



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2022 – 06:05 PM JST

PDB ID : 7EZP
Title : Indole-2-carboxylic acid derivatives as allosteric inhibitors of fructose-1,6-bisphosphatase
Authors : Wang, X.Y.; Zhou, J.; Xu, B.L.
Deposited on : 2021-06-01
Resolution : 2.80 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)
Xtriage (Phenix) : 1.13
EDS : 2.28.1
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) : 2.28.1

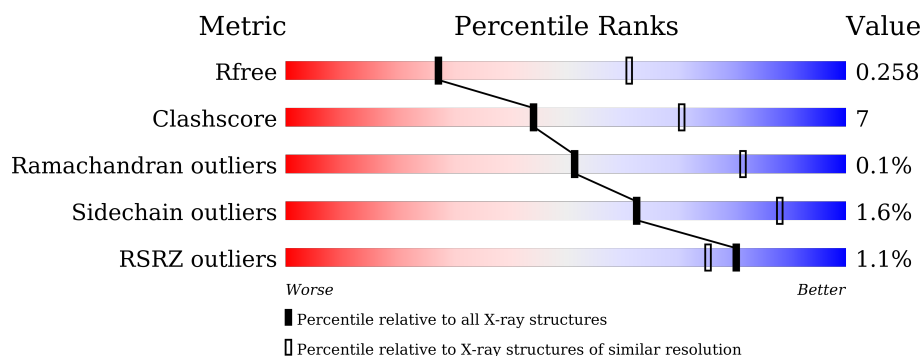
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	<div> <div>2%</div> <div>78% 16% • 5%</div> </div>
1	B	338	<div> <div>2%</div> <div>78% 17% 5%</div> </div>
1	C	338	<div> <div>82% 12% • 6%</div> </div>
1	D	338	<div> <div>2%</div> <div>81% 13% • 5%</div> </div>

2 Entry composition [i](#)

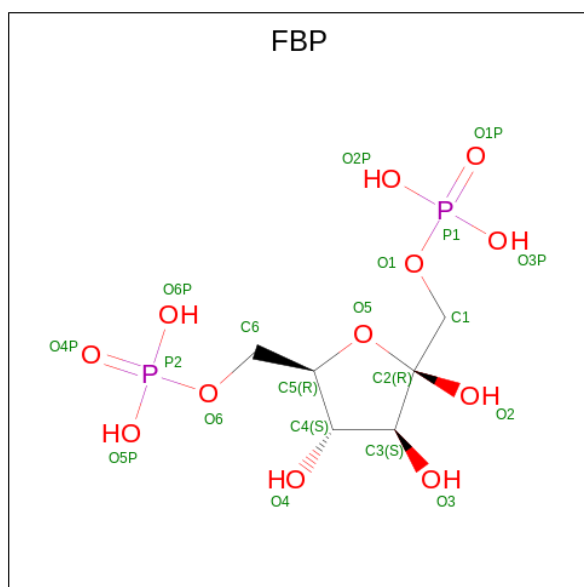
There are 4 unique types of molecules in this entry. The entry contains 10090 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
1	A	320	Total	C	N	O	S	0	0	0
			2445	1555	411	462	17			
1	B	320	Total	C	N	O	S	0	0	1
			2441	1552	411	461	17			
1	C	319	Total	C	N	O	S	0	0	1
			2434	1548	410	459	17			
1	D	320	Total	C	N	O	S	0	0	0
			2445	1555	411	462	17			

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (three-letter code: FBP) (formula: $C_6H_{14}O_{12}P_2$) (labeled as "Ligand of Interest" by depositor).



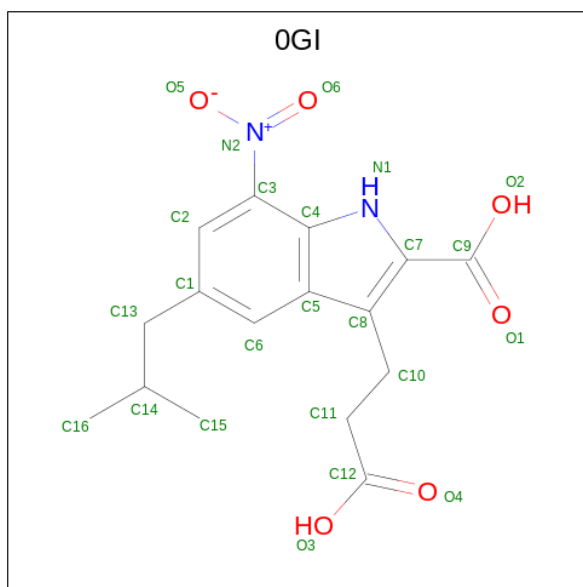
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	O	P	0	0
			20	6	12	2		
2	D	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 3 is 3-(3-hydroxy-3-oxopropyl)-5-(2-methylpropyl)-7-nitro-1H-indole-2-carboxylic acid (three-letter code: 0GI) (formula: C₁₆H₁₈N₂O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			24	16	2	6		
3	B	1	Total	C	N	O	0	0
			24	16	2	6		
3	C	1	Total	C	N	O	0	0
			24	16	2	6		
3	D	1	Total	C	N	O	0	0
			24	16	2	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	29	Total	O	0	0
			29	29		
4	B	32	Total	O	0	0
			32	32		

Continued on next page...

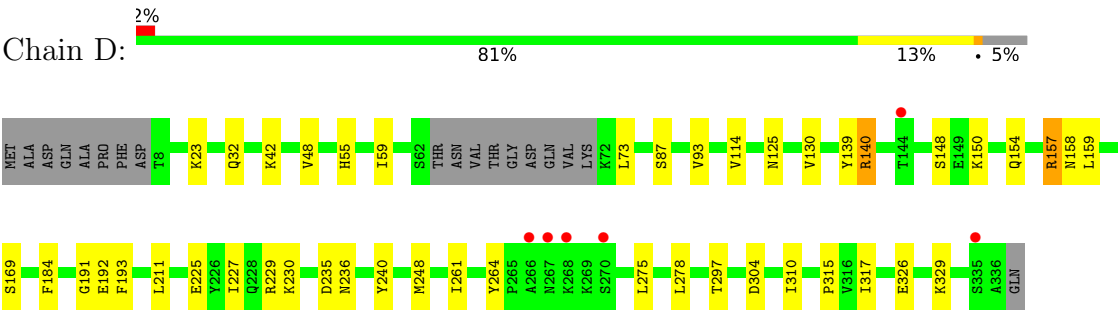
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	47	Total	O	0	0
			47	47		
4	D	41	Total	O	0	0
			41	41		

- 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	67.10Å 143.00Å 73.60Å 90.00° 107.74° 90.00°	Depositor
Resolution (Å)	70.20 – 2.80 70.10 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (70.20-2.80) 99.7 (70.10-2.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.210 , 0.256 0.211 , 0.258	Depositor DCC
R_{free} test set	1644 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.5	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	10090	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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: 0GI, FBP

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	0.51	0/2488	0.54	0/3361
1	B	0.52	0/2484	0.54	0/3356
1	C	0.54	0/2477	0.65	0/3346
1	D	0.52	0/2488	0.57	0/3361
All	All	0.52	0/9937	0.58	0/13424

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2445	0	2486	46	0
1	B	2441	0	2481	40	0
1	C	2434	0	2474	24	1
1	D	2445	0	2486	33	1
2	A	20	0	10	0	0
2	B	20	0	10	0	0
2	C	20	0	10	0	0
2	D	20	0	10	0	0
3	A	24	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	24	0	0	0	0
3	C	24	0	0	0	0
3	D	24	0	0	0	0
4	A	29	0	0	2	0
4	B	32	0	0	1	0
4	C	47	0	0	1	0
4	D	41	0	0	0	0
All	All	10090	0	9967	131	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 7.

All (131) 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:B:15:ARG:HG2	1:D:32:GLN:HE22	1.33	0.93
1:A:14:THR:HG23	1:A:35:ASN:HD21	1.35	0.90
1:A:225:GLU:OE2	1:A:334:HIS:HE1	1.54	0.90
1:D:229:ARG:NH2	1:D:326:GLU:OE2	2.05	0.90
1:A:299:LYS:HE2	1:A:299:LYS:HA	1.53	0.89
1:D:157:ARG:NH1	1:D:304:ASP:OD1	2.06	0.88
1:A:235:ASP:OD1	4:A:501:HOH:O	1.92	0.88
1:A:320:SER:O	1:A:324:VAL:HG23	1.82	0.79
1:A:235:ASP:OD2	1:A:235:ASP:O	2.00	0.78
1:B:328:LEU:O	1:B:332:GLU:HG3	1.86	0.75
1:D:297:THR:HB	1:D:315:PRO:HG2	1.69	0.75
1:A:235:ASP:O	1:A:235:ASP:CG	2.29	0.70
1:B:144:THR:HG22	1:B:144:THR:O	1.91	0.70
1:A:113:TYR:CE1	1:A:140:ARG:HG3	2.28	0.69
1:D:230:LYS:NZ	1:D:240:TYR:CZ	2.58	0.69
1:C:277:LEU:HA	1:C:281:CYS:HB2	1.73	0.69
1:D:211:LEU:HD11	1:D:227:ILE:HD11	1.74	0.68
1:A:225:GLU:OE2	1:A:334:HIS:CE1	2.44	0.67
1:C:112:LYS:N	1:C:112:LYS:HD2	2.10	0.66
1:A:299:LYS:HA	1:A:299:LYS:CE	2.22	0.65
1:D:229:ARG:NH2	1:D:326:GLU:CD	2.48	0.65
1:A:221:PRO:HB2	1:A:334:HIS:HD2	1.62	0.64
1:B:42:LYS:NZ	1:D:192:GLU:OE1	2.26	0.64
1:B:14:THR:O	1:B:18:MET:HG3	1.97	0.64
1:D:229:ARG:HH21	1:D:326:GLU:CD	2.00	0.64
1:B:220:ASP:HB3	1:B:223:VAL:HG23	1.77	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:GLU:OE1	1:B:334:HIS:NE2	2.29	0.62
1:A:198:LYS:HD2	1:A:199:ASP:H	1.64	0.62
1:A:224:THR:O	1:A:228:GLN:HG3	2.00	0.61
1:B:15:ARG:HH12	1:D:87:SER:HB3	1.65	0.61
1:C:146:GLU:HG3	1:C:147:PRO:HD2	1.82	0.60
1:B:107:PRO:HA	1:B:110:ARG:HG3	1.84	0.60
1:A:277:LEU:HA	1:A:281:CYS:HB2	1.84	0.59
1:C:72:LYS:N	4:C:501:HOH:O	2.36	0.59
1:A:258:TYR:OH	1:B:127:ASP:HB2	2.02	0.59
1:A:191:GLY:HA3	1:C:191:GLY:HA3	1.86	0.58
1:C:133:GLY:HA3	1:C:249:VAL:HG21	1.86	0.58
1:A:183:CYS:HB2	1:A:197:ASP:HB2	1.85	0.57
1:D:48:VAL:HA	1:D:73:LEU:HD21	1.86	0.57
1:C:218:ASP:HB3	1:C:268:LYS:HB2	1.87	0.55
1:D:148:SER:OG	1:D:150:LYS:HG2	2.07	0.55
1:B:251:ASP:OD1	1:B:254:ARG:NH1	2.38	0.54
1:D:154:GLN:OE1	1:D:158:ASN:HB3	2.08	0.54
1:C:302:VAL:HG21	1:C:316:VAL:HG23	1.90	0.53
1:B:123:SER:HA	1:B:126:ILE:HB	1.90	0.52
1:C:262:PHE:HB3	1:C:318:LEU:HD11	1.92	0.51
1:D:225:GLU:O	1:D:229:ARG:HG3	2.10	0.51
1:A:57:TYR:CZ	1:B:196:VAL:HG12	2.46	0.51
1:A:182:ASN:OD1	1:A:198:LYS:HD2	2.10	0.51
1:A:35:ASN:ND2	1:C:14:THR:HG21	2.26	0.51
1:B:82:MET:HE1	1:B:103:ILE:HG13	1.92	0.51
1:B:144:THR:O	1:B:144:THR:CG2	2.58	0.51
1:B:222:ALA:HB1	1:B:330:VAL:HG12	1.92	0.50
1:D:229:ARG:NH2	1:D:326:GLU:CG	2.74	0.50
1:A:278:LEU:HD12	1:A:310:ILE:HA	1.93	0.50
1:A:215:TYR:HB2	1:A:219:PHE:CE1	2.46	0.50
1:C:104:ILE:HG23	1:C:147:PRO:HB2	1.93	0.50
1:A:188:PRO:HD2	1:B:51:ALA:HA	1.94	0.49
1:B:25:ARG:HH11	1:B:25:ARG:HG3	1.78	0.49
1:B:216:ALA:HA	1:B:219:PHE:CD1	2.47	0.49
1:B:15:ARG:CG	1:D:32:GLN:HE22	2.16	0.49
1:A:45:SER:O	1:A:49:ARG:HD3	2.13	0.48
1:A:328:LEU:O	1:A:332:GLU:HG3	2.13	0.48
1:D:261:ILE:HD11	1:D:317:ILE:CG2	2.43	0.48
1:A:299:LYS:HE2	1:A:299:LYS:CA	2.28	0.48
1:D:93:VAL:HB	1:D:114:VAL:HG22	1.94	0.48
1:A:13:LEU:O	1:A:17:VAL:HG23	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:111:GLY:C	1:C:112:LYS:HD2	2.34	0.48
1:C:262:PHE:HB3	1:C:318:LEU:CD1	2.44	0.48
1:A:29:GLU:OE1	1:A:112:LYS:HG2	2.14	0.47
1:A:106:GLU:OE1	1:A:109:LYS:HD2	2.13	0.47
1:A:127:ASP:OD1	4:A:502:HOH:O	2.20	0.47
1:A:221:PRO:HB2	1:A:334:HIS:CD2	2.46	0.47
1:B:25:ARG:HG3	1:B:25:ARG:NH1	2.30	0.47
1:D:230:LYS:NZ	1:D:240:TYR:CE2	2.75	0.47
1:D:229:ARG:NH2	1:D:326:GLU:HG2	2.30	0.47
1:B:55:HIS:HA	1:B:59:ILE:HG22	1.97	0.46
1:A:228:GLN:HE21	1:A:228:GLN:HB3	1.54	0.46
1:A:15:ARG:HG3	1:C:32:GLN:OE1	2.15	0.46
1:A:17:VAL:HG12	1:A:31:THR:HG23	1.98	0.45
1:A:299:LYS:O	1:A:299:LYS:HD3	2.16	0.45
1:A:172:MET:HE2	1:A:185:MET:HE2	1.97	0.45
1:B:277:LEU:HA	1:B:281:CYS:HB2	1.98	0.45
1:B:159:LEU:HD23	1:B:159:LEU:HA	1.84	0.45
1:C:233:PRO:HA	1:C:234:PRO:HD3	1.86	0.45
1:A:166:LEU:O	1:A:171:THR:HA	2.17	0.45
1:B:143:SER:OG	1:B:151:ASP:OD1	2.35	0.45
1:B:166:LEU:O	1:B:171:THR:HA	2.16	0.44
1:D:278:LEU:HD12	1:D:310:ILE:HA	1.98	0.44
1:D:154:GLN:CG	1:D:158:ASN:HD22	2.31	0.44
1:D:55:HIS:HA	1:D:59:ILE:HG22	1.99	0.44
1:D:139:TYR:CE2	1:D:159:LEU:HG	2.53	0.44
1:D:154:GLN:HG2	1:D:158:ASN:HB2	1.99	0.44
1:D:125:ASN:O	1:D:130:VAL:HG12	2.18	0.44
1:A:329:LYS:HD2	1:A:329:LYS:N	2.31	0.43
1:C:278:LEU:HD22	1:C:310:ILE:HA	2.00	0.43
1:A:186:LEU:HB2	1:A:193:PHE:CE1	2.54	0.43
1:B:10:VAL:O	1:B:10:VAL:HG13	2.19	0.43
1:C:103:ILE:N	1:C:103:ILE:HD12	2.34	0.43
1:A:48:VAL:HA	1:A:73:LEU:HD21	2.01	0.43
1:A:14:THR:O	1:A:18:MET:HG3	2.20	0.42
1:B:198:LYS:HE3	1:B:198:LYS:HB3	1.84	0.42
1:B:203:LYS:HG3	4:B:516:HOH:O	2.19	0.42
1:D:248:MET:CE	1:D:275:LEU:HD13	2.50	0.42
1:B:157:ARG:HD3	1:B:304:ASP:OD1	2.20	0.42
1:B:290:LYS:HD3	1:B:290:LYS:HA	1.76	0.42
1:A:14:THR:CG2	1:A:35:ASN:HD21	2.19	0.42
1:B:90:ALA:HA	1:B:111:GLY:HA3	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:166:LEU:O	1:C:171:THR:HA	2.19	0.42
1:C:300:GLU:HG3	1:C:305:VAL:HG22	2.00	0.42
1:B:106:GLU:HA	1:B:107:PRO:HD3	1.89	0.42
1:C:49:ARG:HD3	1:D:169:SER:O	2.20	0.42
1:D:184:PHE:HB3	1:D:193:PHE:HB3	2.02	0.42
1:C:178:ASP:HA	1:C:290:LYS:HE3	2.01	0.42
1:B:219:PHE:HD2	1:B:219:PHE:HA	1.79	0.41
1:B:278:LEU:HD12	1:B:310:ILE:HA	2.02	0.41
1:C:131:SER:HB3	1:C:166:LEU:HD11	2.02	0.41
1:C:228:GLN:NE2	1:C:232:PHE:CD1	2.85	0.41
1:B:17:VAL:HG13	1:B:31:THR:HG23	2.02	0.41
1:D:23:LYS:HB3	1:D:23:LYS:HE3	1.78	0.41
1:D:154:GLN:HE21	1:D:154:GLN:HB2	1.71	0.41
1:B:191:GLY:HA3	1:D:191:GLY:HA3	2.03	0.41
1:B:120:LEU:HD12	1:B:120:LEU:HA	1.82	0.41
1:A:59:ILE:HD12	1:A:59:ILE:HA	1.89	0.41
1:A:73:LEU:HD12	1:A:73:LEU:HA	1.96	0.41
1:C:216:ALA:HA	1:C:219:PHE:CD2	2.56	0.41
1:B:183:CYS:HB2	1:B:197:ASP:HB2	2.02	0.41
1:D:42:LYS:HB2	1:D:42:LYS:HE3	1.95	0.40
1:A:120:LEU:HD12	1:A:120:LEU:HA	1.87	0.40
1:A:219:PHE:HD1	1:A:219:PHE:HA	1.76	0.40
1:B:184:PHE:HB3	1:B:193:PHE:HB3	2.04	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:C:24:ALA:O	1:D:140:ARG:NH1[1_554]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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%)	301 (95%)	15 (5%)	0	100	100
1	B	316/338 (94%)	297 (94%)	19 (6%)	0	100	100
1	C	315/338 (93%)	293 (93%)	21 (7%)	1 (0%)	41	72
1	D	316/338 (94%)	299 (95%)	17 (5%)	0	100	100
All	All	1263/1352 (93%)	1190 (94%)	72 (6%)	1 (0%)	51	81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	234	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	266/281 (95%)	261 (98%)	5 (2%)	57	85
1	B	266/281 (95%)	265 (100%)	1 (0%)	91	97
1	C	265/281 (94%)	260 (98%)	5 (2%)	57	85
1	D	266/281 (95%)	260 (98%)	6 (2%)	50	82
All	All	1063/1124 (95%)	1046 (98%)	17 (2%)	62	88

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	LYS
1	A	197	ASP
1	A	220	ASP
1	A	232	PHE
1	A	265	PRO
1	B	145	ASP
1	C	9	ASP
1	C	130	VAL
1	C	221	PRO
1	C	268	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	269	LYS
1	D	140	ARG
1	D	157	ARG
1	D	235	ASP
1	D	236	ASN
1	D	264	TYR
1	D	329	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	228	GLN
1	A	236	ASN
1	A	334	HIS
1	B	125	ASN
1	D	32	GLN
1	D	55	HIS
1	D	158	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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
3	0GI	A	402	-	14,25,25	2.75	2 (14%)	16,36,36	1.09	1 (6%)
2	FBP	A	401	-	18,20,20	3.28	5 (27%)	23,32,32	0.70	0
3	0GI	B	402	-	14,25,25	2.91	2 (14%)	16,36,36	0.77	0
3	0GI	D	402	-	14,25,25	2.91	2 (14%)	16,36,36	0.87	0
2	FBP	D	401	-	18,20,20	3.27	5 (27%)	23,32,32	0.75	0
3	0GI	C	402	-	14,25,25	2.85	2 (14%)	16,36,36	1.05	0
2	FBP	C	401	-	18,20,20	3.36	7 (38%)	23,32,32	0.83	1 (4%)
2	FBP	B	401	-	18,20,20	3.39	5 (27%)	23,32,32	0.77	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	0GI	A	402	-	-	1/9/17/17	0/2/2/2
2	FBP	A	401	-	-	1/13/32/32	0/1/1/1
3	0GI	B	402	-	-	1/9/17/17	0/2/2/2
3	0GI	D	402	-	-	0/9/17/17	0/2/2/2
2	FBP	D	401	-	-	9/13/32/32	0/1/1/1
3	0GI	C	402	-	-	3/9/17/17	0/2/2/2
2	FBP	C	401	-	-	1/13/32/32	0/1/1/1
2	FBP	B	401	-	-	2/13/32/32	0/1/1/1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	402	0GI	O6-N2	10.07	1.39	1.22
3	B	402	0GI	O6-N2	10.02	1.39	1.22
3	C	402	0GI	O6-N2	10.00	1.39	1.22
2	B	401	FBP	O5-C2	-9.75	1.28	1.43
3	A	402	0GI	O6-N2	9.59	1.39	1.22
2	C	401	FBP	O5-C2	-9.35	1.28	1.43
2	D	401	FBP	O5-C2	-9.21	1.29	1.43
2	A	401	FBP	O5-C2	-9.12	1.29	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	FBP	C4-C3	-6.61	1.25	1.52
2	C	401	FBP	C4-C3	-6.46	1.25	1.52
2	A	401	FBP	C4-C3	-6.38	1.26	1.52
2	D	401	FBP	C4-C3	-6.34	1.26	1.52
2	A	401	FBP	O5-C5	4.72	1.54	1.43
2	A	401	FBP	C6-C5	-4.59	1.37	1.51
2	D	401	FBP	C6-C5	-4.53	1.37	1.51
2	B	401	FBP	C6-C5	-4.52	1.37	1.51
2	C	401	FBP	O5-C5	4.50	1.53	1.43
2	C	401	FBP	C6-C5	-4.49	1.37	1.51
2	D	401	FBP	O5-C5	4.25	1.53	1.43
2	B	401	FBP	O5-C5	4.23	1.53	1.43
3	B	402	0GI	C7-N1	-3.53	1.30	1.37
3	D	402	0GI	C7-N1	-3.32	1.31	1.37
3	A	402	0GI	C7-N1	-3.17	1.31	1.37
3	C	402	0GI	C7-N1	-3.12	1.31	1.37
2	B	401	FBP	O4-C4	3.05	1.50	1.43
2	C	401	FBP	O4-C4	3.00	1.50	1.43
2	D	401	FBP	O4-C4	2.92	1.49	1.43
2	A	401	FBP	O4-C4	2.87	1.49	1.43
2	C	401	FBP	C4-C5	2.22	1.58	1.53
2	C	401	FBP	O2-C2	-2.17	1.37	1.40

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	0GI	C10-C8-C7	-2.85	124.38	127.48
2	C	401	FBP	O1-P1-O1P	2.14	112.46	106.47

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	FBP	C1-O1-P1-O3P
2	D	401	FBP	C1-O1-P1-O1P
2	D	401	FBP	C1-O1-P1-O2P
2	D	401	FBP	C6-O6-P2-O4P
2	D	401	FBP	C6-O6-P2-O6P
3	A	402	0GI	C8-C10-C11-C12
3	B	402	0GI	C8-C10-C11-C12
3	C	402	0GI	C8-C10-C11-C12
2	B	401	FBP	C1-O1-P1-O1P

Continued on next page...

Continued from previous page...

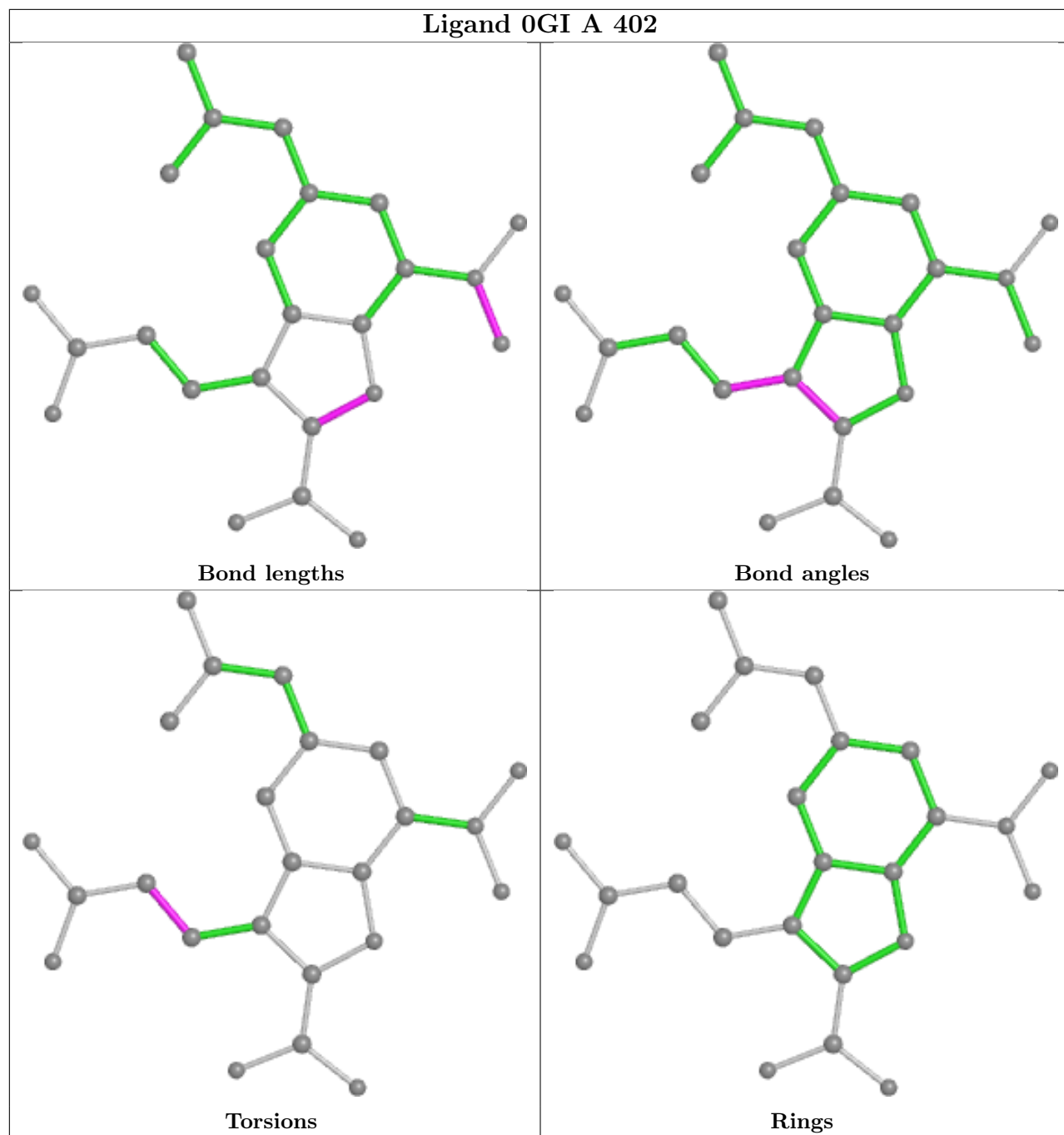
Mol	Chain	Res	Type	Atoms
3	C	402	0GI	C1-C13-C14-C15
3	C	402	0GI	C1-C13-C14-C16
2	D	401	FBP	O1-C1-C2-O5
2	D	401	FBP	O1-C1-C2-O2
2	B	401	FBP	O1-C1-C2-C3
2	C	401	FBP	O1-C1-C2-C3
2	D	401	FBP	O1-C1-C2-C3
2	D	401	FBP	C1-O1-P1-O3P
2	D	401	FBP	C6-O6-P2-O5P

There are no ring outliers.

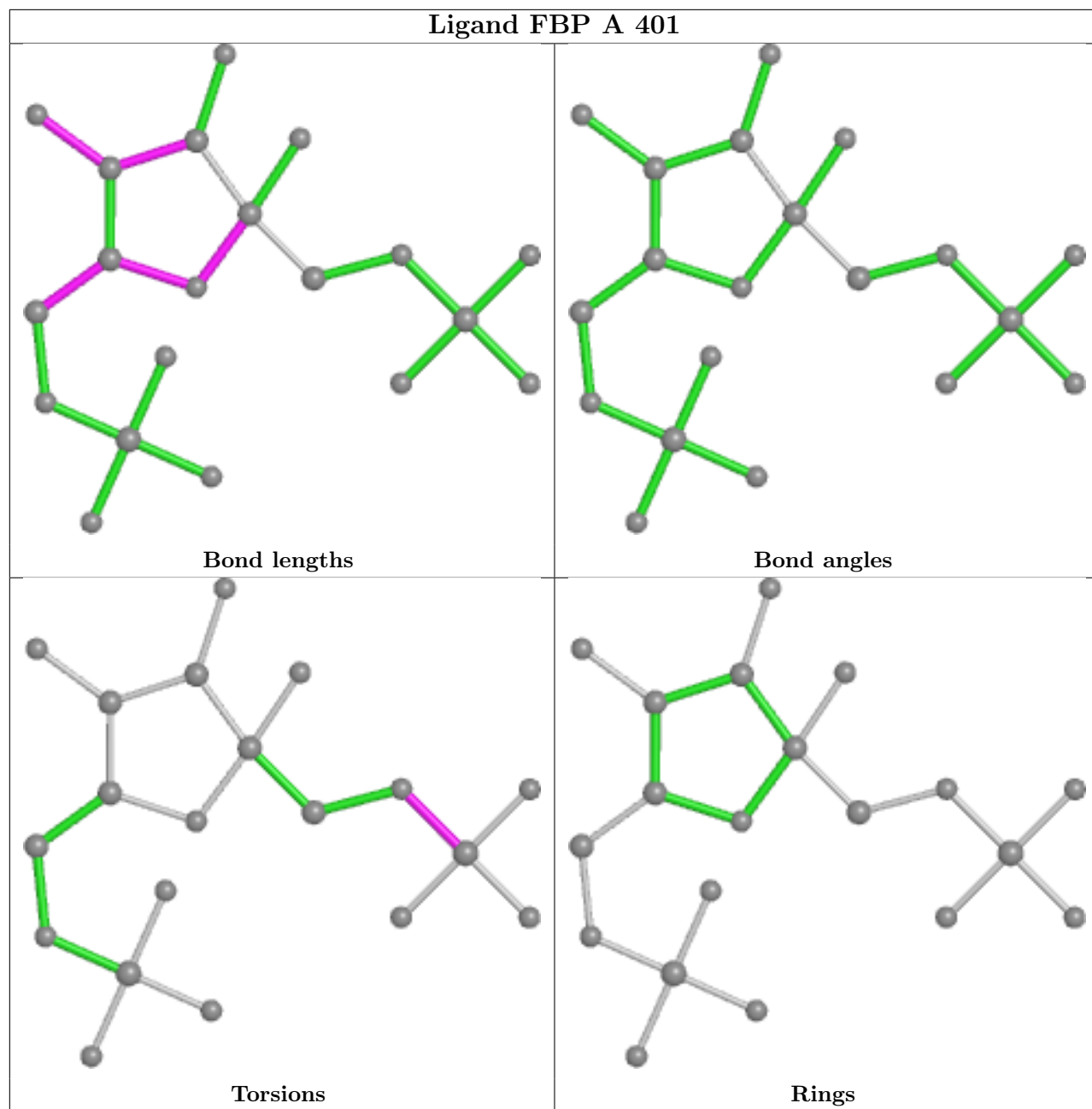
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

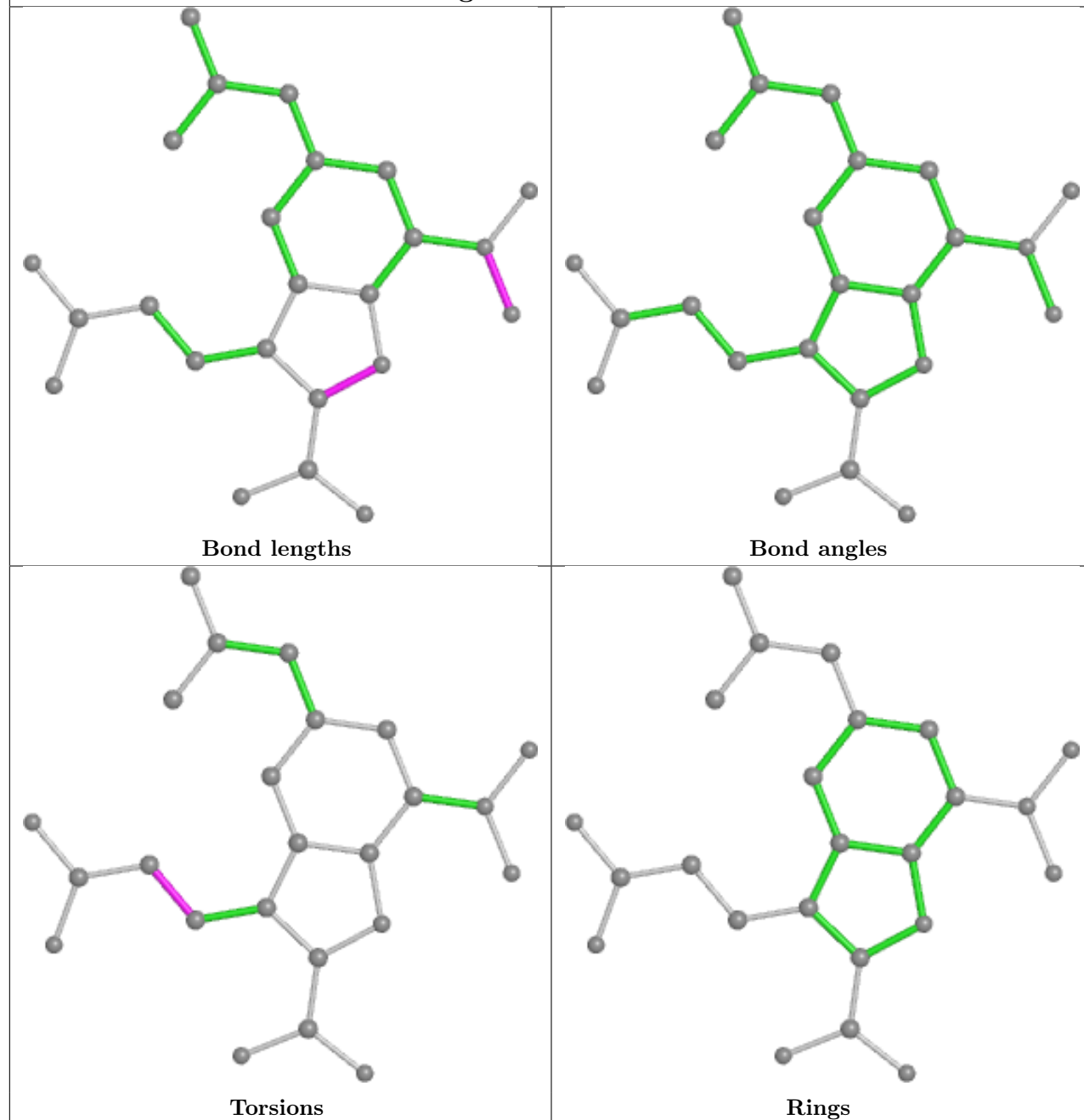
Ligand 0GI A 402



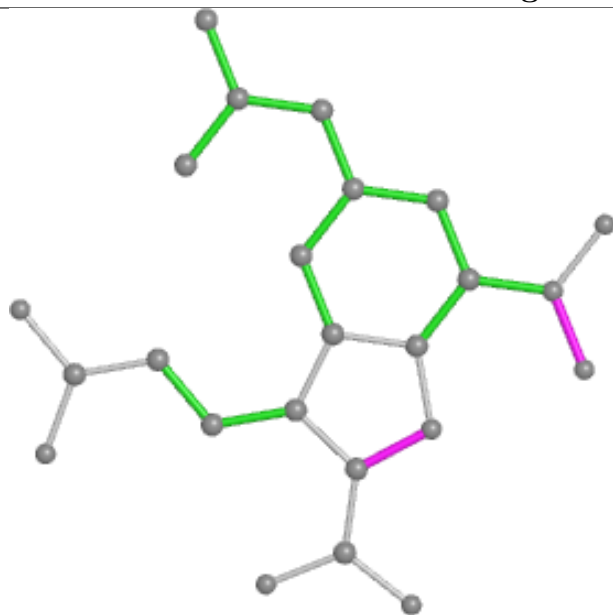
Ligand FBP A 401



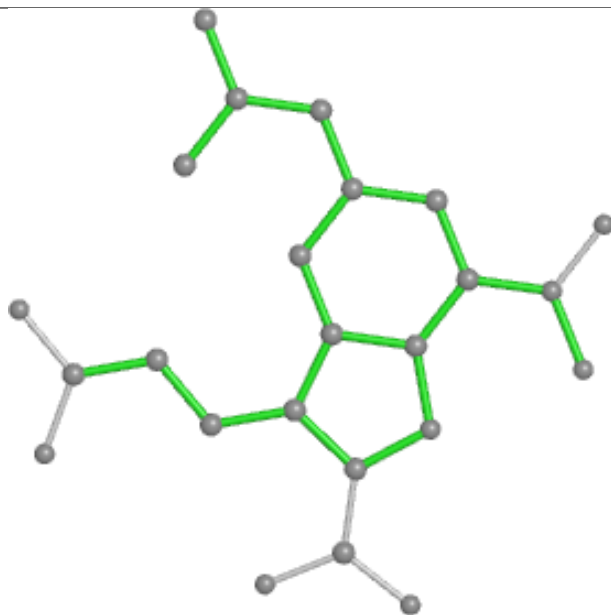
Ligand 0GI B 402



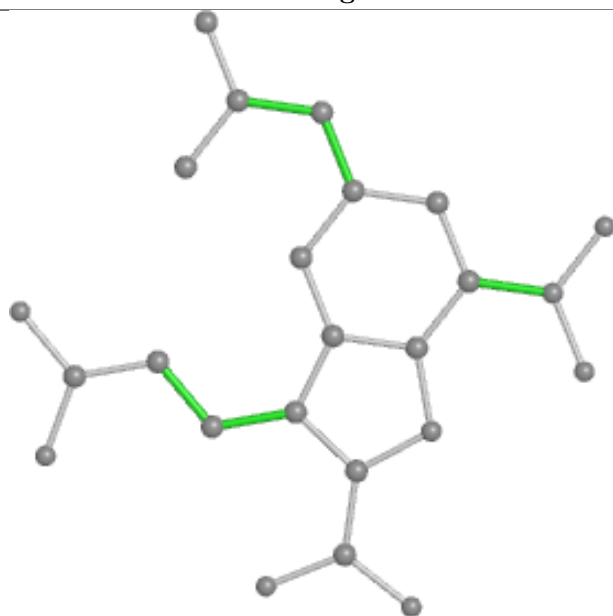
Ligand 0GI D 402



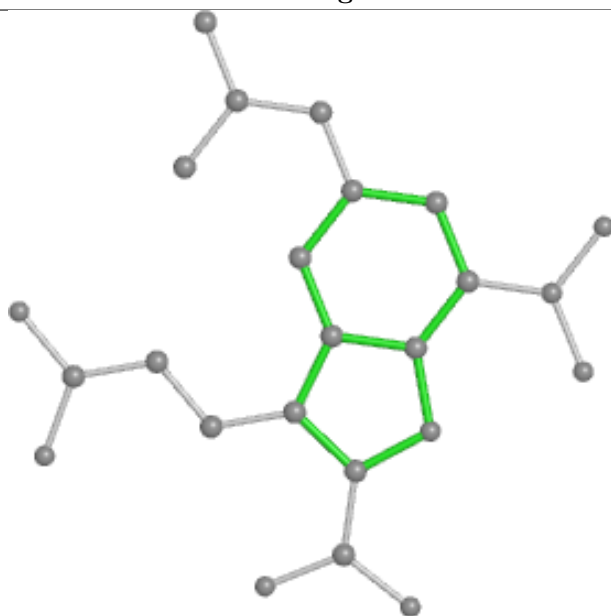
Bond lengths



Bond angles

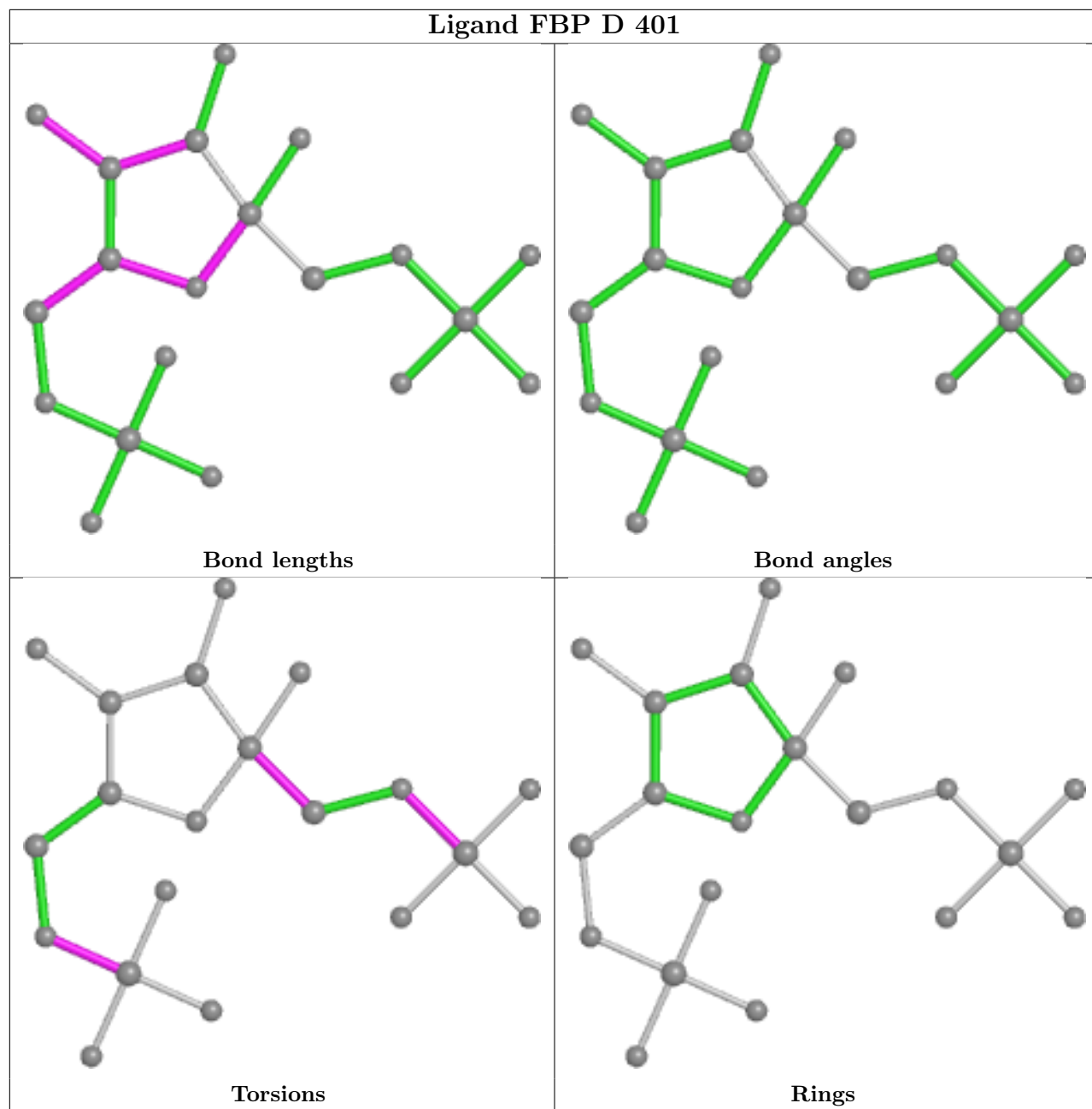


Torsions

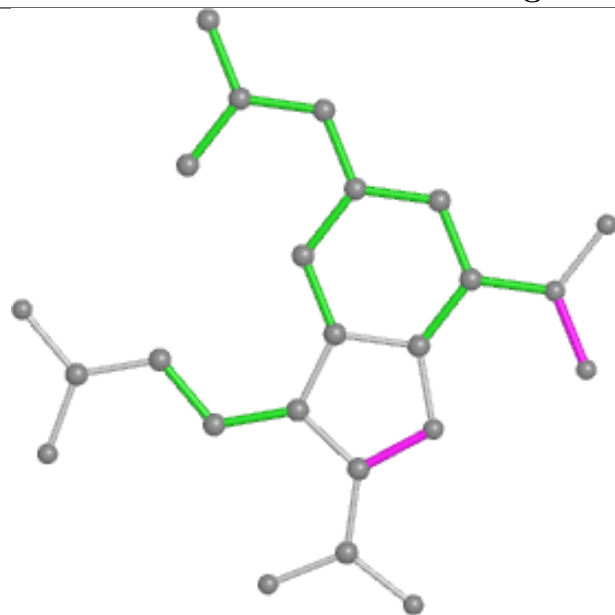


Rings

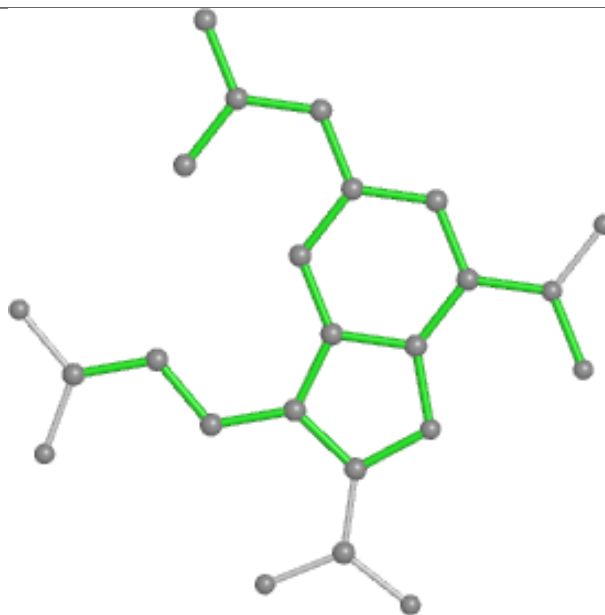
Ligand FBP D 401



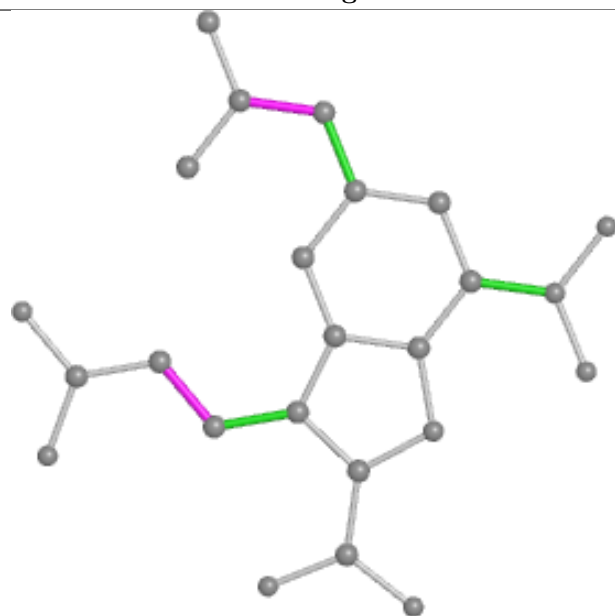
Ligand 0GI C 402



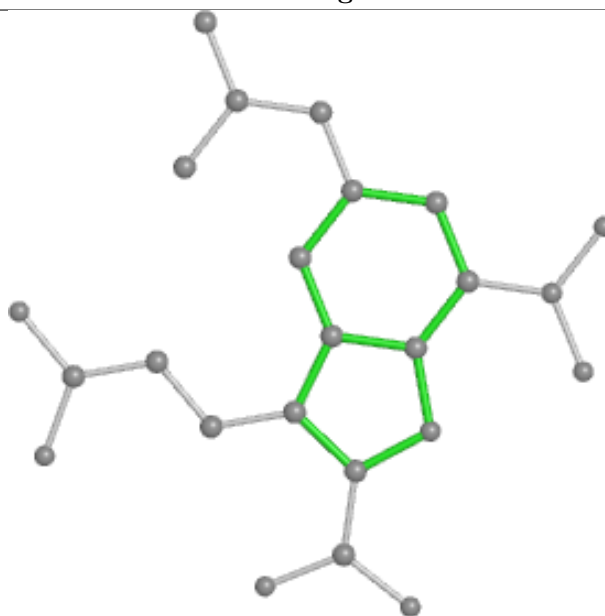
Bond lengths



Bond angles

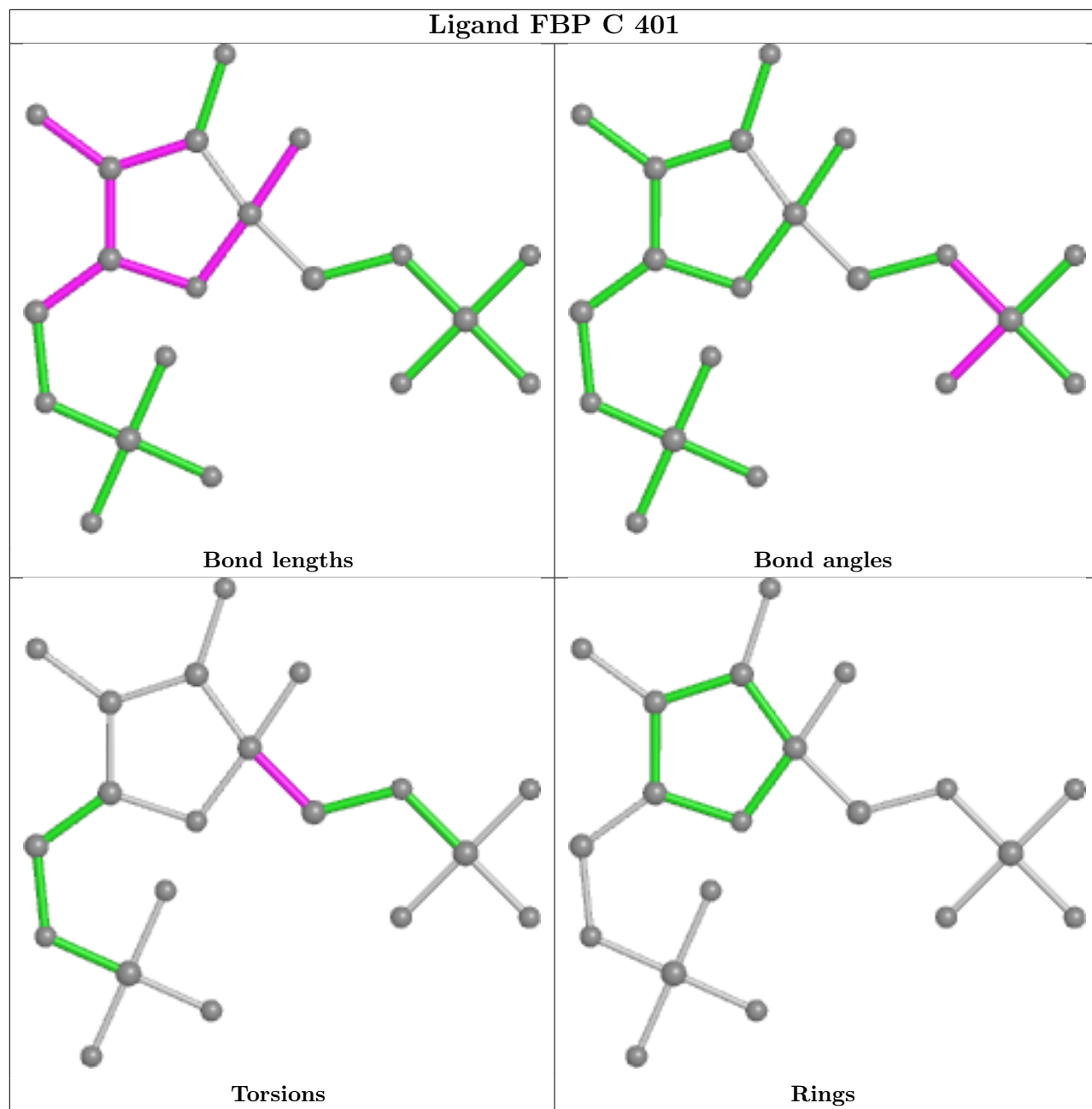


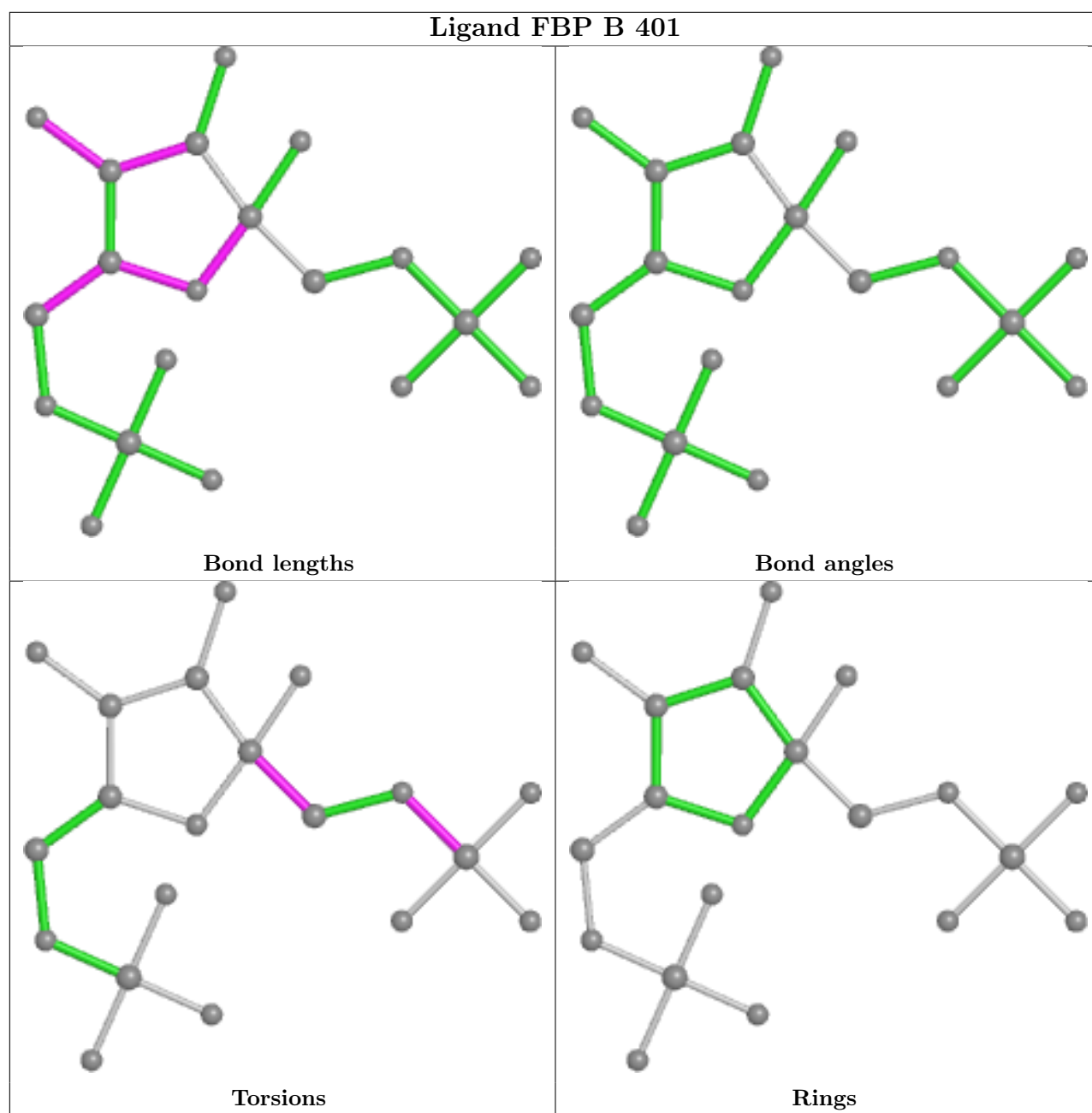
Torsions



Rings

Ligand FBP C 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, 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	320/338 (94%)	-0.02	2 (0%) 89 86	19, 37, 66, 81	0
1	B	320/338 (94%)	-0.10	5 (1%) 72 66	17, 35, 66, 89	0
1	C	319/338 (94%)	-0.06	1 (0%) 94 93	19, 36, 64, 82	0
1	D	320/338 (94%)	-0.10	6 (1%) 66 59	17, 34, 61, 91	0
All	All	1279/1352 (94%)	-0.07	14 (1%) 80 75	17, 36, 65, 91	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	267	ASN	3.4
1	B	235	ASP	3.4
1	B	335	SER	3.2
1	A	234	PRO	2.6
1	D	144	THR	2.6
1	A	270	SER	2.5
1	B	236	ASN	2.5
1	D	266	ALA	2.4
1	C	9	ASP	2.2
1	B	234	PRO	2.1
1	D	270	SER	2.1
1	D	335	SER	2.1
1	D	268	LYS	2.1
1	B	8	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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](#)

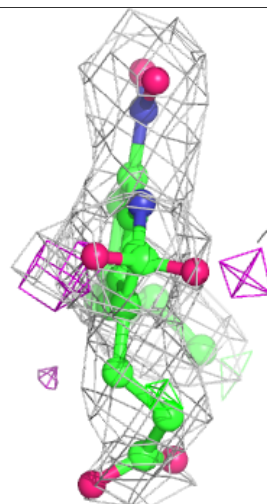
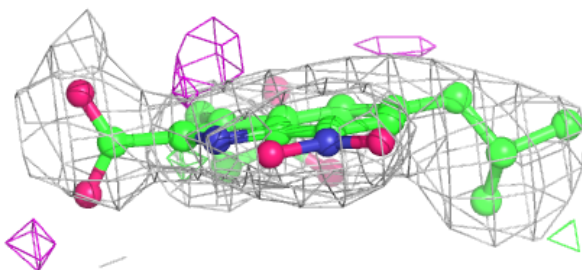
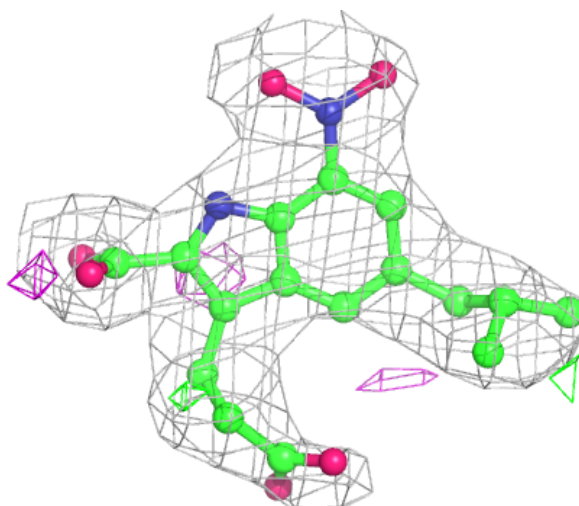
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	0GI	A	402	24/24	0.88	0.25	31,36,54,59	0
3	0GI	B	402	24/24	0.92	0.25	30,38,50,51	0
3	0GI	C	402	24/24	0.92	0.22	38,46,52,65	0
2	FBP	D	401	20/20	0.93	0.18	33,41,44,45	0
2	FBP	A	401	20/20	0.93	0.18	32,40,52,53	0
3	0GI	D	402	24/24	0.93	0.22	29,32,45,46	0
2	FBP	B	401	20/20	0.95	0.16	32,38,47,48	0
2	FBP	C	401	20/20	0.96	0.15	34,40,48,52	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.

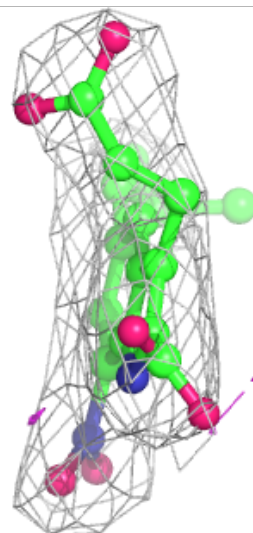
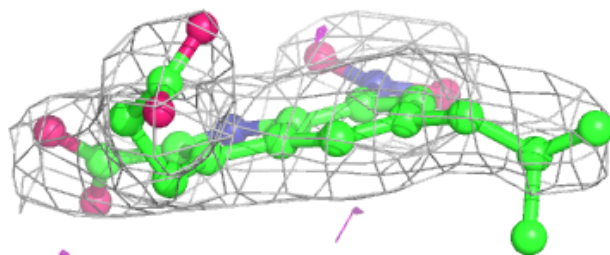
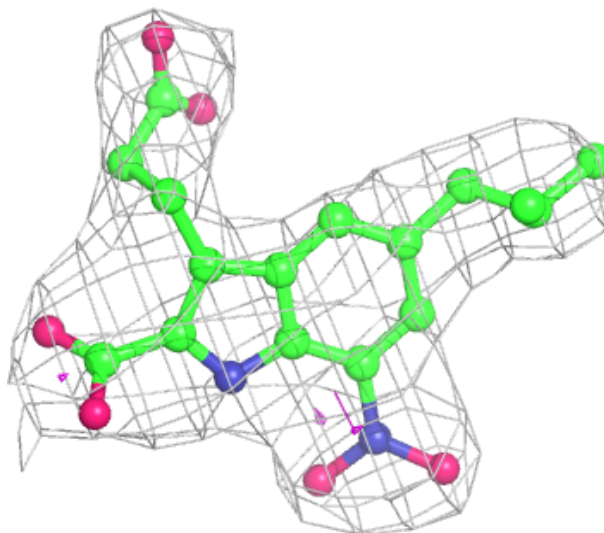
Electron density around 0GI A 402:

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



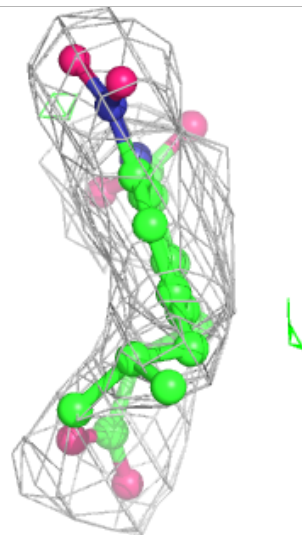
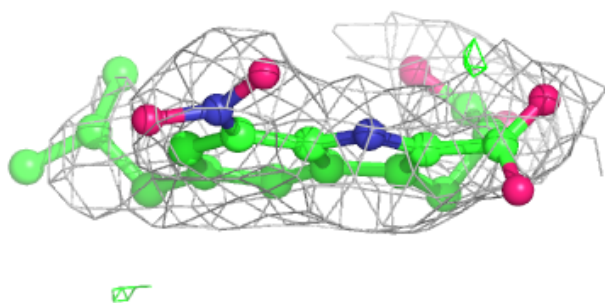
Electron density around 0GI B 402:

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



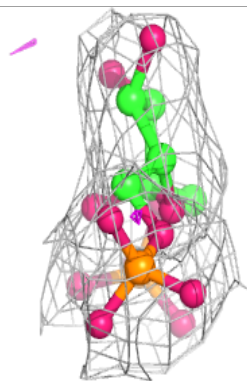
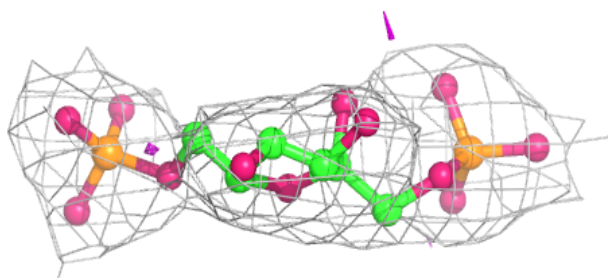
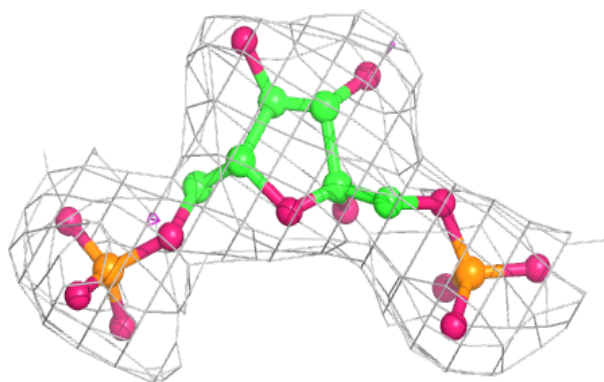
Electron density around 0GI C 402:

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

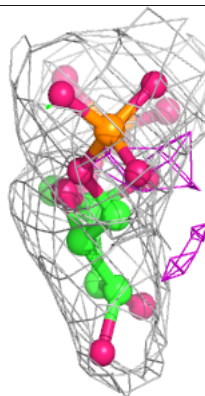
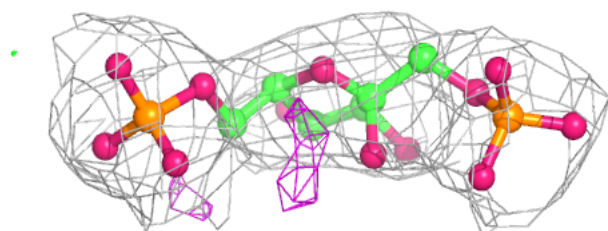
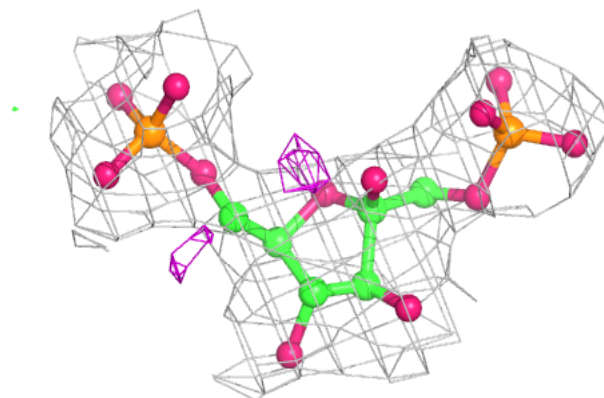


Electron density around FBP D 401:

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

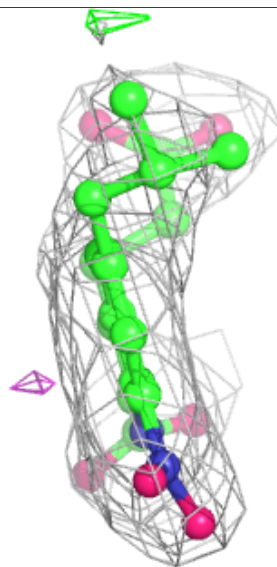
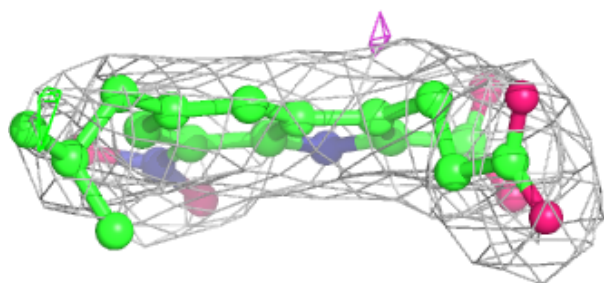
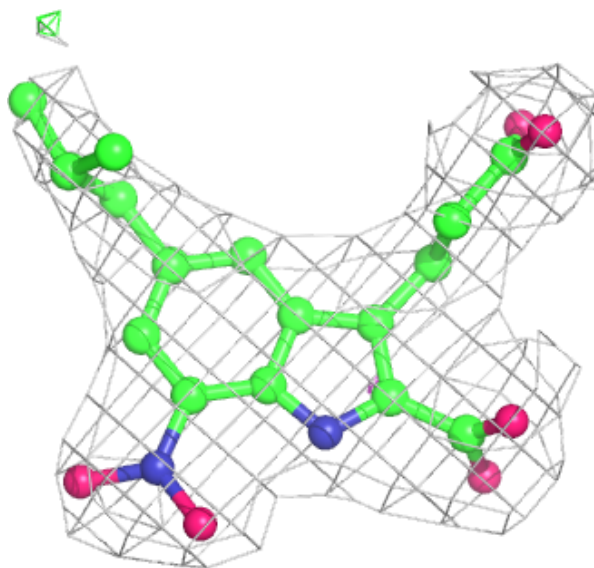
**Electron density around FBP A 401:**

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



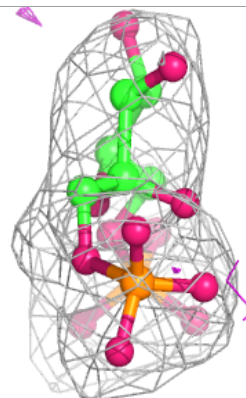
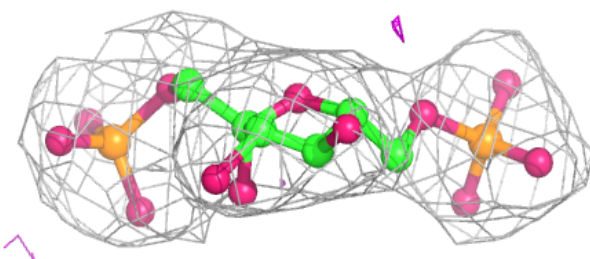
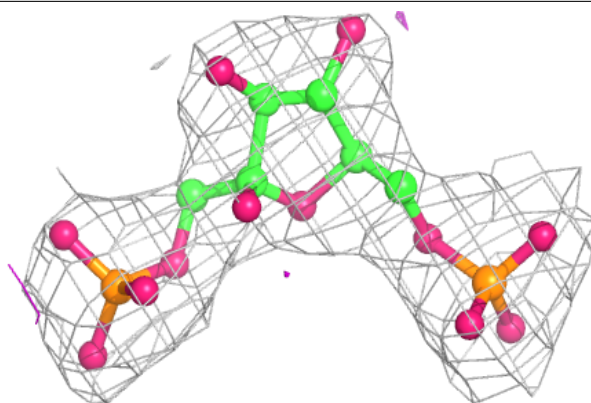
Electron density around 0GI D 402:

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

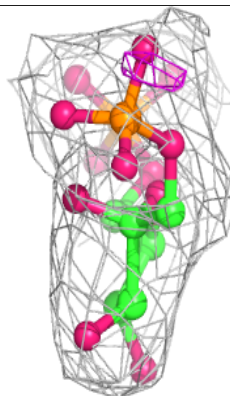
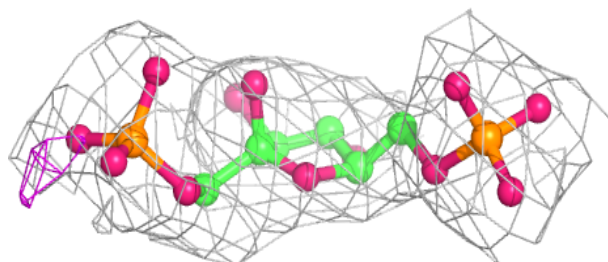
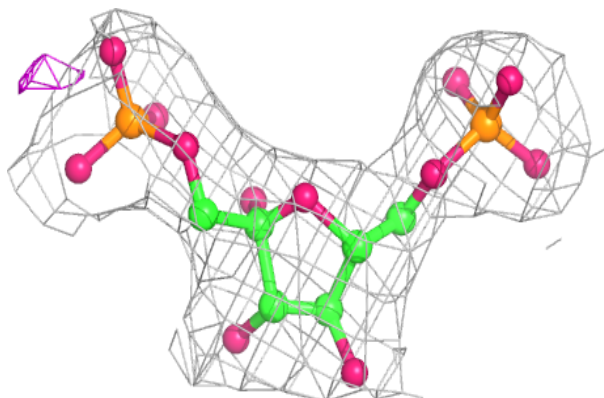


Electron density around FBP B 401:

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

**Electron density around FBP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.