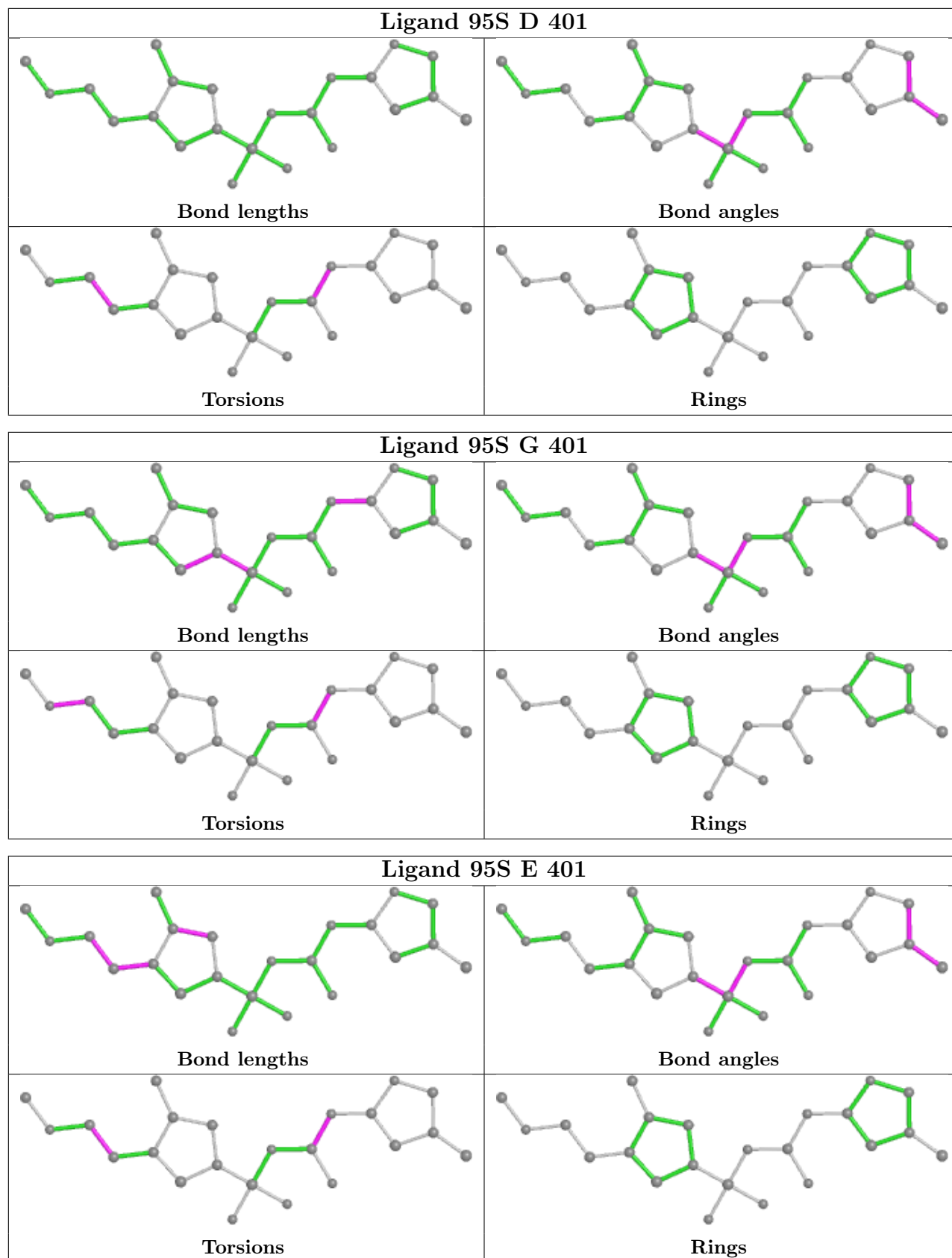
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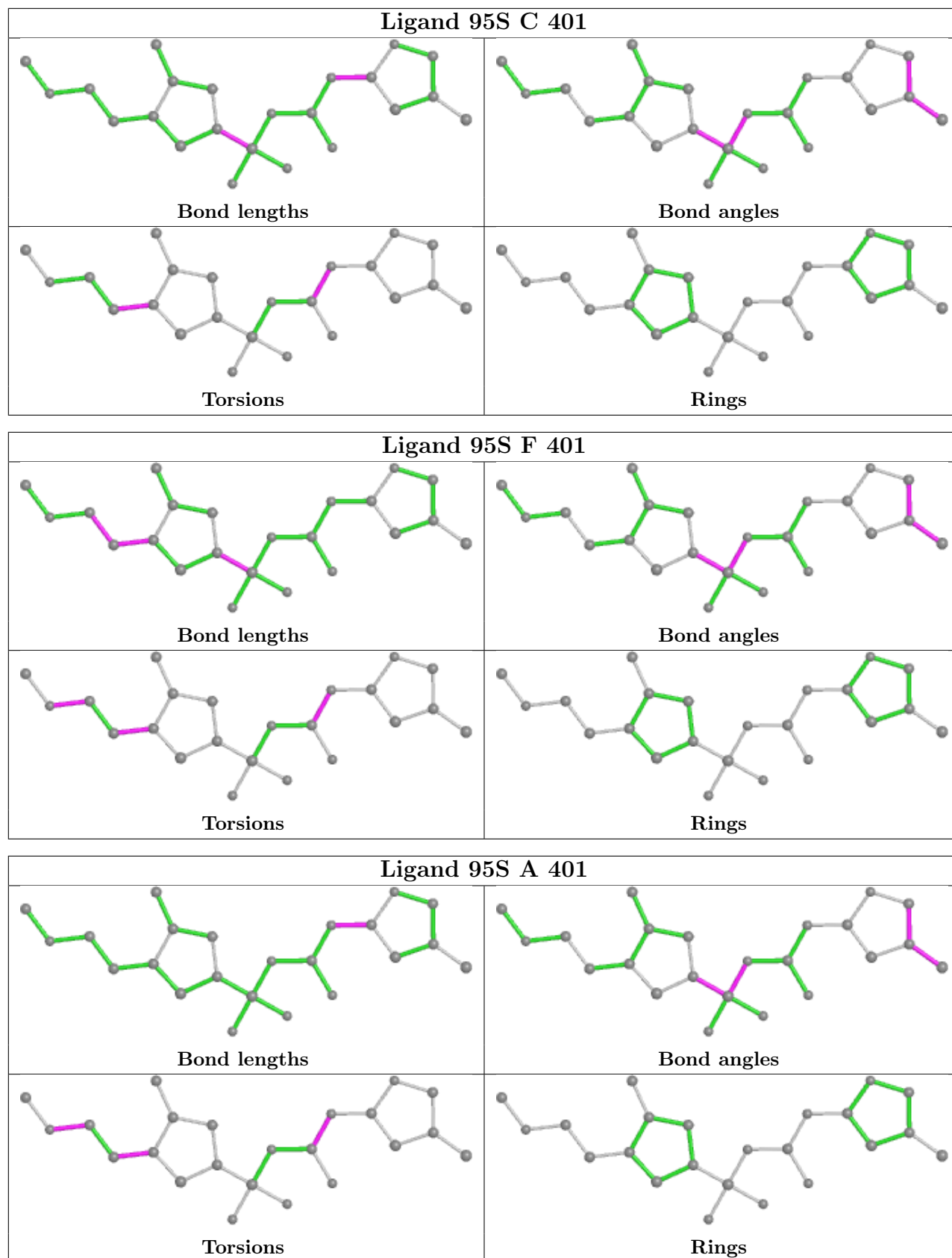
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	95S	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	319/338 (94%)	-0.12	8 (2%) 57 55	26, 39, 70, 120	0
1	B	317/338 (93%)	-0.05	9 (2%) 53 51	22, 41, 75, 139	0
1	C	317/338 (93%)	-0.25	3 (0%) 84 83	23, 35, 61, 104	0
1	D	317/338 (93%)	-0.33	2 (0%) 89 88	19, 34, 63, 96	0
1	E	319/338 (94%)	0.00	9 (2%) 53 51	27, 46, 86, 126	0
1	F	317/338 (93%)	-0.01	14 (4%) 34 32	27, 45, 85, 116	0
1	G	317/338 (93%)	-0.19	5 (1%) 72 70	23, 37, 64, 100	0
1	H	317/338 (93%)	-0.23	4 (1%) 77 75	22, 37, 68, 88	0
All	All	2540/2704 (93%)	-0.15	54 (2%) 63 61	19, 39, 73, 139	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	VAL	11.3
1	F	268	LYS	6.6
1	A	71	LYS	4.9
1	C	236	ASN	4.7
1	E	235	ASP	4.6
1	E	70	VAL	4.4
1	G	236	ASN	4.3
1	A	234	PRO	3.7
1	B	236	ASN	3.4
1	G	126	ILE	3.4
1	F	236	ASN	3.3
1	B	145	ASP	3.3
1	A	107	PRO	3.2
1	B	269	LYS	3.2
1	E	268	LYS	3.2
1	E	269	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	F	150	LYS	3.0
1	F	9	ASP	3.0
1	H	236	ASN	3.0
1	E	266	ALA	2.9
1	F	144	THR	2.8
1	E	236	ASN	2.8
1	G	235	ASP	2.8
1	A	38	CYS	2.7
1	F	145	ASP	2.7
1	B	37	LEU	2.6
1	D	236	ASN	2.6
1	A	25	ARG	2.6
1	F	269	LYS	2.6
1	B	56	LEU	2.5
1	E	238	ALA	2.5
1	A	268	LYS	2.4
1	F	38	CYS	2.4
1	C	126	ILE	2.4
1	B	235	ASP	2.4
1	F	109	LYS	2.3
1	H	268	LYS	2.3
1	H	271	PRO	2.3
1	F	271	PRO	2.2
1	F	234	PRO	2.2
1	C	237	SER	2.2
1	H	25	ARG	2.2
1	G	123	SER	2.2
1	B	126	ILE	2.2
1	E	234	PRO	2.2
1	B	144	THR	2.1
1	B	335	SER	2.1
1	A	146	GLU	2.1
1	F	25	ARG	2.1
1	D	237	SER	2.0
1	F	107	PRO	2.0
1	G	335	SER	2.0
1	F	138	ILE	2.0
1	E	237	SER	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

There are no monosaccharides in this entry.

6.4 Ligands

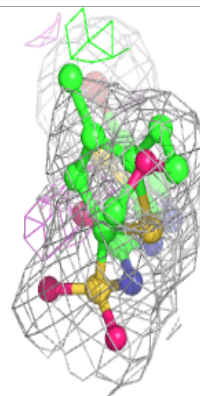
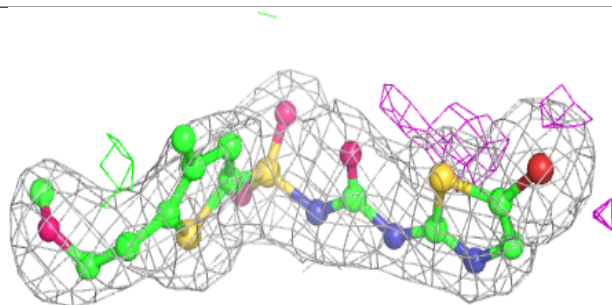
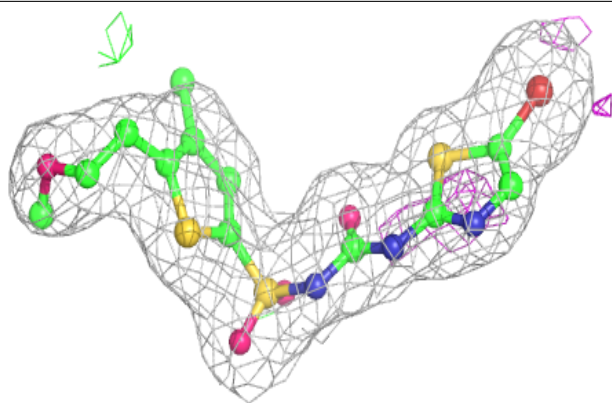
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	95S	H	401	23/23	0.97	0.09	36,45,55,59	1
2	95S	G	401	23/23	0.97	0.10	31,42,52,64	1
2	95S	A	401	23/23	0.97	0.10	19,42,57,65	1
2	95S	B	401	23/23	0.98	0.09	28,41,52,53	1
2	95S	E	401	23/23	0.98	0.09	23,37,50,59	1
2	95S	C	401	23/23	0.98	0.09	30,41,52,61	1
2	95S	F	401	23/23	0.98	0.08	36,48,58,61	1
2	95S	D	401	23/23	0.98	0.08	21,36,50,54	1

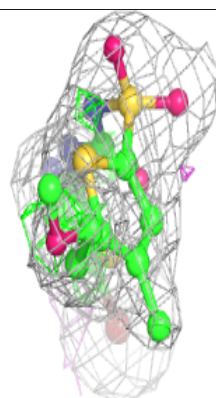
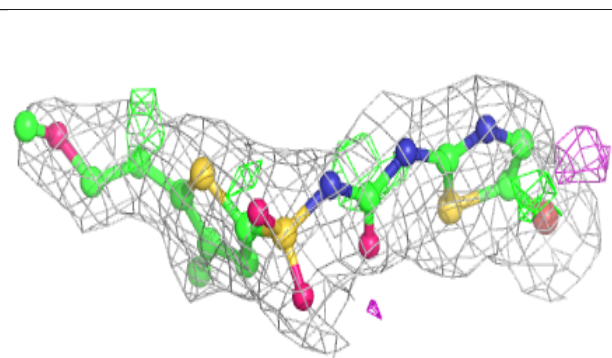
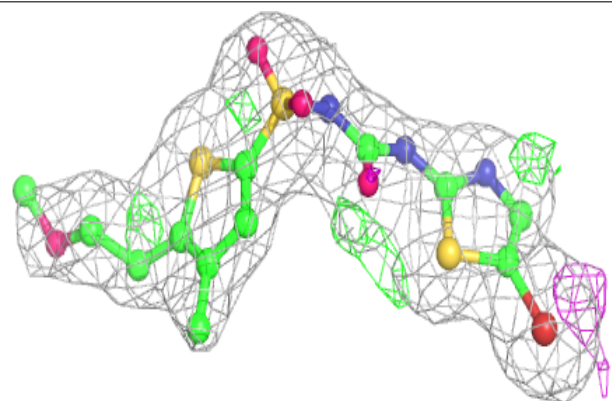
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 95S H 401:

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

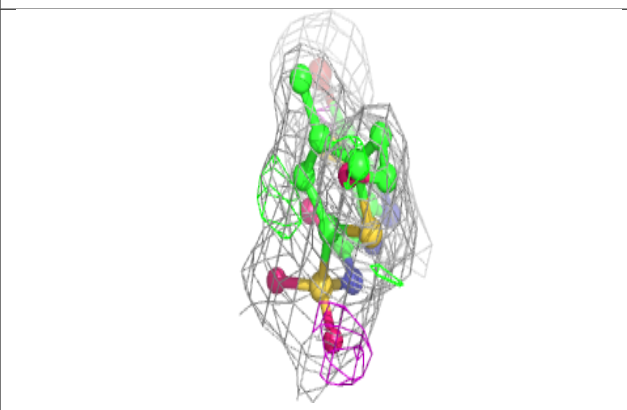
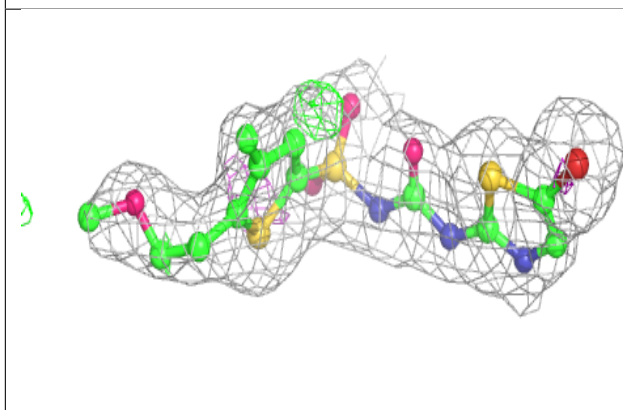
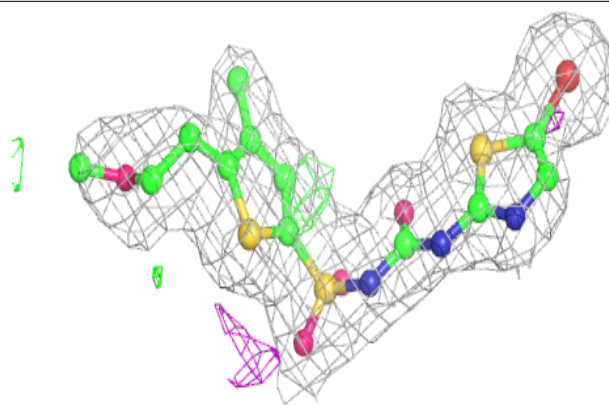
**Electron density around 95S G 401:**

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

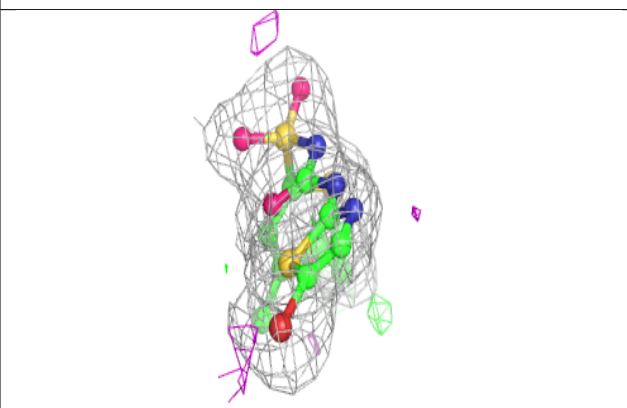
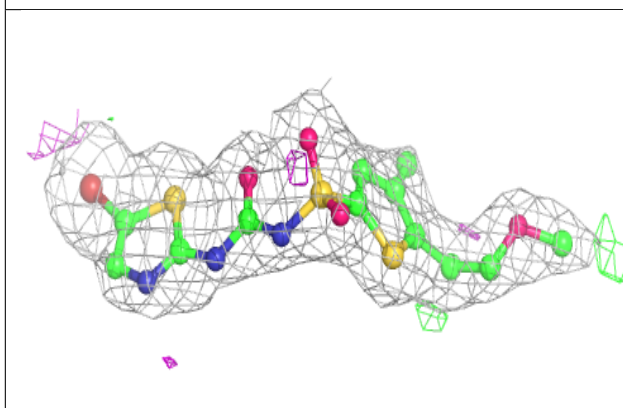
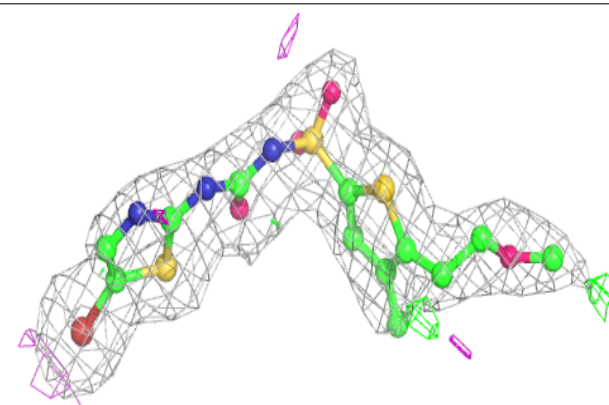


Electron density around 95S A 401:

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

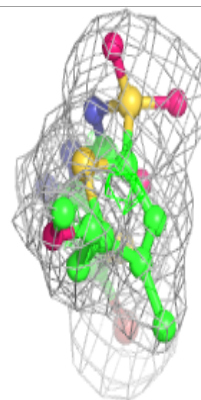
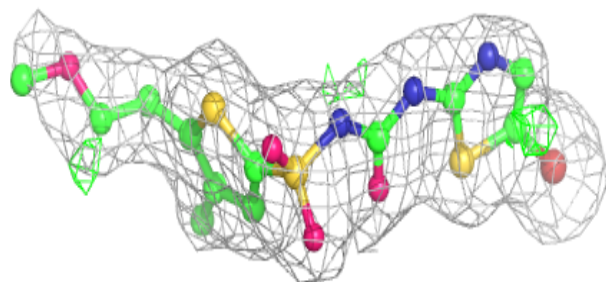
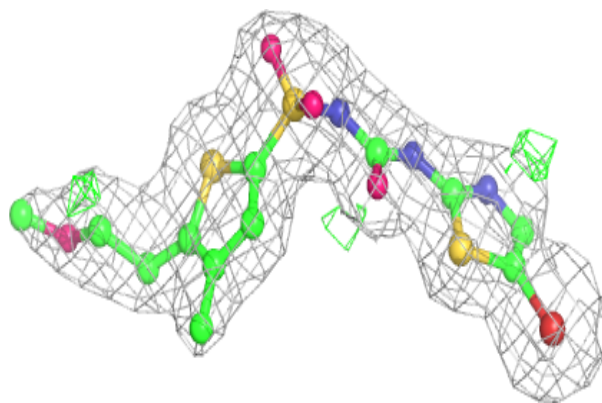
**Electron density around 95S B 401:**

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

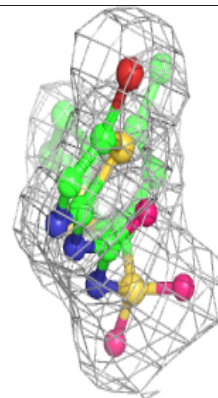
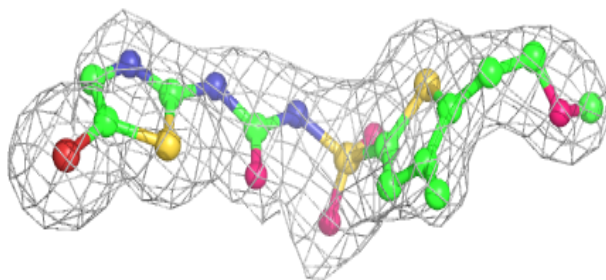
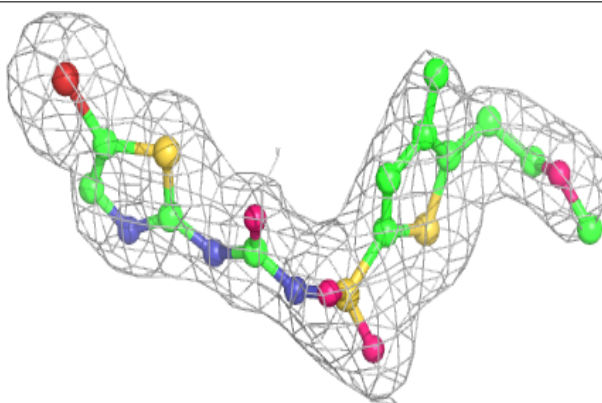


Electron density around 95S E 401:

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

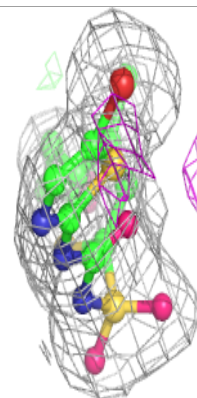
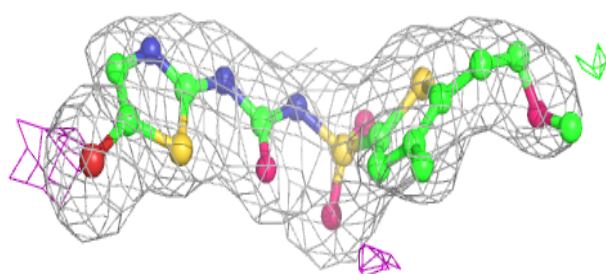
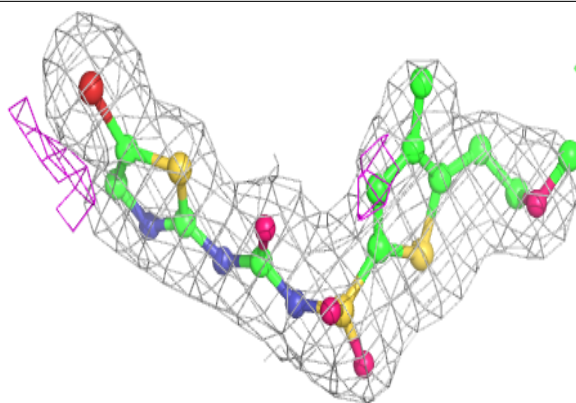
**Electron density around 95S C 401:**

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

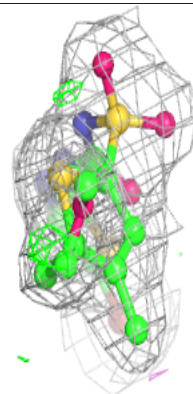
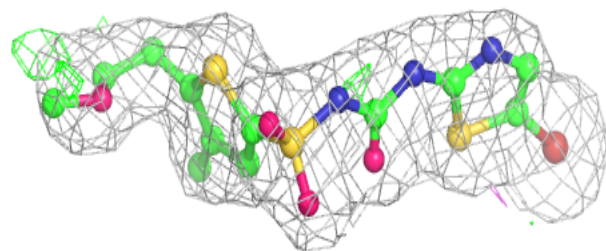
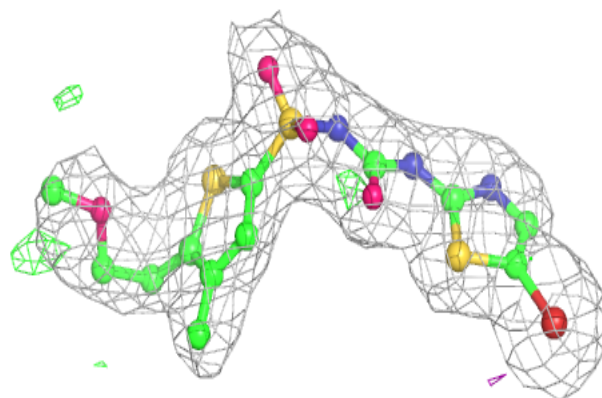


Electron density around 95S F 401:

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

**Electron density around 95S D 401:**

$2mF_o-DF_c$ (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.