



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 06:20 AM JST

PDB ID : 7F20
Title : L-lactate oxidase with L-lactate
Authors : Morimoto, Y.; Inaka, K.
Deposited on : 2021-06-10
Resolution : 1.30 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

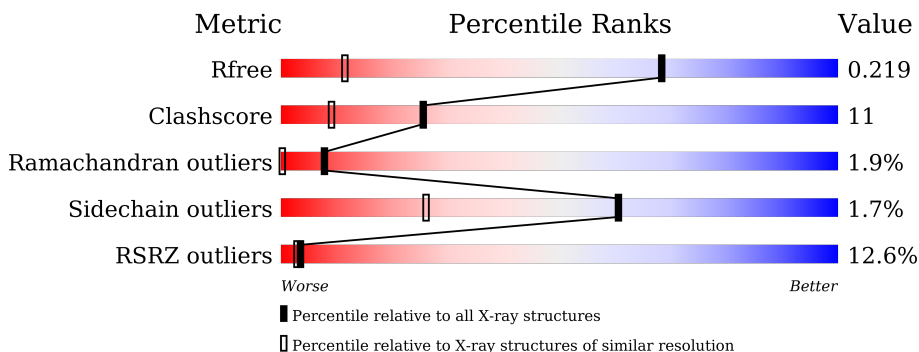
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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	720	 5% 47% 1% 49%
1	B	720	 8% 42% 7% 49%

2 Entry composition i

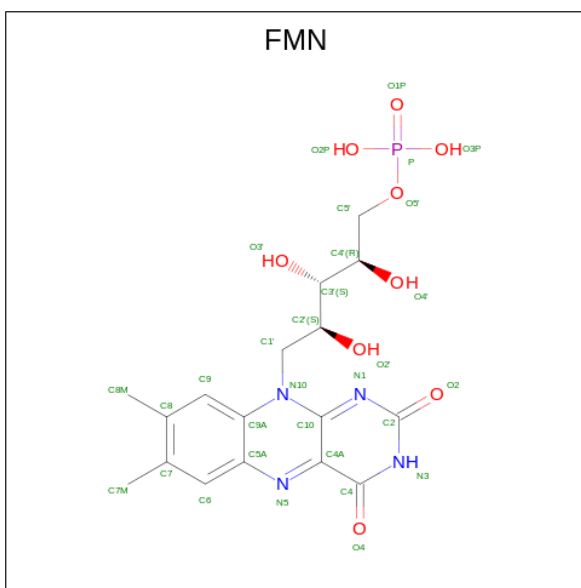
There are 4 unique types of molecules in this entry. The entry contains 6496 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 L-lactate oxidase.

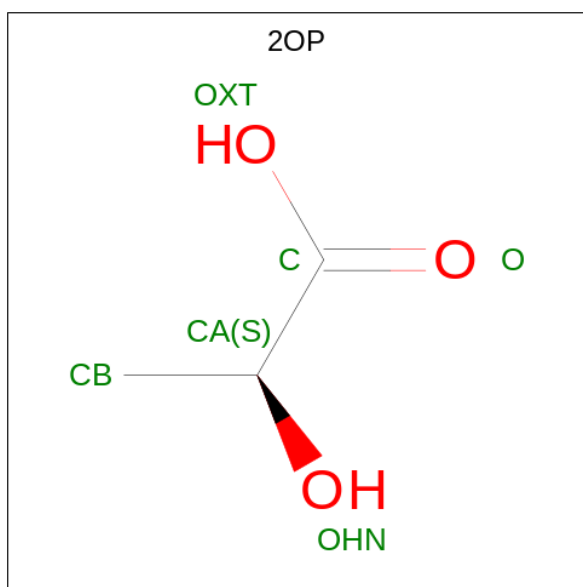
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	368	Total 2841	C 1795	N 494	O 545	S 7	0	0	0
1	B	368	Total 2841	C 1795	N 494	O 545	S 7	0	0	0

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 31	C 17	N 4	O 9	P 1	0	0
2	B	1	Total 31	C 17	N 4	O 9	P 1	0	0

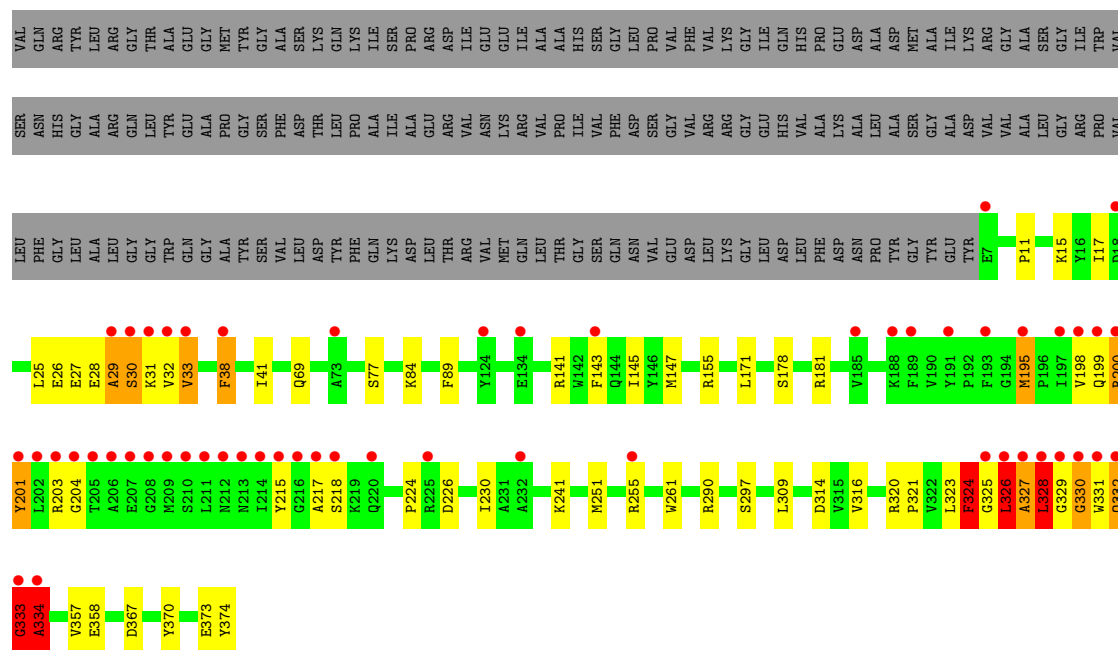
- Molecule 3 is (2S)-2-HYDROXYPROPANOIC ACID (three-letter code: 2OP) (formula: $C_3H_6O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	360	Total O 360 360	0	0
4	B	380	Total O 380 380	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	132.97Å 132.97Å 92.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.05 – 1.30 47.01 – 1.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.05-1.30) 99.9 (47.01-1.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.23 (at 1.30Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.194 , 0.211 0.203 , 0.219	Depositor DCC
R_{free} test set	9948 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	16.7	Xtrriage
Anisotropy	0.176	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.015 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6496	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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: 2OP, FMN

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	0.74	0/2905	0.88	4/3937 (0.1%)
1	B	0.80	1/2905 (0.0%)	0.96	10/3937 (0.3%)
All	All	0.77	1/5810 (0.0%)	0.92	14/7874 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	333	GLY	C-O	10.21	1.40	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	324	PHE	CB-CA-C	7.85	126.10	110.40
1	B	334	ALA	N-CA-CB	-7.11	100.15	110.10
1	B	333	GLY	CA-C-O	-6.75	108.44	120.60
1	B	333	GLY	O-C-N	6.56	133.20	122.70
1	A	370	TYR	CB-CG-CD1	6.42	124.85	121.00
1	A	370	TYR	CB-CG-CD2	-6.36	117.19	121.00
1	B	370	TYR	CB-CG-CD1	6.27	124.76	121.00
1	B	370	TYR	CB-CG-CD2	-6.20	117.28	121.00
1	B	334	ALA	N-CA-C	5.37	125.49	111.00
1	B	374	TYR	CA-C-O	-5.34	108.89	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	334	ALA	O-C-N	-5.34	114.16	122.70
1	B	333	GLY	C-N-CA	5.19	134.67	121.70
1	A	255	ARG	CB-CG-CD	5.13	124.93	111.60
1	A	374	TYR	CA-C-O	-5.01	109.58	120.10

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	32	VAL	Peptide
1	B	33	VAL	Peptide
1	B	330	GLY	Peptide
1	B	333	GLY	Peptide

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	2841	0	2772	16	0
1	B	2841	0	2772	105	1
2	A	31	0	19	1	0
2	B	31	0	19	0	0
3	A	6	0	5	1	0
3	B	6	0	5	1	0
4	A	360	0	0	3	1
4	B	380	0	0	19	4
All	All	6496	0	5592	120	5

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 11.

All (120) 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:331:TRP:CE3	4:B:771:HOH:O	1.83	1.25

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:GLU:O	1:B:329:GLY:N	1.76	1.17
1:B:29:ALA:CB	1:B:326:LEU:O	1.91	1.17
1:B:29:ALA:HB3	1:B:326:LEU:O	1.44	1.16
1:B:334:ALA:HB3	4:B:553:HOH:O	1.47	1.13
1:B:33:VAL:HG11	1:B:326:LEU:CA	1.80	1.12
1:B:33:VAL:HG11	1:B:326:LEU:HA	1.09	1.07
1:B:38:PHE:HA	1:B:326:LEU:HD21	1.12	1.04
1:B:69:GLN:O	4:B:501:HOH:O	1.75	1.04
1:B:224:PRO:HB3	1:B:255:ARG:HG3	1.42	1.01
1:B:325:GLY:HA2	4:B:771:HOH:O	1.57	1.01
1:B:143:PHE:CE2	1:B:145:ILE:HG22	1.97	0.98
1:B:331:TRP:HE3	4:B:771:HOH:O	1.28	0.98
1:B:41:ILE:HD12	1:B:326:LEU:HD22	1.46	0.98
1:A:251:MET:O	1:A:255:ARG:HG2	1.64	0.96
1:B:29:ALA:HB3	1:B:326:LEU:C	1.90	0.93
1:B:251:MET:O	1:B:255:ARG:HG2	1.68	0.92
1:B:33:VAL:CG1	1:B:326:LEU:HA	1.98	0.92
1:B:38:PHE:HA	1:B:326:LEU:CD2	1.98	0.92
1:B:323:LEU:O	1:B:325:GLY:N	2.02	0.91
1:B:30:SER:HB3	1:B:328:LEU:N	1.87	0.90
1:B:30:SER:HB2	4:B:667:HOH:O	1.74	0.87
1:B:38:PHE:CE2	1:B:327:ALA:HB1	2.10	0.86
1:B:155:ARG:NH2	4:B:503:HOH:O	1.97	0.86
1:B:33:VAL:CG2	4:B:523:HOH:O	2.22	0.86
1:B:314:ASP:OD2	4:B:502:HOH:O	1.95	0.85
1:B:143:PHE:HE2	1:B:145:ILE:HG22	1.36	0.84
1:A:224:PRO:HB3	1:A:255:ARG:HG3	1.59	0.84
1:B:201:TYR:CZ	1:B:217:ALA:O	2.33	0.82
1:B:17:ILE:HD12	1:B:332:GLN:HB3	1.63	0.81
1:B:38:PHE:CD2	1:B:327:ALA:CB	2.64	0.81
1:B:324:PHE:O	1:B:332:GLN:C	2.21	0.79
1:B:29:ALA:H	1:B:330:GLY:N	1.80	0.79
1:B:77:SER:HB2	1:B:84:LYS:HE3	1.65	0.78
1:B:29:ALA:HB1	1:B:326:LEU:O	1.83	0.77
1:B:333:GLY:N	4:B:504:HOH:O	2.05	0.75
1:B:29:ALA:CB	1:B:326:LEU:C	2.51	0.74
1:B:226:ASP:O	1:B:230:ILE:HG12	1.87	0.74
1:B:333:GLY:O	4:B:504:HOH:O	2.06	0.72
1:B:38:PHE:CA	1:B:326:LEU:HD21	2.07	0.72
1:B:143:PHE:CE2	1:B:145:ILE:CG2	2.72	0.71
1:B:29:ALA:HB2	1:B:331:TRP:HA	1.70	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:GLY:C	4:B:504:HOH:O	2.28	0.71
1:B:200:ARG:NH1	4:B:507:HOH:O	2.23	0.70
1:B:33:VAL:HG23	4:B:523:HOH:O	1.86	0.69
1:B:331:TRP:O	1:B:332:GLN:CG	2.40	0.69
1:B:27:GLU:HA	1:B:328:LEU:CB	2.23	0.69
1:B:30:SER:H	1:B:329:GLY:N	1.91	0.68
1:B:38:PHE:CE2	1:B:327:ALA:CB	2.75	0.68
1:B:28:GLU:HB3	1:B:330:GLY:O	1.94	0.67
1:B:373:GLU:OE2	4:B:505:HOH:O	2.12	0.67
1:B:195:MET:HB3	1:B:198:VAL:HG22	1.77	0.67
1:B:367:ASP:OD2	4:B:506:HOH:O	2.13	0.67
1:B:11:PRO:HG3	1:B:33:VAL:HG22	1.76	0.66
1:B:27:GLU:HA	1:B:328:LEU:HB2	1.78	0.66
1:B:41:ILE:HD12	1:B:326:LEU:CD2	2.24	0.66
1:B:41:ILE:CD1	1:B:326:LEU:HD22	2.22	0.66
1:B:29:ALA:H	1:B:329:GLY:C	1.98	0.66
1:A:61:LYS:NZ	1:A:367:ASP:OD1	2.21	0.64
1:B:331:TRP:O	1:B:332:GLN:HG3	1.96	0.64
1:B:38:PHE:CD2	1:B:327:ALA:HB2	2.31	0.64
1:B:30:SER:H	1:B:328:LEU:C	2.01	0.64
1:A:225:ARG:NH1	4:A:503:HOH:O	2.32	0.61
1:B:324:PHE:O	1:B:332:GLN:O	2.17	0.61
1:B:201:TYR:CE1	1:B:217:ALA:O	2.53	0.61
1:B:26:GLU:O	1:B:327:ALA:HB3	2.00	0.60
1:B:326:LEU:O	1:B:327:ALA:O	2.21	0.59
1:B:38:PHE:CD2	1:B:327:ALA:HA	2.38	0.59
1:B:30:SER:HB3	1:B:328:LEU:CA	2.33	0.58
1:B:181:ARG:NH1	1:B:215:TYR:OH	2.37	0.57
1:B:38:PHE:CZ	1:B:328:LEU:HD13	2.40	0.57
1:B:33:VAL:HG21	4:B:523:HOH:O	1.95	0.56
1:B:30:SER:OG	1:B:30:SER:O	2.23	0.56
1:B:147:MET:HG2	1:B:171:LEU:HD11	1.88	0.55
1:B:33:VAL:HG11	1:B:326:LEU:C	2.27	0.55
1:B:38:PHE:CG	1:B:327:ALA:HA	2.42	0.54
1:B:331:TRP:O	1:B:332:GLN:HG2	2.09	0.53
1:B:27:GLU:HA	1:B:328:LEU:HB3	1.91	0.52
1:B:29:ALA:O	1:B:33:VAL:CG2	2.59	0.51
1:B:38:PHE:CD2	1:B:327:ALA:CA	2.93	0.51
1:B:29:ALA:HB3	1:B:327:ALA:O	2.10	0.50
1:B:323:LEU:O	1:B:324:PHE:C	2.50	0.50
1:B:324:PHE:O	1:B:333:GLY:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:401:FMN:N5	3:A:402:2OP:HA	2.28	0.48
1:B:143:PHE:CD2	1:B:145:ILE:CG2	2.97	0.47
1:A:251:MET:HG2	1:A:255:ARG:HE	1.79	0.47
1:B:31:LYS:HB3	4:B:832:HOH:O	2.15	0.47
1:B:33:VAL:CG1	1:B:326:LEU:HD12	2.45	0.47
1:B:201:TYR:CE2	1:B:218:SER:HA	2.50	0.47
1:B:29:ALA:C	1:B:326:LEU:O	2.55	0.46
1:B:28:GLU:O	1:B:29:ALA:CB	2.65	0.45
1:B:77:SER:HB2	1:B:84:LYS:CE	2.42	0.45
1:A:199:GLN:NE2	4:A:519:HOH:O	2.49	0.45
1:B:38:PHE:HB2	1:B:326:LEU:HG	1.98	0.45
1:A:212:ASN:ND2	4:A:523:HOH:O	2.49	0.45
1:B:224:PRO:HB3	1:B:255:ARG:CG	2.30	0.45
1:A:273:ALA:HB1	1:A:274:PRO:HD2	1.99	0.44
1:B:215:TYR:OH	3:B:402:2OP:HB3	2.18	0.44
1:B:29:ALA:CA	1:B:326:LEU:O	2.63	0.43
1:B:38:PHE:HD2	1:B:327:ALA:HB2	1.82	0.43
1:A:358:GLU:O	1:B:290:ARG:HD3	2.18	0.43
1:B:178:SER:HB2	1:B:215:TYR:CD2	2.54	0.43
1:B:334:ALA:CA	4:B:553:HOH:O	2.65	0.42
1:B:241:LYS:HA	1:B:261:TRP:HB3	2.02	0.42
1:A:89:PHE:HA	1:A:316:VAL:O	2.19	0.42
1:B:29:ALA:HB1	1:B:326:LEU:C	2.34	0.42
1:A:66:ARG:C	1:A:67:LEU:HD23	2.40	0.42
1:B:29:ALA:O	1:B:33:VAL:HG23	2.19	0.42
1:A:22:THR:HB	1:A:321:PRO:HB3	2.02	0.42
1:B:89:PHE:HA	1:B:316:VAL:O	2.20	0.42
1:B:15:LYS:HG2	1:B:331:TRP:CZ2	2.55	0.41
1:B:77:SER:CB	1:B:84:LYS:HE3	2.43	0.41
1:B:309:LEU:HD23	1:B:357:VAL:HG13	2.03	0.41
1:A:199:GLN:HE21	1:A:199:GLN:HB3	1.69	0.41
1:B:33:VAL:O	1:B:33:VAL:HG12	2.19	0.41
1:B:320:ARG:N	1:B:321:PRO:HD2	2.36	0.41
1:B:25:LEU:HD22	1:B:332:GLN:CB	2.51	0.40
1:A:153:GLN:OE1	1:A:201:TYR:HB3	2.21	0.40
1:A:290:ARG:HD3	1:B:358:GLU:O	2.21	0.40
1:A:241:LYS:HA	1:A:261:TRP:HB3	2.04	0.40

All (5) 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)
4:B:836:HOH:O	4:B:836:HOH:O[3_655]	1.91	0.29
4:A:814:HOH:O	4:A:814:HOH:O[3_655]	2.00	0.20
4:B:601:HOH:O	4:B:601:HOH:O[3_655]	2.03	0.17
4:B:505:HOH:O	4:B:733:HOH:O[4_565]	2.14	0.06
1:B:373:GLU:OE1	4:B:505:HOH:O[3_655]	2.17	0.03

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	366/720 (51%)	354 (97%)	9 (2%)	3 (1%)	19	2
1	B	366/720 (51%)	338 (92%)	17 (5%)	11 (3%)	4	0
All	All	732/1440 (51%)	692 (94%)	26 (4%)	14 (2%)	8	0

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	205	THR
1	A	297	SER
1	B	200	ARG
1	B	297	SER
1	B	324	PHE
1	B	327	ALA
1	B	328	LEU
1	B	334	ALA
1	B	29	ALA
1	B	326	LEU
1	A	203	ARG
1	B	203	ARG
1	B	204	GLY
1	B	332	GLN

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	292/569 (51%)	290 (99%)	2 (1%)	84	61
1	B	292/569 (51%)	284 (97%)	8 (3%)	44	9
All	All	584/1138 (51%)	574 (98%)	10 (2%)	60	26

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	141	ARG
1	A	199	GLN
1	B	30	SER
1	B	38	PHE
1	B	141	ARG
1	B	195	MET
1	B	199	GLN
1	B	201	TYR
1	B	326	LEU
1	B	328	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	199	GLN
1	B	35	HIS
1	B	153	GLN

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](#)

4 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	2OP	A	402	-	5,5,5	1.18	1 (20%)	4,6,6	0.55	0
2	FMN	B	401	-	33,33,33	0.82	0	48,50,50	0.81	1 (2%)
2	FMN	A	401	-	33,33,33	0.76	0	48,50,50	0.80	0
3	2OP	B	402	-	5,5,5	0.94	0	4,6,6	0.47	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	2OP	A	402	-	-	2/4/4/4	-
2	FMN	B	401	-	-	1/18/18/18	0/3/3/3
2	FMN	A	401	-	-	2/18/18/18	0/3/3/3
3	2OP	B	402	-	-	1/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	2OP	OXT-C	-2.45	1.22	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	FMN	C4A-C4-N3	2.18	118.71	113.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

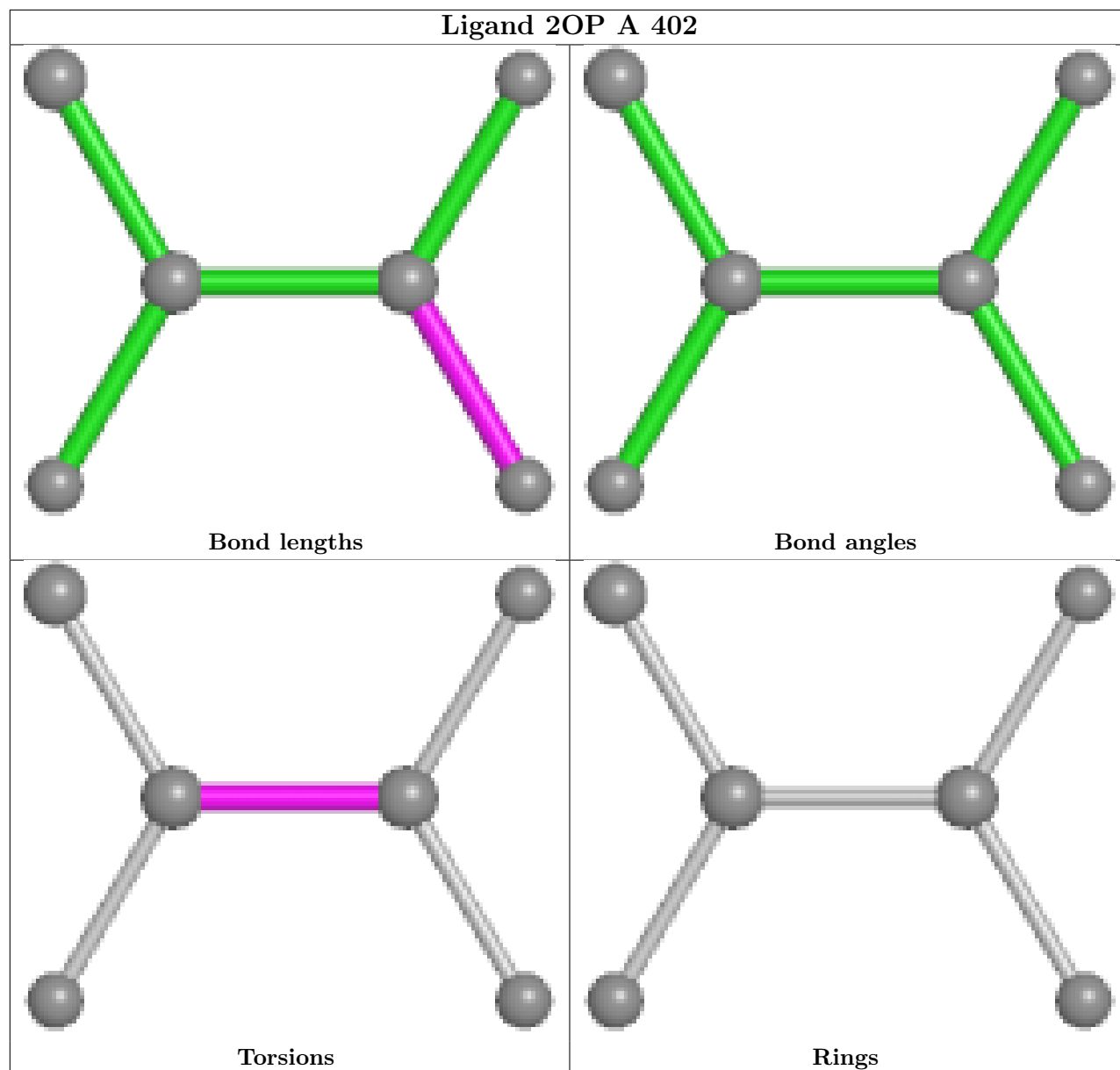
Mol	Chain	Res	Type	Atoms
2	B	401	FMN	C4'-C5'-O5'-P
3	A	402	2OP	O-C-CA-CB
3	A	402	2OP	OXT-C-CA-CB
2	A	401	FMN	C4'-C5'-O5'-P
2	A	401	FMN	C5'-O5'-P-O1P
3	B	402	2OP	OXT-C-CA-CB

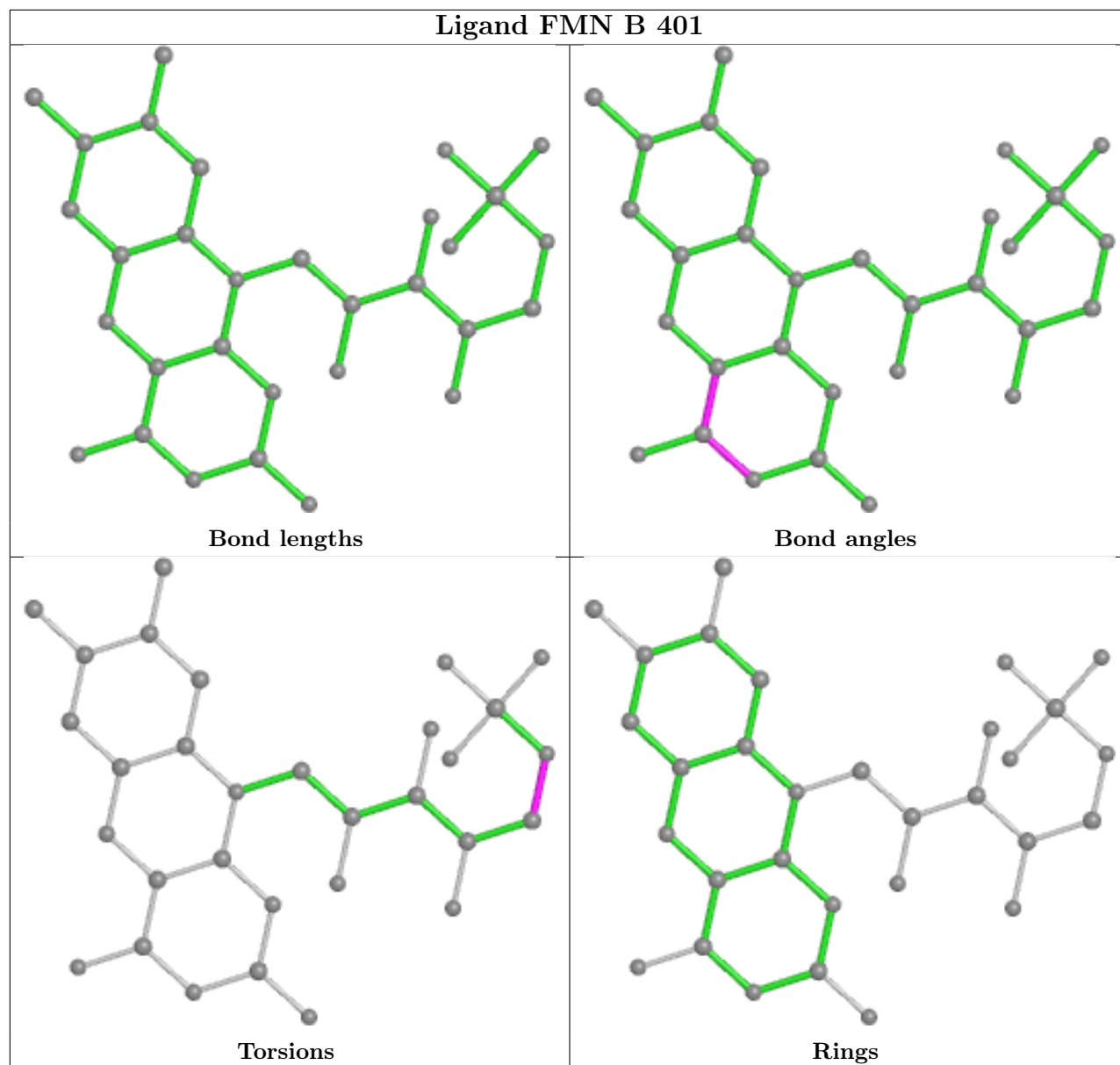
There are no ring outliers.

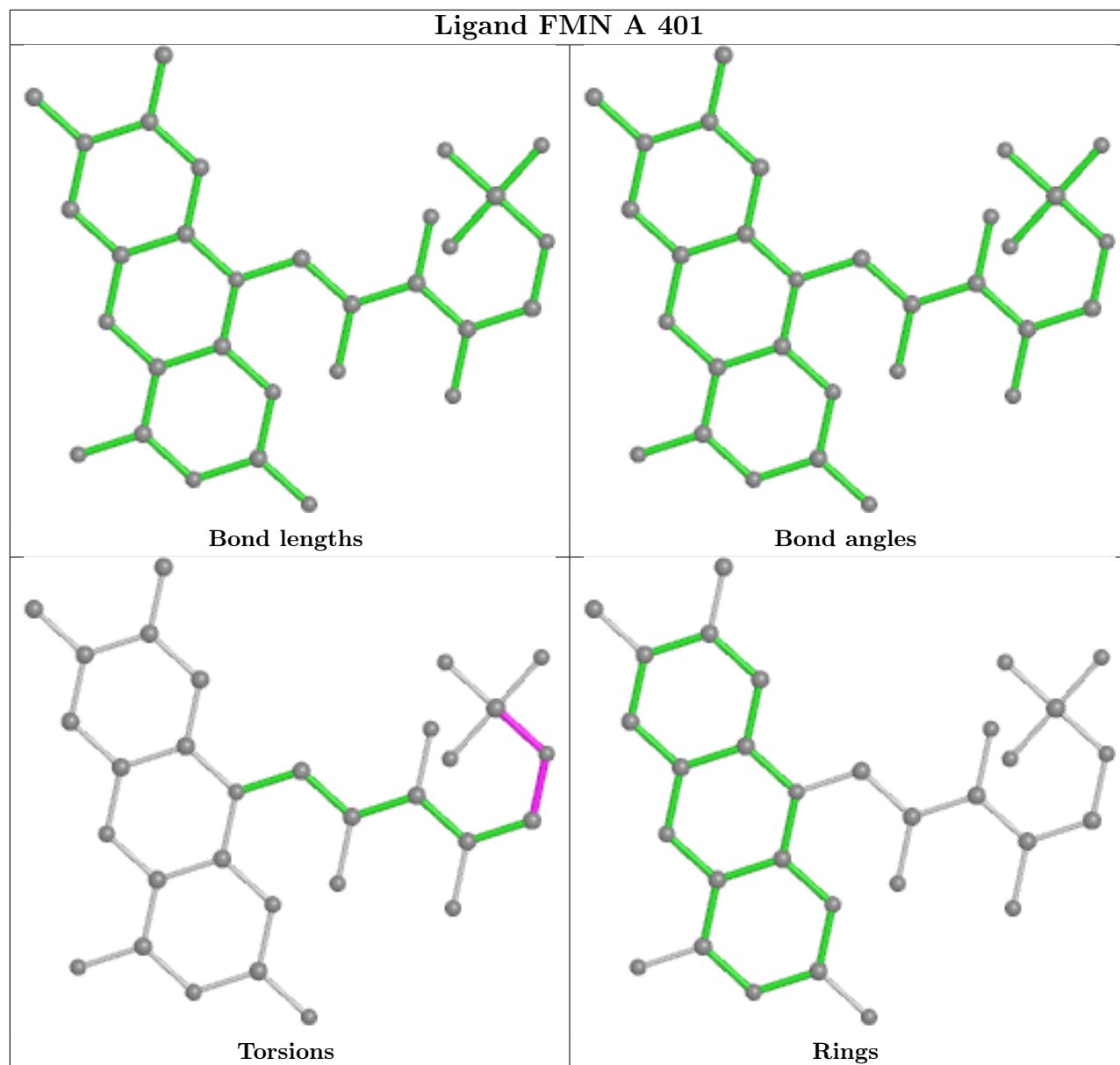
3 monomers are involved in 2 short contacts:

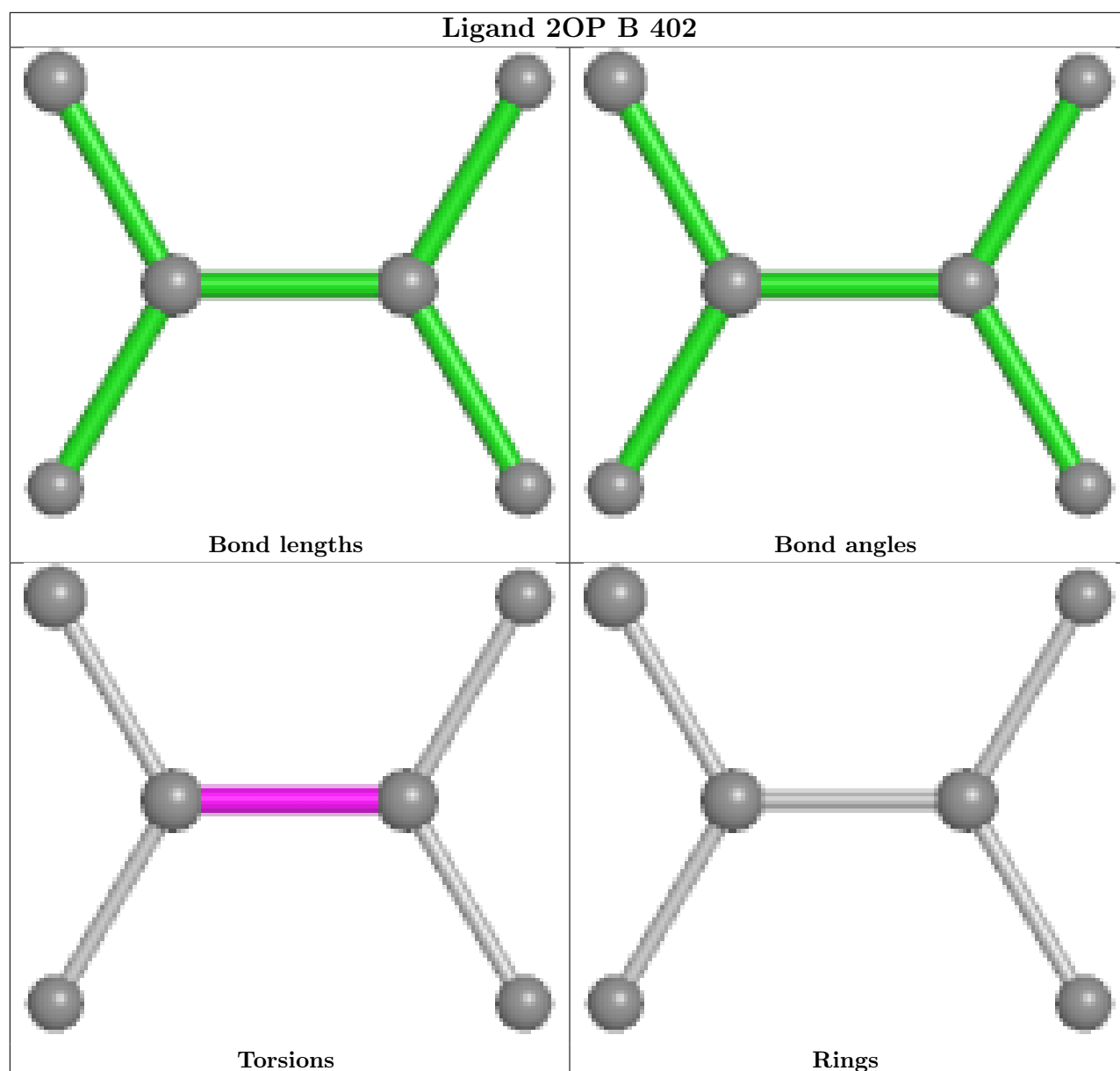
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	2OP	1	0
2	A	401	FMN	1	0
3	B	402	2OP	1	0

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	368/720 (51%)	0.59	39 (10%) 6 4	12, 18, 39, 97	0
1	B	368/720 (51%)	1.76	54 (14%) 2 2	12, 20, 73, 187	0
All	All	736/1440 (51%)	1.18	93 (12%) 3 3	12, 19, 52, 187	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	206	ALA	45.8
1	B	211	LEU	44.5
1	B	326	LEU	31.0
1	B	32	VAL	28.8
1	B	209	MET	25.8
1	B	331	TRP	22.5
1	B	332	GLN	21.1
1	B	330	GLY	19.6
1	B	212	ASN	19.4
1	B	201	TYR	18.8
1	B	208	GLY	18.1
1	B	204	GLY	17.8
1	B	207	GLU	17.7
1	B	210	SER	17.6
1	B	328	LEU	16.7
1	B	205	THR	16.2
1	B	203	ARG	14.4
1	B	327	ALA	12.7
1	A	203	ARG	12.5
1	A	205	THR	12.5
1	A	204	GLY	11.0
1	B	31	LYS	10.9
1	B	325	GLY	10.7
1	B	30	SER	10.6

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Mol	Chain	Res	Type	RSRZ
1	B	214	ILE	10.3
1	B	202	LEU	10.3
1	B	29	ALA	10.0
1	B	329	GLY	10.0
1	B	191	TYR	8.7
1	B	213	ASN	8.3
1	A	208	GLY	7.7
1	B	33	VAL	7.6
1	A	215	TYR	7.5
1	B	216	GLY	6.9
1	B	198	VAL	6.9
1	A	201	TYR	6.9
1	A	198	VAL	6.5
1	B	215	TYR	6.5
1	A	214	ILE	6.5
1	B	199	GLN	6.2
1	A	206	ALA	6.0
1	B	200	ARG	6.0
1	A	197	ILE	5.9
1	A	202	LEU	5.9
1	B	333	GLY	5.9
1	A	211	LEU	5.8
1	A	199	GLN	5.5
1	B	124	TYR	5.4
1	A	7	GLU	5.3
1	A	200	ARG	5.1
1	B	38	PHE	4.9
1	A	212	ASN	4.8
1	B	143	PHE	4.8
1	A	129	PHE	4.7
1	A	157	ILE	4.7
1	B	334	ALA	4.0
1	A	216	GLY	3.8
1	A	209	MET	3.8
1	B	7	GLU	3.5
1	B	185	VAL	3.5
1	B	217	ALA	3.4
1	A	213	ASN	3.4
1	A	207	GLU	3.2
1	A	163	SER	3.2
1	A	210	SER	3.2
1	A	143	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	73	ALA	3.1
1	B	197	ILE	3.1
1	B	189	PHE	3.0
1	B	255	ARG	2.9
1	B	18	ASP	2.9
1	A	230	ILE	2.9
1	B	195	MET	2.8
1	A	231	ALA	2.7
1	A	232	ALA	2.6
1	B	188	LYS	2.6
1	A	195	MET	2.5
1	A	161	ALA	2.5
1	A	158	LEU	2.5
1	A	189	PHE	2.4
1	B	193	PHE	2.4
1	A	124	TYR	2.4
1	A	191	TYR	2.4
1	A	193	PHE	2.4
1	A	145	ILE	2.3
1	B	232	ALA	2.3
1	B	218	SER	2.3
1	A	225	ARG	2.3
1	A	9	ASN	2.2
1	B	134	GLU	2.2
1	B	225	ARG	2.2
1	A	148	ALA	2.1
1	B	220	GLN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

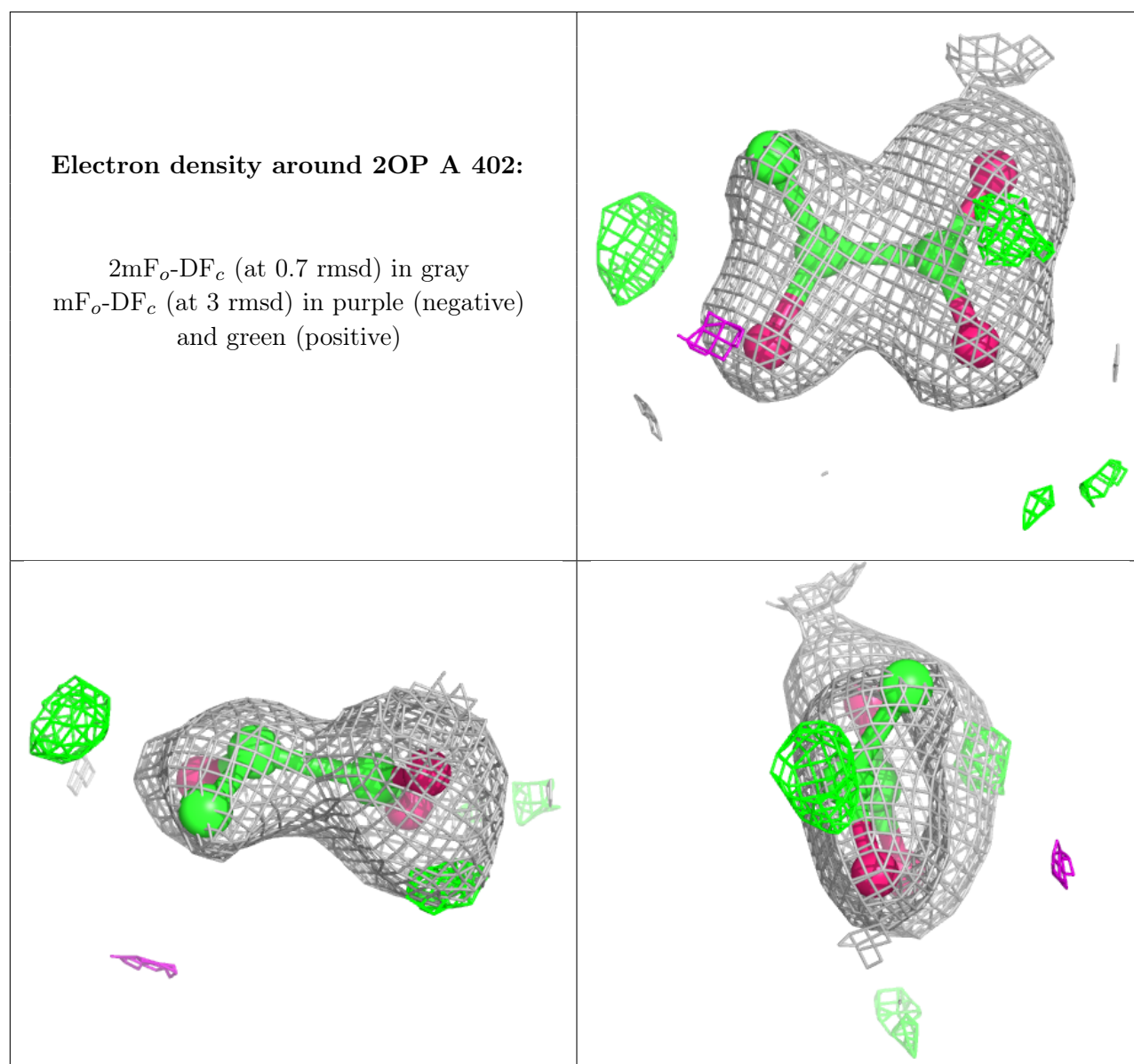
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 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.

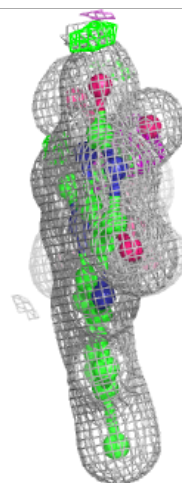
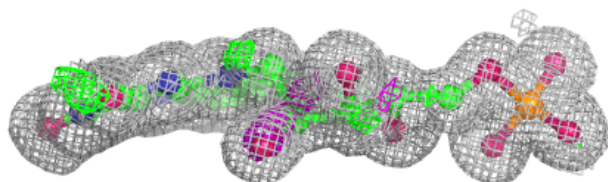
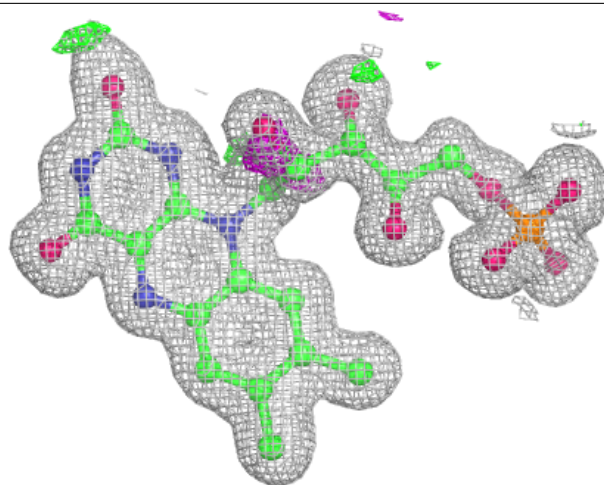
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	2OP	A	402	6/6	0.94	0.11	26,28,30,32	0
2	FMN	B	401	31/31	0.97	0.06	13,14,19,20	0
3	2OP	B	402	6/6	0.97	0.14	27,32,33,37	0
2	FMN	A	401	31/31	0.98	0.06	13,14,18,20	0

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.



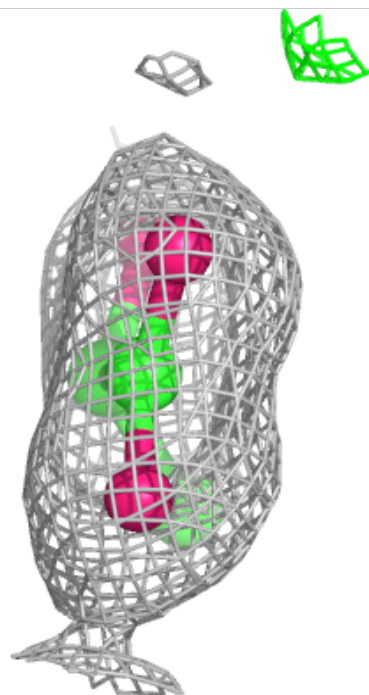
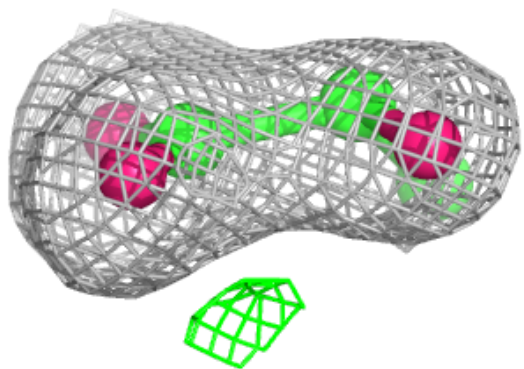
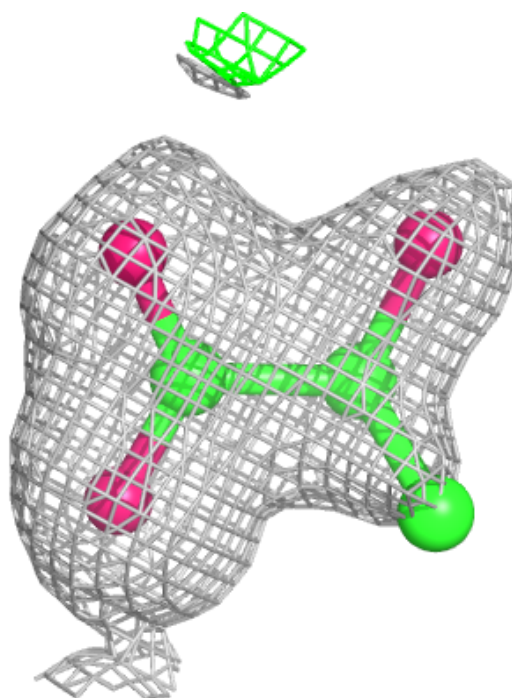
Electron density around FMN B 401:

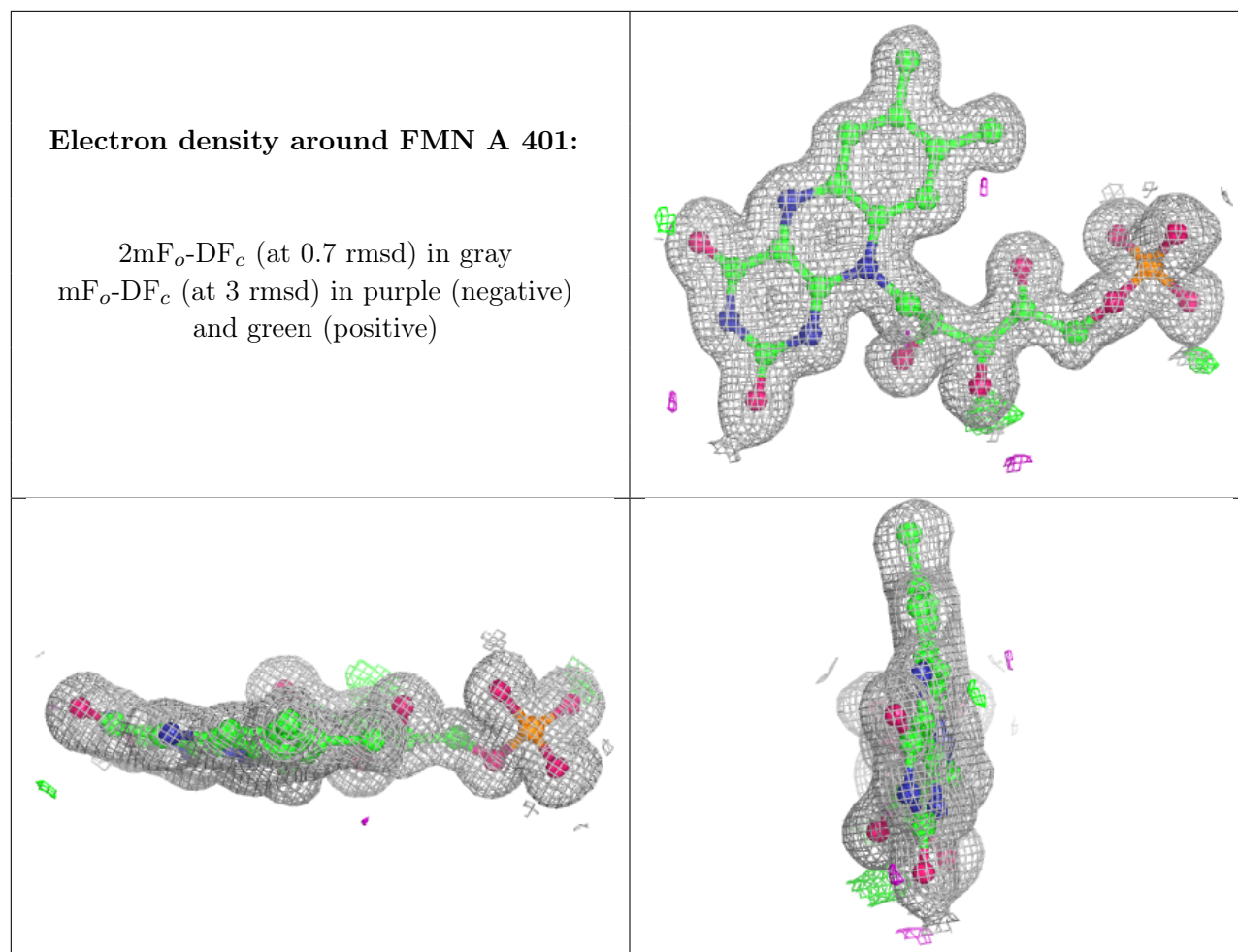
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 2OP 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)





6.5 Other polymers [i](#)

There are no such residues in this entry.