



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:29 PM UTC

PDB ID : 7O0T / pdb_00007o0t
Title : Crystal structure of Chloroflexus aggregans ene-reductase CaOYE holoenzyme
Authors : Robescu, M.S.; Niero, M.; Loprete, G.; Bergantino, E.; Cendron, L.
Deposited on : 2021-03-26
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

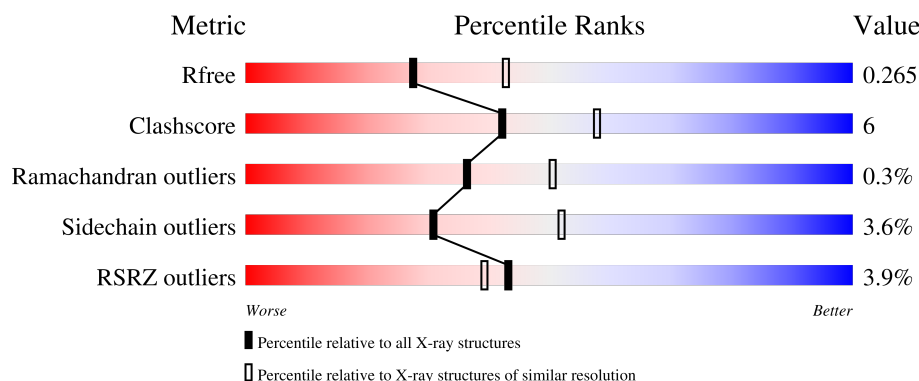
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	354	2% 87% 12% .
1	B	354	% 86% 14% .
1	C	354	2% 86% 13% .
1	D	354	10% 80% 18% .

2 Entry composition [i](#)

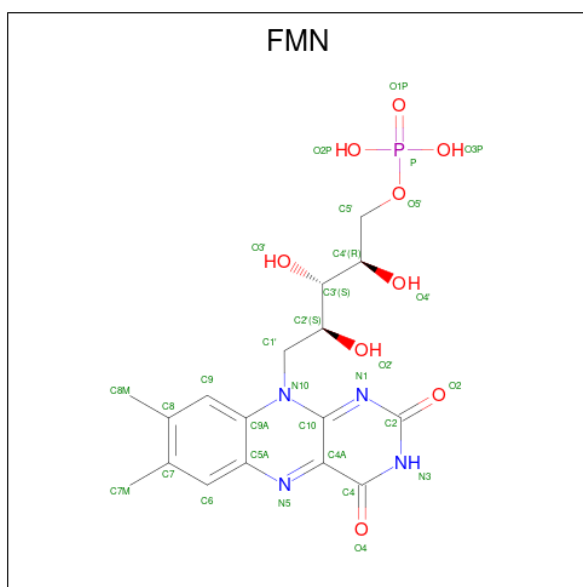
There are 4 unique types of molecules in this entry. The entry contains 11101 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NADH:flavin oxidoreductase/NADH oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	3	0
			2696	1708	497	483	8			
1	B	354	Total	C	N	O	S	0	2	0
			2685	1702	493	482	8			
1	C	354	Total	C	N	O	S	0	4	0
			2707	1714	501	484	8			
1	D	354	Total	C	N	O	S	0	0	0
			2666	1692	487	479	8			

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



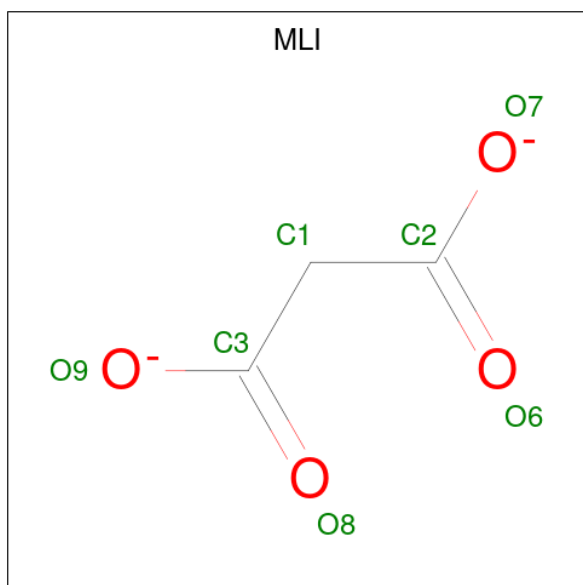
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

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

- Molecule 3 is MALONATE ION (CCD ID: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	3	4		
3	B	1	Total	C	O	0	0
			7	3	4		
3	C	1	Total	C	O	0	0
			7	3	4		
3	D	1	Total	C	O	0	0
			7	3	4		
3	D	1	Total	C	O	0	0
			7	3	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	57	Total	O	0	0
			57	57		
4	B	51	Total	O	0	0
			51	51		

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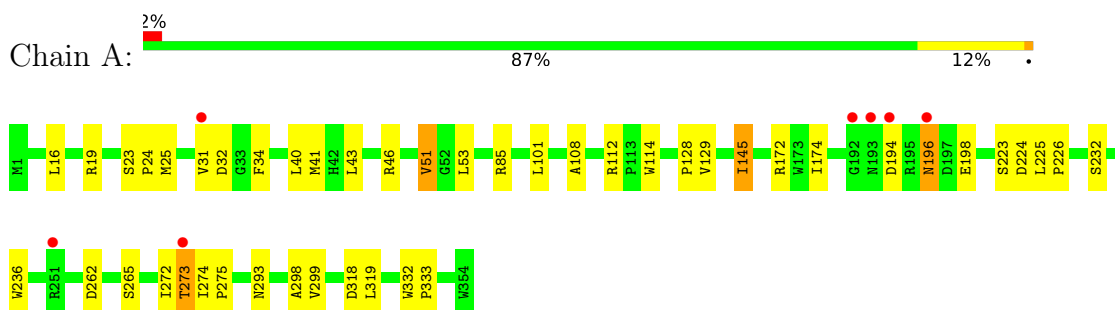
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	51	Total	O	0	0
			51	51		
4	D	29	Total	O	0	0
			29	29		

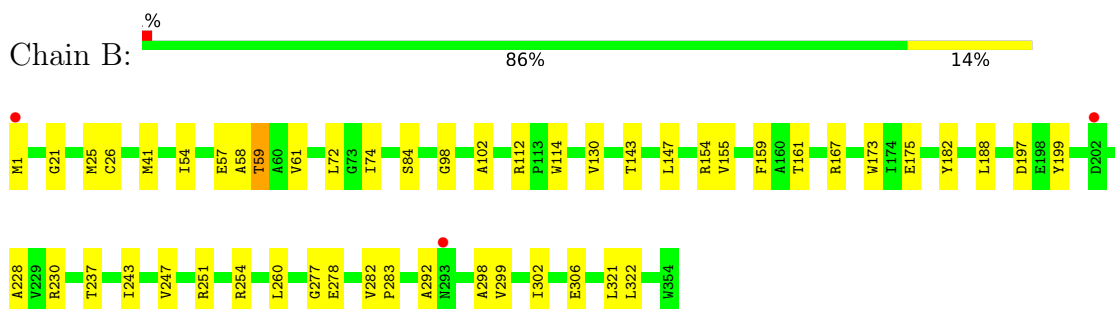
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

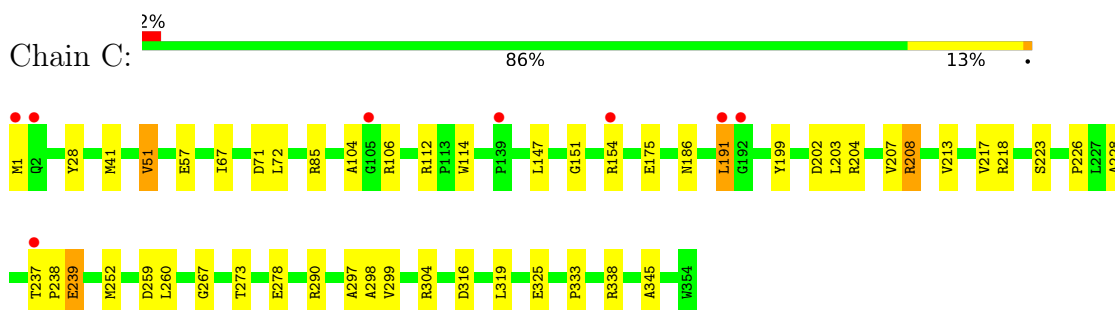
- Molecule 1: NADH:flavin oxidoreductase/NADH oxidase



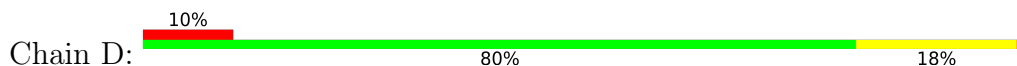
- Molecule 1: NADH:flavin oxidoreductase/NADH oxidase

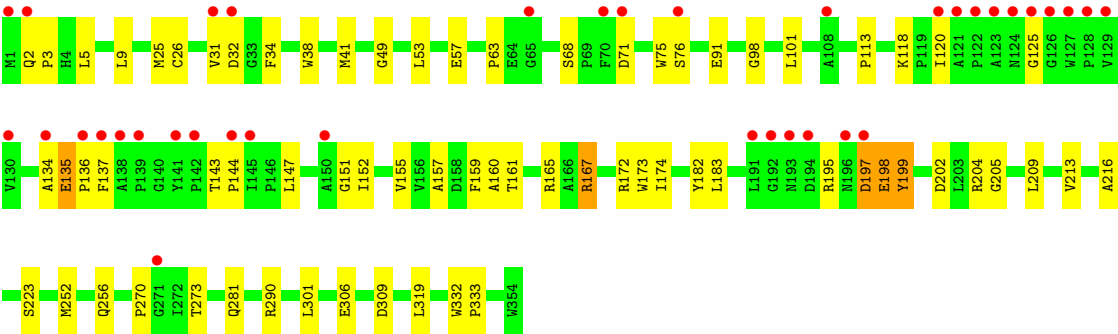


- Molecule 1: NADH:flavin oxidoreductase/NADH oxidase



- Molecule 1: NADH:flavin oxidoreductase/NADH oxidase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	108.19Å 162.27Å 216.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.01 – 2.40 38.01 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.2 (38.01-2.40) 99.2 (38.01-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.220 , 0.263 (Not available) , 0.265	Depositor DCC
R_{free} test set	3677 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 23.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11101	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% 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: MLI, 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	0.65	0/2767	1.15	8/3780 (0.2%)
1	B	0.61	0/2756	1.13	1/3766 (0.0%)
1	C	0.64	0/2778	1.11	2/3794 (0.1%)
1	D	0.62	0/2737	1.15	5/3741 (0.1%)
All	All	0.63	0/11038	1.14	16/15081 (0.1%)

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	A	0	2

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	194	ASP	N-CA-C	-8.64	100.69	114.09
1	C	316	ASP	CA-CB-CG	7.31	119.91	112.60
1	D	223	SER	CA-C-N	6.84	130.13	120.28
1	D	223	SER	C-N-CA	6.84	130.13	120.28
1	A	223	SER	CA-C-N	6.82	132.11	120.72
1	A	223	SER	C-N-CA	6.82	132.11	120.72
1	B	59	THR	CA-CB-OG1	-6.15	100.38	109.60
1	A	262	ASP	CA-CB-CG	5.56	118.16	112.60
1	D	281	GLN	CA-C-N	5.40	123.64	120.24
1	D	281	GLN	C-N-CA	5.40	123.64	120.24
1	C	238	PRO	CB-CA-C	-5.30	104.23	113.20
1	A	196	ASN	N-CA-C	-5.26	106.88	113.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	ASP	N-CA-CB	-5.10	101.92	110.39
1	D	273	THR	CB-CA-C	5.06	118.67	110.22
1	A	262	ASP	CB-CA-C	5.02	118.60	110.22
1	A	273	THR	CB-CA-C	5.01	117.77	110.26

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	293[A]	ASN	Mainchain
1	A	293[B]	ASN	Mainchain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2696	0	2690	30	0
1	B	2685	0	2678	29	0
1	C	2707	0	2702	33	0
1	D	2666	0	2661	49	0
2	A	31	0	19	4	0
2	B	31	0	19	4	0
2	C	31	0	19	2	0
2	D	31	0	19	2	0
3	A	7	0	2	0	0
3	B	7	0	2	1	0
3	C	7	0	2	0	0
3	D	14	0	4	1	0
4	A	57	0	0	2	0
4	B	51	0	0	1	0
4	C	51	0	0	4	0
4	D	29	0	0	1	0
All	All	11101	0	10817	136	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (136) 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:A:85:ARG:NH2	1:C:85:ARG:HH12	1.49	1.10
1:A:85:ARG:HH22	1:C:85:ARG:NH1	1.66	0.94
1:A:85:ARG:HH22	1:C:85:ARG:HH12	0.90	0.89
1:A:31:VAL:HG13	1:A:32:ASP:H	1.47	0.79
1:C:112:ARG:HD3	1:C:114:TRP:CZ2	2.19	0.78
1:D:120:ILE:HG21	1:D:125:GLY:O	1.87	0.74
1:D:31:VAL:HG22	1:D:34:PHE:HB2	1.70	0.74
1:D:167:ARG:HA	1:D:167:ARG:HE	1.51	0.74
1:B:299:VAL:HG22	2:B:401:FMN:H5'1	1.71	0.73
1:D:31:VAL:HG23	1:D:32:ASP:H	1.55	0.72
1:D:134:ALA:O	1:D:144:PRO:HD2	1.89	0.72
1:B:112:ARG:HD3	1:B:114:TRP:CZ2	2.28	0.69
1:D:161:THR:HG22	1:D:165:ARG:CZ	2.23	0.69
1:C:199:TYR:CE1	1:C:208[A]:ARG:HG2	2.29	0.66
1:A:31:VAL:HG13	1:A:32:ASP:N	2.10	0.66
1:D:147:LEU:HD22	1:D:151:GLY:HA3	1.77	0.66
1:C:106:ARG:HD3	1:C:186:ASN:OD1	1.96	0.66
1:A:299:VAL:HG22	2:A:401:FMN:H5'2	1.76	0.65
1:B:175:GLU:OE2	1:B:230:ARG:HB2	1.96	0.65
1:D:198:GLU:O	1:D:199:TYR:HB2	1.96	0.65
1:D:91:GLU:OE2	1:D:172:ARG:HD3	1.97	0.65
1:A:85:ARG:NH2	1:C:85:ARG:NH1	2.33	0.64
1:C:28:TYR:HB3	1:C:112:ARG:HG2	1.79	0.64
1:D:161:THR:HG22	1:D:165:ARG:NH1	2.11	0.64
1:A:25:MET:HA	2:A:401:FMN:N5	2.12	0.64
1:D:202:ASP:HB3	1:D:205:GLY:H	1.60	0.64
1:C:299:VAL:HG22	2:C:401:FMN:H5'2	1.81	0.63
1:D:101:LEU:HD11	1:D:174:ILE:HD11	1.80	0.63
1:C:112:ARG:HD3	1:C:114:TRP:CH2	2.34	0.62
1:D:120:ILE:CG2	1:D:125:GLY:O	2.48	0.61
1:D:2:GLN:NE2	1:D:3:PRO:HD2	2.17	0.60
1:C:299:VAL:HG22	2:C:401:FMN:C5'	2.32	0.60
1:D:137:PHE:CE1	1:D:270:PRO:HD3	2.36	0.59
1:D:31:VAL:HG23	1:D:32:ASP:N	2.18	0.59
1:D:25:MET:HA	2:D:401:FMN:N5	2.19	0.58
1:A:31:VAL:CG1	1:A:32:ASP:H	2.15	0.58
1:A:298:ALA:HA	4:A:536:HOH:O	2.02	0.58
1:B:197:ASP:HB3	1:B:199:TYR:H	1.70	0.57
1:B:58:ALA:HB1	1:B:102:ALA:CB	2.34	0.57
1:D:53:LEU:HD23	1:D:319:LEU:HD21	1.86	0.57
1:C:67:ILE:HG13	1:C:104:ALA:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:GLU:H	1:B:306:GLU:CD	2.14	0.55
1:D:167:ARG:HA	1:D:167:ARG:NE	2.19	0.55
1:B:130:VAL:HB	1:B:147:LEU:HD21	1.87	0.55
1:D:157:ALA:O	1:D:161:THR:OG1	2.22	0.55
1:B:175:GLU:HA	1:B:228:ALA:O	2.06	0.55
1:C:202:ASP:OD2	1:C:204:ARG:HB3	2.06	0.55
1:C:218:ARG:NH1	1:C:223:SER:HA	2.22	0.54
1:D:195:ARG:NH2	1:D:197:ASP:OD2	2.41	0.54
1:D:306:GLU:H	1:D:306:GLU:CD	2.16	0.54
1:B:243:ILE:O	1:B:247:VAL:HG23	2.08	0.53
1:A:101:LEU:HD11	1:A:174:ILE:HD11	1.89	0.53
1:C:298:ALA:HA	4:C:502:HOH:O	2.09	0.53
1:D:161:THR:CG2	1:D:165:ARG:NH1	2.71	0.53
1:D:152:ILE:HD13	1:D:199:TYR:CE1	2.44	0.53
1:B:58:ALA:HB1	1:B:102:ALA:HB2	1.91	0.53
1:C:278:GLU:OE2	1:C:304:ARG:HD3	2.09	0.53
1:A:172:ARG:O	1:A:226:PRO:HD2	2.08	0.52
1:C:213:VAL:O	1:C:217:VAL:HG23	2.09	0.52
1:A:236:TRP:CD1	1:A:274:ILE:HG12	2.44	0.52
1:B:182:TYR:OH	3:B:402:MLI:H12	2.10	0.52
1:C:325:GLU:HB2	4:C:520:HOH:O	2.09	0.52
1:D:31:VAL:CG2	1:D:32:ASP:H	2.21	0.52
1:D:25:MET:HA	2:D:401:FMN:C5A	2.40	0.51
1:B:298:ALA:HA	4:B:529:HOH:O	2.10	0.51
1:A:19:ARG:HD2	1:A:318:ASP:O	2.12	0.50
1:B:57:GLU:O	1:B:59:THR:HG23	2.11	0.50
1:D:160:ALA:HB2	1:D:216:ALA:HB1	1.92	0.50
1:A:299:VAL:HG22	2:A:401:FMN:C5'	2.42	0.50
1:C:207:VAL:HG22	1:C:252:MET:HE1	1.94	0.50
1:A:31:VAL:CG1	1:A:32:ASP:N	2.76	0.49
1:B:143:THR:HG21	1:C:345:ALA:HB3	1.95	0.49
1:C:147:LEU:HD22	1:C:151:GLY:C	2.38	0.49
1:D:38:TRP:CZ3	1:D:41:MET:HE1	2.48	0.48
1:B:59:THR:HG22	1:B:72:LEU:HD23	1.95	0.48
1:D:31:VAL:CG2	1:D:34:PHE:HB2	2.42	0.47
1:C:226:PRO:HA	1:C:259:ASP:OD2	2.15	0.47
1:B:277:GLY:O	1:B:278:GLU:C	2.58	0.47
1:A:43:LEU:HD21	4:A:550:HOH:O	2.13	0.47
1:C:51:VAL:HG12	1:C:333:PRO:HG2	1.96	0.47
1:D:182:TYR:O	1:D:183:LEU:C	2.58	0.46
1:D:63:PRO:HD3	1:D:75:TRP:CG	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:VAL:HG12	1:A:34:PHE:HB2	1.97	0.46
1:D:161:THR:CG2	1:D:165:ARG:CZ	2.94	0.46
1:D:98:GLY:HA2	1:D:173:TRP:O	2.16	0.46
1:D:252:MET:HE3	1:D:256:GLN:OE1	2.15	0.46
1:D:113:PRO:O	3:D:403:MLI:H12	2.16	0.45
1:D:101:LEU:CD1	1:D:174:ILE:HD11	2.47	0.45
1:C:290:ARG:NH2	4:C:504:HOH:O	2.45	0.45
1:D:137:PHE:CZ	1:D:270:PRO:HD3	2.52	0.45
1:A:112:ARG:HD3	1:A:114:TRP:CZ2	2.51	0.45
1:A:51:VAL:HG12	1:A:333:PRO:HG2	1.98	0.45
1:A:232:SER:HA	1:A:265:SER:HA	1.99	0.45
1:B:21:GLY:HA3	1:B:321:LEU:CD2	2.47	0.45
1:D:159:PHE:CG	1:D:213:VAL:HG13	2.51	0.45
1:D:118:LYS:HD3	1:D:118:LYS:HA	1.66	0.45
1:C:175:GLU:HA	1:C:228:ALA:O	2.17	0.44
1:A:16:LEU:HD22	1:A:53:LEU:HB2	1.99	0.44
1:A:25:MET:HA	2:A:401:FMN:C5A	2.47	0.44
1:B:299:VAL:HG22	2:B:401:FMN:C5'	2.42	0.44
1:D:332:TRP:HB3	1:D:333:PRO:HD3	2.00	0.44
1:C:267:GLY:HA2	4:C:528:HOH:O	2.18	0.44
1:D:5:LEU:HB2	1:D:309:ASP:HA	1.99	0.44
1:B:254:ARG:HD2	1:B:292:ALA:O	2.18	0.43
1:B:282:VAL:N	1:B:283:PRO:CD	2.81	0.43
1:B:61:VAL:HG11	1:B:159:PHE:CD1	2.53	0.43
1:B:98:GLY:HA2	1:B:173:TRP:O	2.18	0.43
1:D:49:GLY:O	1:D:333:PRO:HB2	2.18	0.43
1:A:23:SER:OG	1:A:24:PRO:HD2	2.19	0.43
1:B:59:THR:CG2	1:B:72:LEU:HD23	2.49	0.43
1:B:299:VAL:CG2	2:B:401:FMN:H5'1	2.45	0.43
1:A:46:ARG:HD2	1:A:46:ARG:N	2.34	0.43
1:B:247:VAL:O	1:B:251[A]:ARG:HG3	2.19	0.43
1:B:25:MET:HA	2:B:401:FMN:N5	2.33	0.43
1:C:67:ILE:HB	1:C:71:ASP:OD2	2.19	0.43
1:D:9:LEU:HD12	4:D:514:HOH:O	2.20	0.42
1:B:260:LEU:C	1:B:260:LEU:HD23	2.44	0.42
1:D:26:CYS:HA	1:D:57:GLU:HB2	2.01	0.42
1:D:209:LEU:O	1:D:213:VAL:HG23	2.20	0.42
1:D:198:GLU:H	1:D:198:GLU:HG3	1.68	0.42
1:C:191:LEU:HD11	1:C:239:GLU:HG2	2.02	0.41
1:D:143:THR:HA	1:D:144:PRO:HD3	1.81	0.41
1:A:319:LEU:HD12	1:A:319:LEU:HA	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:31:VAL:CG2	1:D:32:ASP:N	2.82	0.41
1:D:135:GLU:HA	1:D:136:PRO:HD3	1.91	0.41
1:C:260:LEU:HD23	1:C:260:LEU:C	2.46	0.41
1:A:128:PRO:HB2	1:A:145:ILE:HD11	2.01	0.41
1:A:225:LEU:HA	1:A:226:PRO:HD3	1.95	0.41
1:C:72:LEU:HD23	1:C:72:LEU:HA	1.90	0.41
1:C:202:ASP:OD1	1:C:203:LEU:N	2.54	0.41
1:C:112:ARG:CD	1:C:114:TRP:CZ2	2.99	0.40
1:A:332:TRP:N	1:A:333:PRO:CD	2.85	0.40
1:C:297:ALA:HA	1:C:319:LEU:O	2.22	0.40
1:A:108:ALA:HB3	1:A:129:VAL:HG12	2.02	0.40
1:B:26:CYS:HA	1:B:57:GLU:HB2	2.03	0.40
1:B:302:ILE:O	1:B:322:LEU:HD22	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	355/354 (100%)	341 (96%)	13 (4%)	1 (0%)	36	50
1	B	354/354 (100%)	340 (96%)	13 (4%)	1 (0%)	36	50
1	C	356/354 (101%)	337 (95%)	18 (5%)	1 (0%)	36	50
1	D	352/354 (99%)	327 (93%)	24 (7%)	1 (0%)	36	50
All	All	1417/1416 (100%)	1345 (95%)	68 (5%)	4 (0%)	36	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	57	GLU
1	D	199	TYR

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Mol	Chain	Res	Type
1	B	74	ILE
1	A	275	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	271/268 (101%)	263 (97%)	8 (3%)	36	58
1	B	270/268 (101%)	260 (96%)	10 (4%)	30	51
1	C	272/268 (102%)	261 (96%)	11 (4%)	28	47
1	D	268/268 (100%)	257 (96%)	11 (4%)	27	46
All	All	1081/1072 (101%)	1041 (96%)	40 (4%)	31	51

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LEU
1	A	41	MET
1	A	51	VAL
1	A	145	ILE
1	A	196	ASN
1	A	198	GLU
1	A	272	ILE
1	A	273	THR
1	B	1	MET
1	B	41	MET
1	B	54	ILE
1	B	84	SER
1	B	154	ARG
1	B	155	VAL
1	B	161	THR
1	B	167	ARG
1	B	188	LEU
1	B	237	THR
1	C	1	MET

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Mol	Chain	Res	Type
1	C	41	MET
1	C	51	VAL
1	C	154	ARG
1	C	191	LEU
1	C	208[A]	ARG
1	C	208[B]	ARG
1	C	237	THR
1	C	239	GLU
1	C	273	THR
1	C	338	ARG
1	D	68	SER
1	D	71	ASP
1	D	76	SER
1	D	135	GLU
1	D	155	VAL
1	D	167	ARG
1	D	197	ASP
1	D	198	GLU
1	D	204	ARG
1	D	290	ARG
1	D	301	LEU

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	2	GLN
1	B	177	HIS
1	B	281	GLN
1	B	307	HIS
1	D	2	GLN
1	D	124	ASN
1	D	193	ASN
1	D	281	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 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	FMN	C	401	-	33,33,33	0.83	1 (3%)	48,50,50	1.32	6 (12%)
3	MLI	A	402	-	6,6,6	1.55	0	7,7,7	0.95	0
3	MLI	D	402	-	6,6,6	1.47	1 (16%)	7,7,7	0.99	0
2	FMN	D	401	-	33,33,33	0.70	0	48,50,50	0.88	1 (2%)
3	MLI	C	402	-	6,6,6	1.37	0	7,7,7	1.10	0
2	FMN	B	401	-	33,33,33	0.93	1 (3%)	48,50,50	1.30	7 (14%)
3	MLI	D	403	-	6,6,6	1.31	0	7,7,7	0.94	0
3	MLI	B	402	-	6,6,6	1.54	2 (33%)	7,7,7	0.90	0
2	FMN	A	401	-	33,33,33	0.77	0	48,50,50	1.23	6 (12%)

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	FMN	C	401	-	-	5/18/18/18	0/3/3/3
3	MLI	A	402	-	-	0/4/4/4	-
3	MLI	D	402	-	-	0/4/4/4	-
2	FMN	D	401	-	-	1/18/18/18	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MLI	C	402	-	-	2/4/4/4	-
2	FMN	B	401	-	-	7/18/18/18	0/3/3/3
3	MLI	D	403	-	-	0/4/4/4	-
3	MLI	B	402	-	-	0/4/4/4	-
2	FMN	A	401	-	-	6/18/18/18	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	FMN	C1'-C2'	2.47	1.56	1.52
3	D	402	MLI	O7-C2	-2.29	1.23	1.30
3	B	402	MLI	O7-C2	-2.16	1.23	1.30
3	B	402	MLI	O9-C3	-2.07	1.24	1.30
2	C	401	FMN	C2-N1	-2.05	1.32	1.36

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	FMN	C5'-C4'-C3'	-4.54	103.65	112.22
2	B	401	FMN	O2P-P-O5'	-4.41	95.18	106.67
2	C	401	FMN	C4'-C3'-C2'	3.69	119.72	113.57
2	A	401	FMN	C5'-C4'-C3'	-3.68	105.27	112.22
2	B	401	FMN	C5'-C4'-C3'	-3.42	105.77	112.22
2	C	401	FMN	O2P-P-O5'	-3.35	97.95	106.67
2	A	401	FMN	O3P-P-O5'	-3.27	98.15	106.67
2	B	401	FMN	O3P-P-O5'	2.91	114.25	106.67
2	A	401	FMN	C1'-C2'-C3'	2.90	117.53	109.66
2	B	401	FMN	O3'-C3'-C2'	2.61	114.87	108.93
2	B	401	FMN	C1'-C2'-C3'	2.32	115.95	109.66
2	A	401	FMN	C4'-C3'-C2'	2.21	117.24	113.57
2	D	401	FMN	C4-N3-C2	-2.19	121.75	125.64
2	B	401	FMN	C4-N3-C2	-2.17	121.79	125.64
2	C	401	FMN	O2'-C2'-C3'	-2.13	104.27	109.25
2	C	401	FMN	C1'-C2'-C3'	2.11	115.39	109.66
2	A	401	FMN	C4A-C4-N3	2.07	118.52	113.25
2	C	401	FMN	C4-N3-C2	-2.04	122.01	125.64
2	B	401	FMN	C4A-C4-N3	2.02	118.40	113.25
2	A	401	FMN	C4-N3-C2	-2.01	122.07	125.64

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	FMN	C2'-C3'-C4'-O4'
2	A	401	FMN	C2'-C3'-C4'-C5'
2	A	401	FMN	O3'-C3'-C4'-O4'
2	A	401	FMN	O3'-C3'-C4'-C5'
2	A	401	FMN	O4'-C4'-C5'-O5'
2	B	401	FMN	C3'-C4'-C5'-O5'
2	B	401	FMN	O4'-C4'-C5'-O5'
2	B	401	FMN	C5'-O5'-P-O1P
2	C	401	FMN	O4'-C4'-C5'-O5'
2	C	401	FMN	C2'-C3'-C4'-O4'
2	C	401	FMN	C2'-C3'-C4'-C5'
2	B	401	FMN	C2'-C3'-C4'-O4'
2	C	401	FMN	C3'-C4'-C5'-O5'
2	B	401	FMN	C4'-C5'-O5'-P
2	B	401	FMN	C2'-C3'-C4'-C5'
2	C	401	FMN	C4'-C5'-O5'-P
2	D	401	FMN	C4'-C5'-O5'-P
2	B	401	FMN	C5'-O5'-P-O2P
2	A	401	FMN	C4'-C5'-O5'-P
3	C	402	MLI	C2-C1-C3-O8
3	C	402	MLI	C2-C1-C3-O9

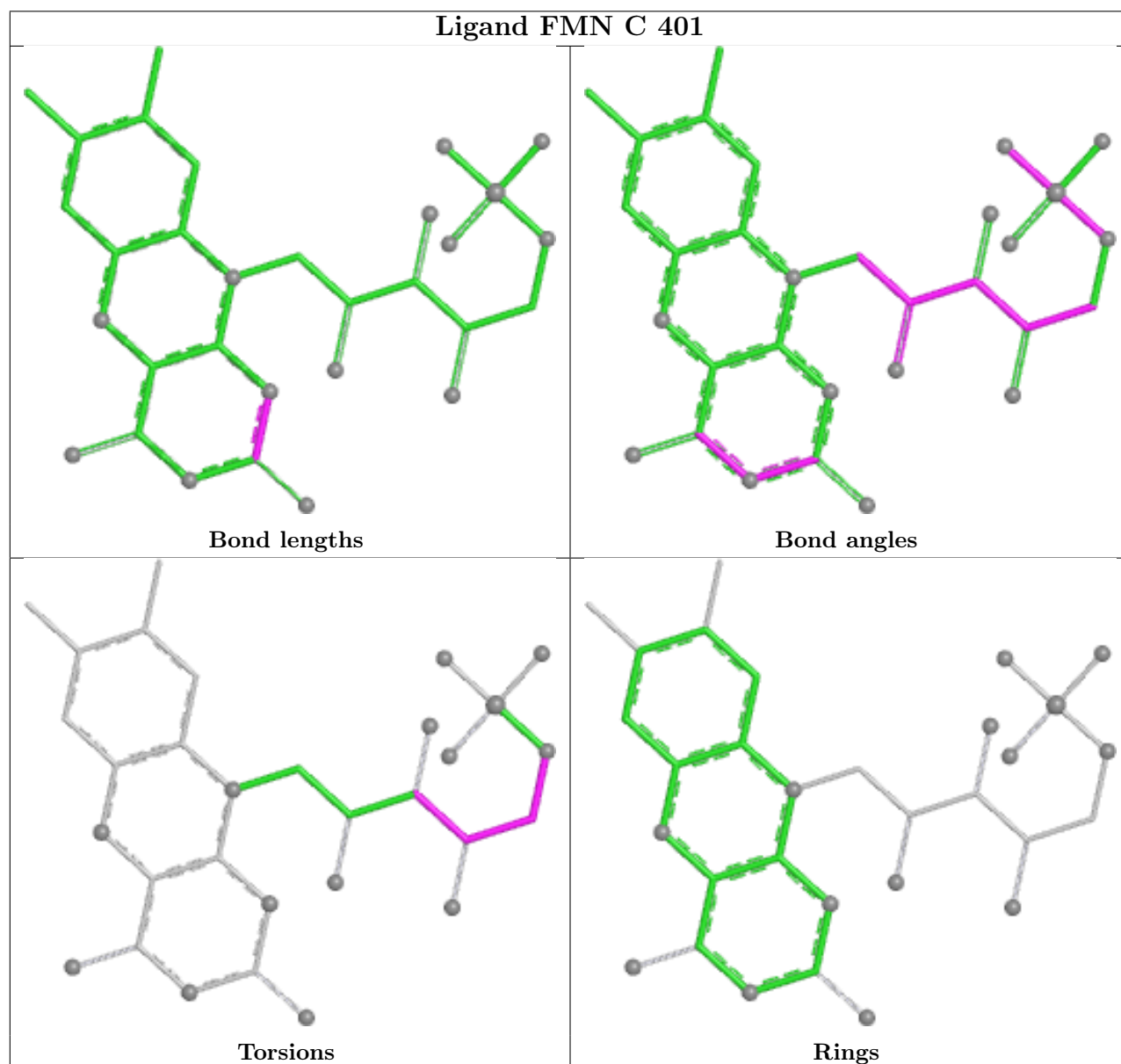
There are no ring outliers.

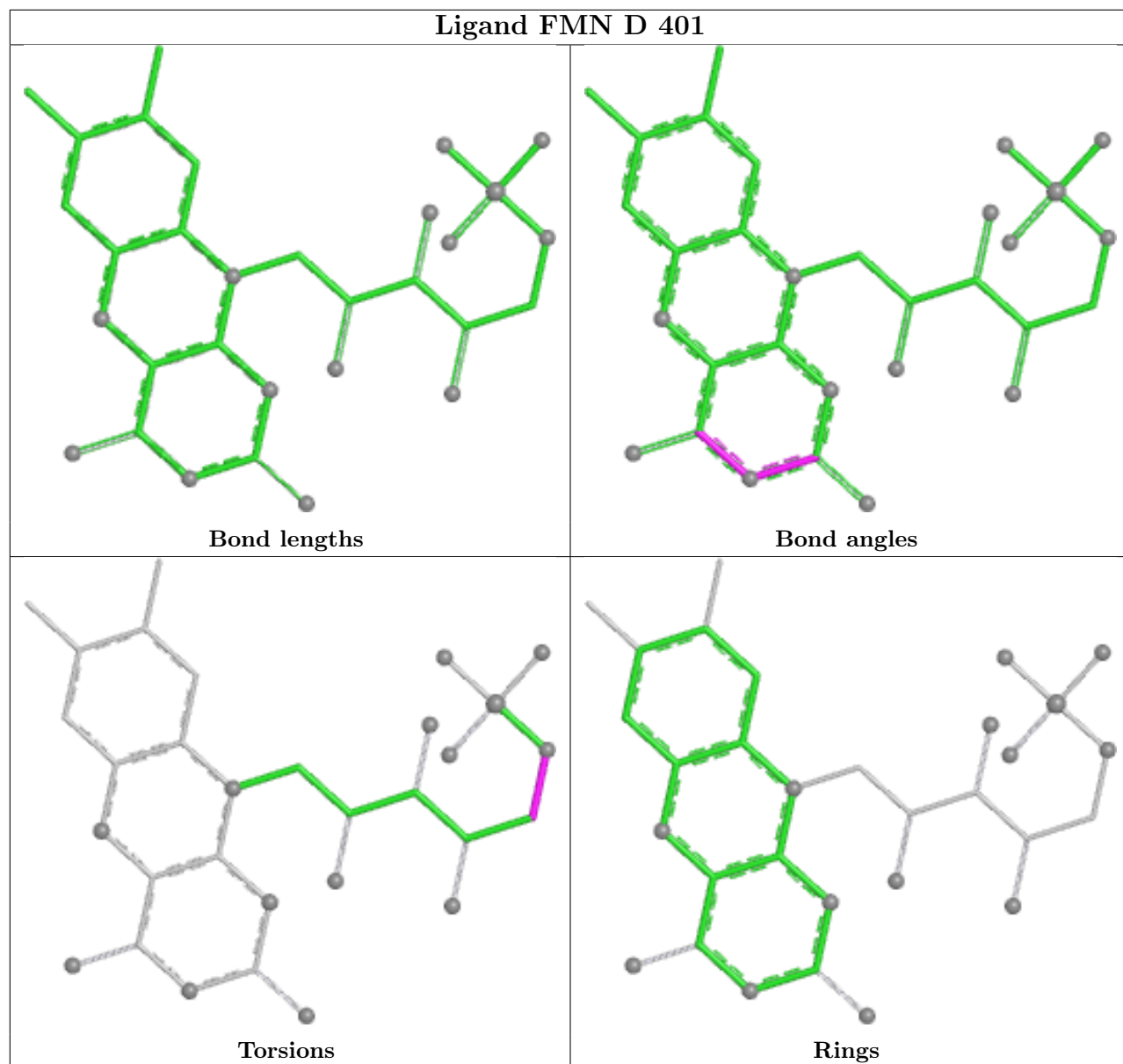
6 monomers are involved in 14 short contacts:

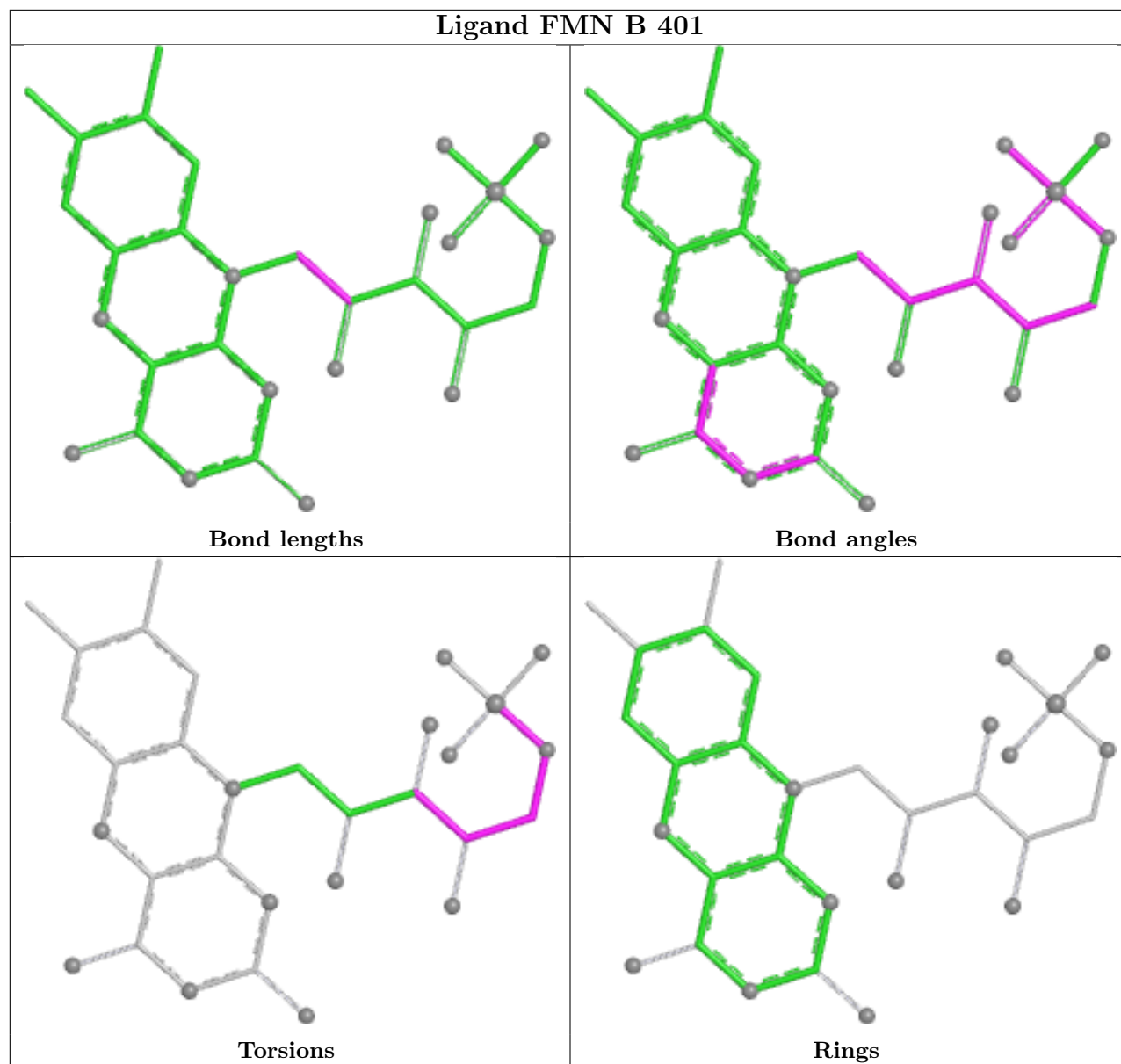
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	FMN	2	0
2	D	401	FMN	2	0
2	B	401	FMN	4	0
3	D	403	MLI	1	0
3	B	402	MLI	1	0
2	A	401	FMN	4	0

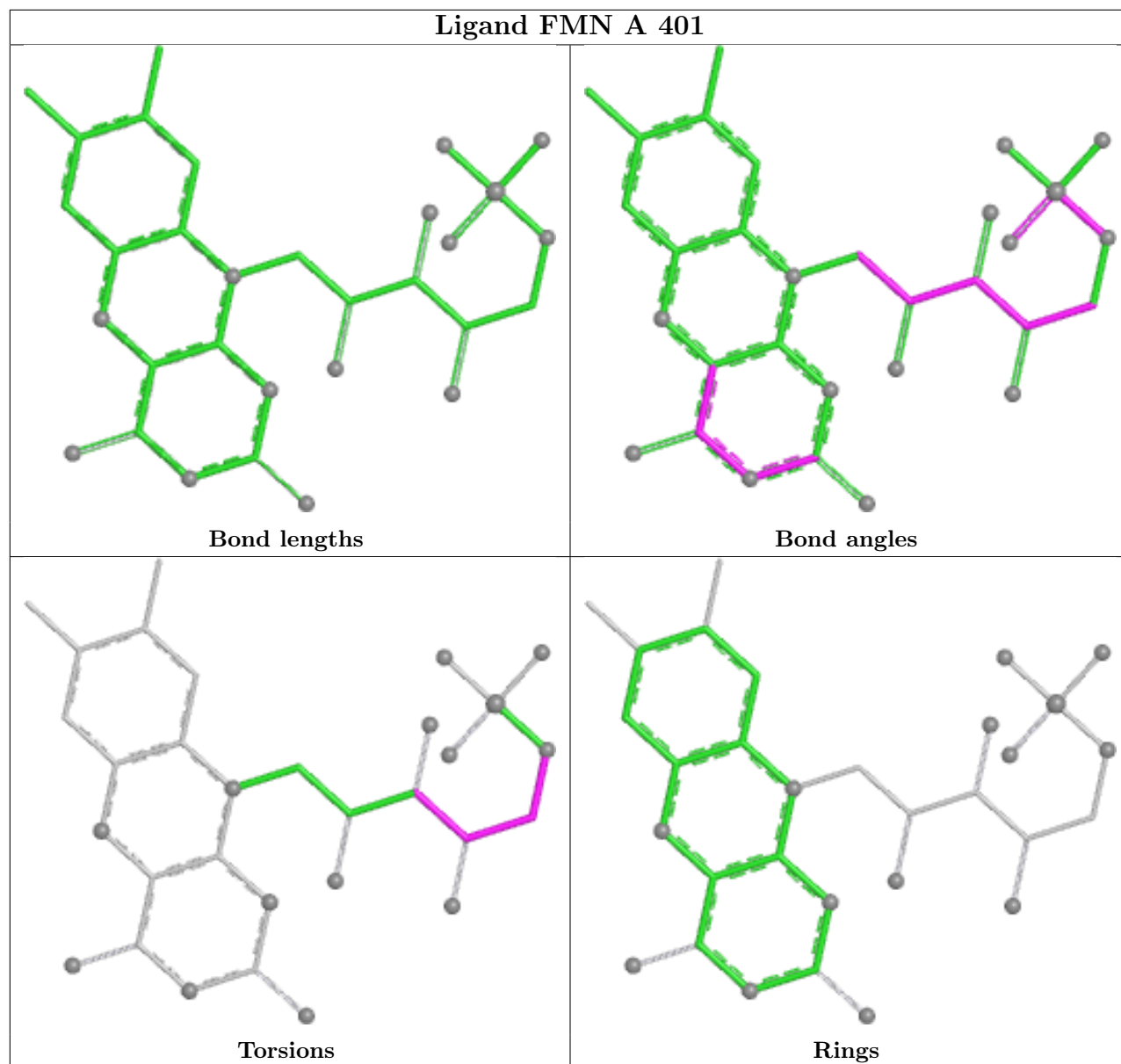
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	354/354 (100%)	-0.21	7 (1%) 65 60	11, 25, 52, 85	3 (0%)
1	B	354/354 (100%)	-0.08	3 (0%) 82 80	17, 32, 49, 97	2 (0%)
1	C	354/354 (100%)	0.02	8 (2%) 61 57	11, 30, 65, 90	4 (1%)
1	D	354/354 (100%)	0.42	37 (10%) 11 8	17, 34, 79, 108	0
All	All	1416/1416 (100%)	0.04	55 (3%) 43 39	11, 30, 64, 108	9 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	122	PRO	6.3
1	D	125	GLY	6.0
1	D	192	GLY	5.4
1	D	121	ALA	5.2
1	D	1	MET	5.1
1	D	123	ALA	4.6
1	B	1	MET	4.4
1	D	128	PRO	4.3
1	D	191	LEU	4.3
1	B	293[A]	ASN	4.1
1	D	124	ASN	4.1
1	A	193	ASN	3.9
1	D	193	ASN	3.8
1	D	134	ALA	3.8
1	C	1	MET	3.7
1	A	192	GLY	3.7
1	C	2	GLN	3.5
1	D	142	PRO	3.4
1	D	194	ASP	3.2
1	B	202	ASP	3.2
1	D	2	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	194	ASP	3.1
1	D	31	VAL	3.1
1	D	145	ILE	3.0
1	D	141	TYR	3.0
1	D	129	VAL	2.9
1	D	150	ALA	2.9
1	D	65	GLY	2.9
1	D	136	PRO	2.8
1	C	192	GLY	2.8
1	D	108	ALA	2.8
1	D	126	GLY	2.7
1	D	127	TRP	2.7
1	C	237	THR	2.7
1	D	137	PHE	2.6
1	D	197	ASP	2.6
1	A	273	THR	2.6
1	D	138	ALA	2.6
1	D	144	PRO	2.5
1	A	31	VAL	2.4
1	D	70	PHE	2.3
1	A	196	ASN	2.3
1	D	139	PRO	2.3
1	D	32	ASP	2.3
1	D	271	GLY	2.2
1	D	71	ASP	2.2
1	D	196	ASN	2.2
1	D	130	VAL	2.2
1	D	76	SER	2.1
1	C	139	PRO	2.1
1	D	120	ILE	2.1
1	C	105	GLY	2.0
1	A	251[A]	ARG	2.0
1	C	154	ARG	2.0
1	C	191	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

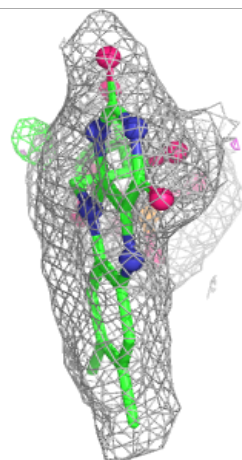
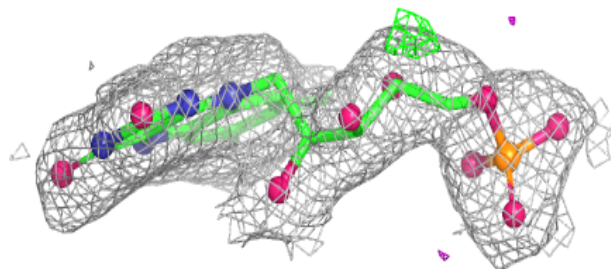
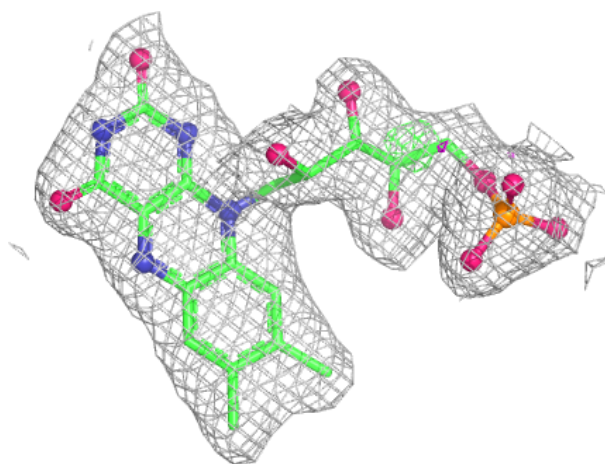
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MLI	D	403	7/7	0.85	0.19	60,62,66,74	0
3	MLI	C	402	7/7	0.86	0.17	46,49,60,66	0
3	MLI	D	402	7/7	0.93	0.09	34,36,41,47	0
3	MLI	A	402	7/7	0.95	0.08	27,31,36,39	0
2	FMN	B	401	31/31	0.96	0.07	16,20,22,24	0
2	FMN	C	401	31/31	0.96	0.07	19,22,26,28	0
2	FMN	D	401	31/31	0.97	0.06	17,20,23,24	0
2	FMN	A	401	31/31	0.97	0.06	19,23,26,27	0
3	MLI	B	402	7/7	0.97	0.05	31,31,33,37	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.

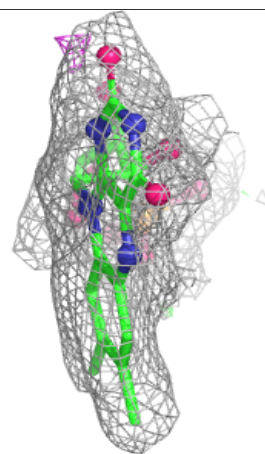
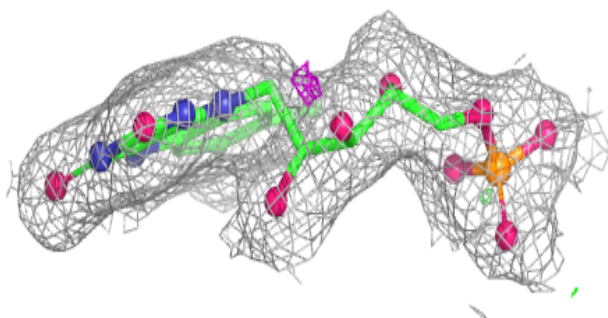
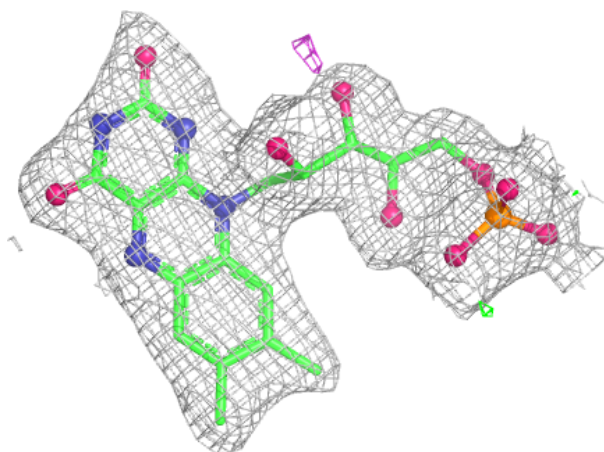
Electron density around FMN B 401:

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



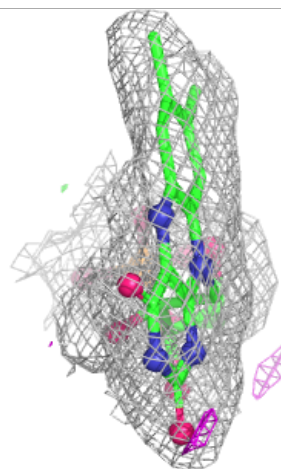
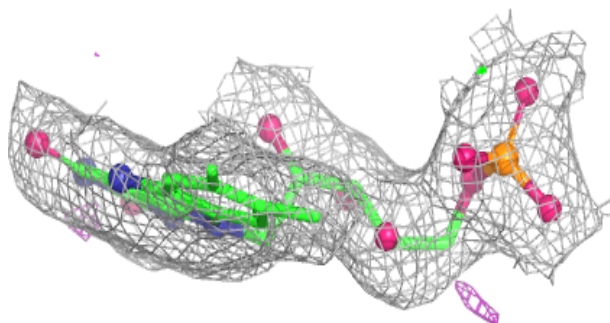
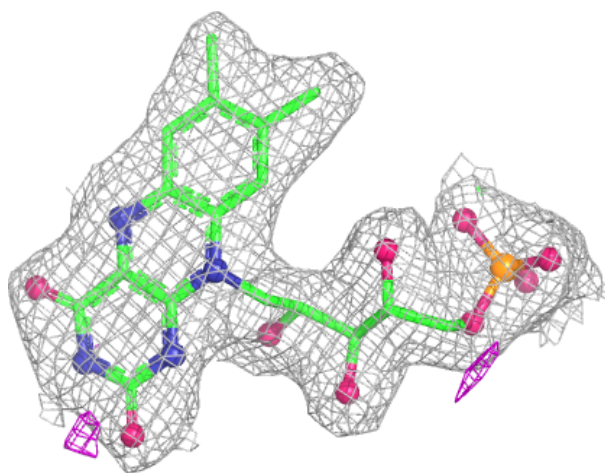
Electron density around FMN C 401:

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



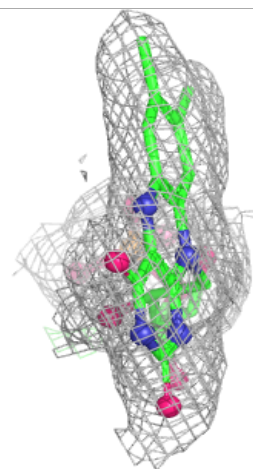
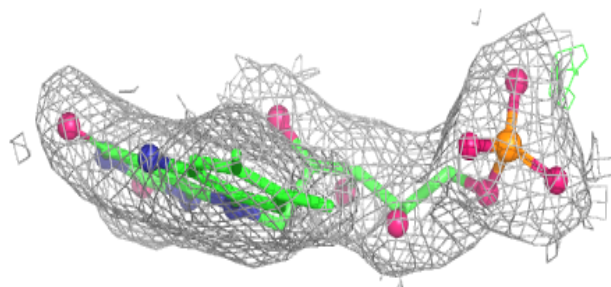
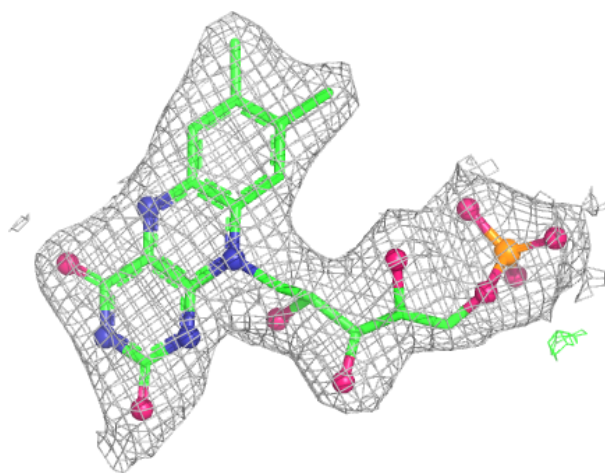
Electron density around FMN D 401:

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



Electron density around FMN 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.