



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:36 PM UTC

PDB ID : 6MVH / pdb_00006mvh
Title : Crystal structure of FMN-binding beta-glucuronidase from Roseburia hominis
Authors : Pellock, S.J.; Redinbo, M.R.
Deposited on : 2018-10-25
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

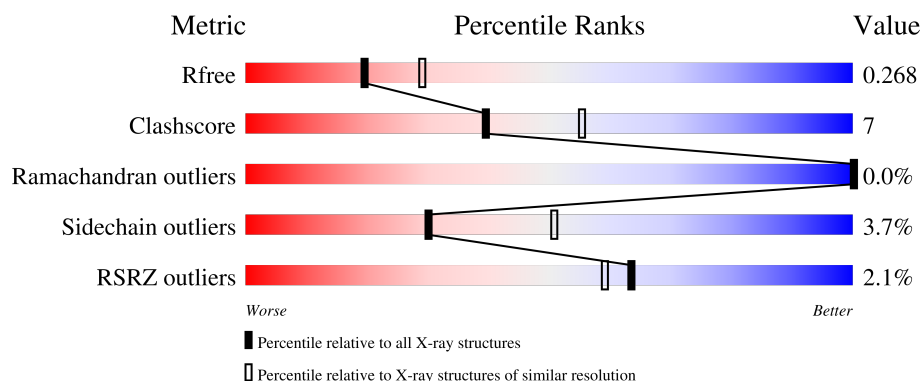
1 Overall quality at a glance

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	780	69% 12% • 18%
1	B	780	67% 13% • 18%
1	C	780	68% 13% • 18%
1	D	780	5% 29% 10% • 61%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18964 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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	0	0
			5175	3290	865	1000	20			
1	B	638	Total	C	N	O	S	0	0	0
			5167	3286	864	997	20			
1	C	638	Total	C	N	O	S	0	1	0
			5174	3291	866	997	20			
1	D	305	Total	C	N	O	S	0	0	0
			2494	1573	429	480	12			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	HIS	-	expression tag	UNP A0A174HGC0
A	-22	HIS	-	expression tag	UNP A0A174HGC0
A	-21	HIS	-	expression tag	UNP A0A174HGC0
A	-20	HIS	-	expression tag	UNP A0A174HGC0
A	-19	HIS	-	expression tag	UNP A0A174HGC0
A	-18	HIS	-	expression tag	UNP A0A174HGC0
A	-17	SER	-	expression tag	UNP A0A174HGC0
A	-16	SER	-	expression tag	UNP A0A174HGC0
A	-15	GLY	-	expression tag	UNP A0A174HGC0
A	-14	VAL	-	expression tag	UNP A0A174HGC0
A	-13	ASP	-	expression tag	UNP A0A174HGC0
A	-12	LEU	-	expression tag	UNP A0A174HGC0
A	-11	GLY	-	expression tag	UNP A0A174HGC0
A	-10	THR	-	expression tag	UNP A0A174HGC0
A	-9	GLU	-	expression tag	UNP A0A174HGC0
A	-8	ASN	-	expression tag	UNP A0A174HGC0
A	-7	LEU	-	expression tag	UNP A0A174HGC0
A	-6	TYR	-	expression tag	UNP A0A174HGC0
A	-5	PHE	-	expression tag	UNP A0A174HGC0
A	-4	GLN	-	expression tag	UNP A0A174HGC0
A	-3	SER	-	expression tag	UNP A0A174HGC0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ASN	-	expression tag	UNP A0A174HGC0
A	-1	ALA	-	expression tag	UNP A0A174HGC0
B	-23	HIS	-	expression tag	UNP A0A174HGC0
B	-22	HIS	-	expression tag	UNP A0A174HGC0
B	-21	HIS	-	expression tag	UNP A0A174HGC0
B	-20	HIS	-	expression tag	UNP A0A174HGC0
B	-19	HIS	-	expression tag	UNP A0A174HGC0
B	-18	HIS	-	expression tag	UNP A0A174HGC0
B	-17	SER	-	expression tag	UNP A0A174HGC0
B	-16	SER	-	expression tag	UNP A0A174HGC0
B	-15	GLY	-	expression tag	UNP A0A174HGC0
B	-14	VAL	-	expression tag	UNP A0A174HGC0
B	-13	ASP	-	expression tag	UNP A0A174HGC0
B	-12	LEU	-	expression tag	UNP A0A174HGC0
B	-11	GLY	-	expression tag	UNP A0A174HGC0
B	-10	THR	-	expression tag	UNP A0A174HGC0
B	-9	GLU	-	expression tag	UNP A0A174HGC0
B	-8	ASN	-	expression tag	UNP A0A174HGC0
B	-7	LEU	-	expression tag	UNP A0A174HGC0
B	-6	TYR	-	expression tag	UNP A0A174HGC0
B	-5	PHE	-	expression tag	UNP A0A174HGC0
B	-4	GLN	-	expression tag	UNP A0A174HGC0
B	-3	SER	-	expression tag	UNP A0A174HGC0
B	-2	ASN	-	expression tag	UNP A0A174HGC0
B	-1	ALA	-	expression tag	UNP A0A174HGC0
C	-23	HIS	-	expression tag	UNP A0A174HGC0
C	-22	HIS	-	expression tag	UNP A0A174HGC0
C	-21	HIS	-	expression tag	UNP A0A174HGC0
C	-20	HIS	-	expression tag	UNP A0A174HGC0
C	-19	HIS	-	expression tag	UNP A0A174HGC0
C	-18	HIS	-	expression tag	UNP A0A174HGC0
C	-17	SER	-	expression tag	UNP A0A174HGC0
C	-16	SER	-	expression tag	UNP A0A174HGC0
C	-15	GLY	-	expression tag	UNP A0A174HGC0
C	-14	VAL	-	expression tag	UNP A0A174HGC0
C	-13	ASP	-	expression tag	UNP A0A174HGC0
C	-12	LEU	-	expression tag	UNP A0A174HGC0
C	-11	GLY	-	expression tag	UNP A0A174HGC0
C	-10	THR	-	expression tag	UNP A0A174HGC0
C	-9	GLU	-	expression tag	UNP A0A174HGC0
C	-8	ASN	-	expression tag	UNP A0A174HGC0
C	-7	LEU	-	expression tag	UNP A0A174HGC0

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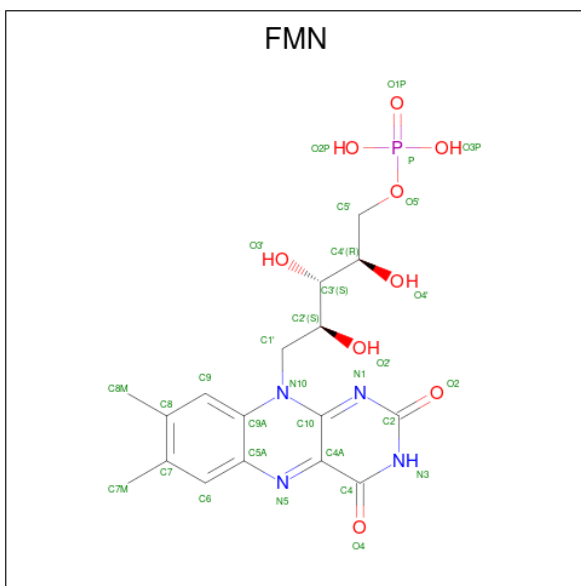
Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	TYR	-	expression tag	UNP A0A174HGC0
C	-5	PHE	-	expression tag	UNP A0A174HGC0
C	-4	GLN	-	expression tag	UNP A0A174HGC0
C	-3	SER	-	expression tag	UNP A0A174HGC0
C	-2	ASN	-	expression tag	UNP A0A174HGC0
C	-1	ALA	-	expression tag	UNP A0A174HGC0
D	-23	HIS	-	expression tag	UNP A0A174HGC0
D	-22	HIS	-	expression tag	UNP A0A174HGC0
D	-21	HIS	-	expression tag	UNP A0A174HGC0
D	-20	HIS	-	expression tag	UNP A0A174HGC0
D	-19	HIS	-	expression tag	UNP A0A174HGC0
D	-18	HIS	-	expression tag	UNP A0A174HGC0
D	-17	SER	-	expression tag	UNP A0A174HGC0
D	-16	SER	-	expression tag	UNP A0A174HGC0
D	-15	GLY	-	expression tag	UNP A0A174HGC0
D	-14	VAL	-	expression tag	UNP A0A174HGC0
D	-13	ASP	-	expression tag	UNP A0A174HGC0
D	-12	LEU	-	expression tag	UNP A0A174HGC0
D	-11	GLY	-	expression tag	UNP A0A174HGC0
D	-10	THR	-	expression tag	UNP A0A174HGC0
D	-9	GLU	-	expression tag	UNP A0A174HGC0
D	-8	ASN	-	expression tag	UNP A0A174HGC0
D	-7	LEU	-	expression tag	UNP A0A174HGC0
D	-6	TYR	-	expression tag	UNP A0A174HGC0
D	-5	PHE	-	expression tag	UNP A0A174HGC0
D	-4	GLN	-	expression tag	UNP A0A174HGC0
D	-3	SER	-	expression tag	UNP A0A174HGC0
D	-2	ASN	-	expression tag	UNP A0A174HGC0
D	-1	ALA	-	expression tag	UNP A0A174HGC0

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0

- Molecule 3 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: C₁₇H₂₁N₄O₉P)

(labeled as "Ligand of Interest" by depositor).



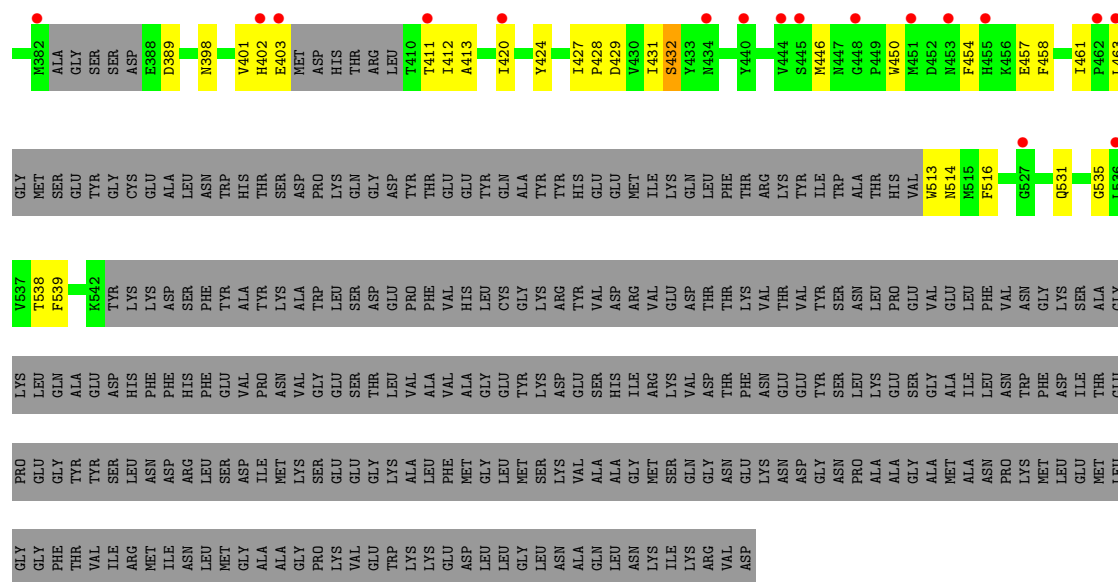
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 31	C 17	N 4	O 9	P 1	0	0
3	B	1	Total 31	C 17	N 4	O 9	P 1	0	0
3	C	1	Total 31	C 17	N 4	O 9	P 1	0	0
3	D	1	Total 31	C 17	N 4	O 9	P 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	296	Total O 296 296	0	0
4	B	224	Total O 224 224	0	0
4	C	234	Total O 234 234	0	0
4	D	72	Total O 72 72	0	0

- Molecule 1: Beta-galactosidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.42Å 137.41Å 108.83Å 90.00° 91.88° 90.00°	Depositor
Resolution (Å)	29.27 – 2.40 29.27 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.3 (29.27-2.40) 98.3 (29.27-2.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.77 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.209 , 0.268 0.209 , 0.268	Depositor DCC
R_{free} test set	1996 reflections (1.84%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	18964	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: CA, 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	0.42	0/5317	0.63	4/7226 (0.1%)
1	B	0.36	0/5309	0.57	0/7215
1	C	0.37	0/5320	0.59	1/7230 (0.0%)
1	D	0.34	0/2560	0.56	0/3468
All	All	0.38	0/18506	0.59	5/25139 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	199	ASP	N-CA-CB	-7.65	99.90	110.38
1	A	198	LYS	CA-C-N	7.53	135.22	121.06
1	A	198	LYS	C-N-CA	7.53	135.22	121.06
1	C	139	GLY	N-CA-C	6.82	119.47	111.63
1	A	199	ASP	CA-CB-CG	6.80	119.40	112.60

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	5175	0	4870	64	0
1	B	5167	0	4866	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	5174	0	4873	63	0
1	D	2494	0	2279	51	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	31	0	19	3	0
3	B	31	0	19	0	0
3	C	31	0	19	2	0
3	D	31	0	19	2	0
4	A	296	0	0	8	0
4	B	224	0	0	6	0
4	C	234	0	0	8	0
4	D	72	0	0	3	0
All	All	18964	0	16964	234	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 7.

All (234) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:157:ASP:OD2	4:C:901:HOH:O	1.92	0.88
1:A:196:THR:HG22	1:A:207:LYS:HG3	1.65	0.79
1:D:162:PRO:HD3	1:D:364:GLN:HA	1.65	0.77
1:C:61:LYS:HA	1:C:64:LEU:HD12	1.71	0.72
1:B:349:ASN:ND2	4:B:901:HOH:O	2.12	0.70
1:B:13:THR:HB	1:B:15:GLU:H	1.57	0.70
1:A:16:ALA:HB3	1:A:54:TYR:HD2	1.56	0.70
1:A:120:GLU:OE1	4:A:901:HOH:O	2.09	0.70
1:C:479:ASP:OD2	1:C:481:LYS:HE2	1.92	0.70
3:D:802:FMN:H2'	3:D:802:FMN:H9	1.74	0.69
1:C:213:GLY:O	4:C:902:HOH:O	2.11	0.68
1:C:304:ASP:OD1	1:C:332:ARG:NH2	2.27	0.68
1:B:227:TRP:CZ2	1:B:333:GLY:HA2	2.29	0.67
1:B:165:LYS:NZ	4:B:904:HOH:O	2.27	0.67
1:A:59:LEU:HD21	1:A:64:LEU:HD11	1.77	0.66
1:B:68:ASP:OD2	1:B:151:LYS:HG2	1.96	0.66
1:C:388:GLU:HA	1:C:391:LEU:HD12	1.79	0.65
1:B:263:ILE:HD12	1:B:409:LEU:HD12	1.80	0.64
1:D:458:PHE:HB3	1:D:461:ILE:HD12	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:LYS:HE3	1:B:209:GLU:OE2	1.99	0.63
1:A:16:ALA:HB3	1:A:54:TYR:CD2	2.33	0.63
1:D:420:ILE:HG22	1:D:458:PHE:HE2	1.64	0.62
1:A:8:THR:O	1:A:57:LYS:HE2	1.98	0.62
1:B:88:LYS:HE2	1:B:107:GLU:OE2	1.99	0.61
1:C:316:LEU:HB3	1:C:321:HIS:CD2	2.36	0.61
1:C:356:SER:OG	4:C:903:HOH:O	2.14	0.60
1:C:439:TRP:CZ2	1:C:471:GLU:HG3	2.37	0.60
1:D:73:GLU:HG3	1:D:101:ARG:HG2	1.82	0.60
1:D:324:TYR:CE2	1:D:328:LEU:HD11	2.35	0.60
1:D:356:SER:OG	4:D:902:HOH:O	2.16	0.60
1:C:13:THR:OG1	1:C:15:GLU:HG3	2.02	0.60
1:D:420:ILE:HG22	1:D:458:PHE:CE2	2.37	0.59
1:C:6:PHE:CZ	1:C:57:LYS:HD3	2.37	0.59
1:A:13:THR:HG21	1:A:16:ALA:HB2	1.84	0.59
1:A:563:LEU:HD13	1:A:580:VAL:HG22	1.84	0.59
1:D:317:ALA:HB1	1:D:318:HIS:CG	2.38	0.58
1:D:325:PHE:HA	1:D:328:LEU:HD12	1.85	0.58
3:D:802:FMN:O4'	3:D:802:FMN:O2'	2.18	0.58
1:A:316:LEU:HB3	1:A:321:HIS:CD2	2.38	0.58
1:D:454:PHE:CD2	1:D:463:LEU:HD22	2.39	0.58
1:A:591:VAL:O	1:A:594:LYS:HE3	2.04	0.57
1:D:360:GLU:O	1:D:364:GLN:HG2	2.04	0.57
1:C:540:ASP:OD2	1:C:542:LYS:HD2	2.04	0.56
1:D:142:ARG:HD3	4:D:937:HOH:O	2.06	0.56
1:C:175:ALA:HB2	1:C:224:VAL:HG11	1.88	0.56
1:D:85:VAL:HG21	1:D:104:ILE:HD11	1.87	0.56
1:D:318:HIS:HA	1:D:339:GLU:OE2	2.06	0.56
1:D:398:ASN:ND2	1:D:427:ILE:O	2.39	0.55
1:C:411:THR:HA	1:C:430:VAL:O	2.06	0.55
1:B:126:ARG:NH2	4:B:916:HOH:O	2.38	0.55
1:D:85:VAL:HG21	1:D:104:ILE:CD1	2.37	0.55
1:B:580:VAL:HG11	1:B:587:VAL:HG11	1.89	0.55
1:B:59:LEU:HD21	1:B:64:LEU:HD21	1.89	0.54
1:C:214:GLU:OE2	1:C:216:LYS:HE2	2.07	0.54
1:A:304:ASP:OD1	1:A:332:ARG:NH2	2.34	0.54
1:C:194:VAL:HG22	1:C:209:GLU:HG2	1.89	0.54
1:B:128:TYR:HB3	1:B:341:PRO:O	2.08	0.54
1:C:412:ILE:HG12	1:C:428:PRO:HG3	1.89	0.54
1:B:94:ASP:OD1	4:B:902:HOH:O	2.19	0.53
1:B:197:VAL:HG23	1:B:219:LEU:HD22	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:588:GLU:HB2	1:A:598:LYS:HG2	1.91	0.53
1:B:58:GLN:HB2	1:B:60:LYS:HE2	1.90	0.52
1:D:317:ALA:HB1	1:D:318:HIS:ND1	2.24	0.52
1:A:203:LYS:HD3	1:A:204:GLU:O	2.10	0.52
1:D:320:GLN:NE2	1:D:364:GLN:OE1	2.35	0.52
1:C:426:GLN:HB3	1:D:403:GLU:OE1	2.09	0.52
1:A:330:ASP:CG	1:A:369:PRO:HD2	2.34	0.52
1:A:26:LYS:NZ	1:D:389:ASP:OD2	2.31	0.52
1:A:170:ILE:HD11	1:A:258:CYS:HB3	1.92	0.52
1:C:15:GLU:OE2	1:C:17:THR:HG23	2.10	0.52
1:D:398:ASN:HD21	1:D:428:PRO:HA	1.75	0.52
1:C:303:ILE:HD13	1:C:328:LEU:HB3	1.92	0.52
1:B:198:LYS:NZ	1:B:239:GLU:OE2	2.32	0.52
1:A:577:LYS:H	1:A:577:LYS:HD2	1.76	0.51
1:A:6:PHE:CZ	1:A:57:LYS:HD3	2.46	0.51
1:C:33:PRO:HB3	1:C:141:TYR:O	2.10	0.51
1:B:65:PRO:O	1:B:70:TYR:OH	2.21	0.51
1:B:17:THR:HG21	1:D:351:ARG:HG2	1.92	0.51
1:C:563:LEU:HD13	1:C:580:VAL:HG22	1.91	0.51
1:D:402:HIS:NE2	1:D:429:ASP:OD2	2.27	0.51
1:C:476:HIS:HA	1:C:488:GLU:OE2	2.11	0.50
1:B:175:ALA:HB3	1:B:221:ILE:HG23	1.93	0.50
1:A:37:ASN:HA	1:A:40:ASP:OD1	2.11	0.50
1:C:1:ARG:NH1	1:C:156:LEU:HD21	2.26	0.50
1:C:419:ASP:OD2	1:C:421:HIS:N	2.39	0.50
1:C:514:ASN:O	1:C:535:GLY:HA2	2.11	0.50
1:C:599:LEU:HD12	1:C:608:PHE:CE1	2.47	0.50
1:A:315:ARG:HH12	1:A:467:GLU:HG3	1.76	0.50
1:A:411:THR:HA	1:A:430:VAL:O	2.11	0.50
1:D:82:ASP:OD2	1:D:92:HIS:HB2	2.12	0.50
1:D:35:SER:O	1:D:38:GLU:HG3	2.12	0.49
1:D:398:ASN:HD22	1:D:427:ILE:HG22	1.77	0.49
1:B:62:SER:HB3	1:B:111:GLU:OE2	2.12	0.49
1:B:14:LYS:NZ	4:B:906:HOH:O	2.30	0.49
1:C:8:THR:O	1:C:57:LYS:HE2	2.13	0.49
1:D:316:LEU:HD12	1:D:326:TYR:HE1	1.78	0.49
1:B:60:LYS:NZ	1:B:112:GLU:OE1	2.42	0.49
1:A:266:GLU:OE1	1:A:455:HIS:NE2	2.33	0.49
1:A:439:TRP:CZ2	1:A:471:GLU:HG3	2.47	0.49
1:A:33:PRO:HB3	1:A:141:TYR:O	2.13	0.49
1:A:351:ARG:HH11	1:C:17:THR:HG22	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:SER:O	1:A:382:MET:HB2	2.13	0.48
1:B:561:VAL:HG23	1:B:582:SER:HB2	1.94	0.48
1:B:352:GLU:CD	1:B:352:GLU:H	2.21	0.48
1:A:21:LYS:HG3	3:C:802:FMN:O1P	2.13	0.48
1:A:182:VAL:HG21	1:A:240:VAL:HG11	1.95	0.48
1:A:616:SER:OG	4:A:902:HOH:O	2.19	0.48
1:C:482:GLN:NE2	4:C:909:HOH:O	2.33	0.48
1:B:191:GLN:NE2	1:B:244:SER:OG	2.44	0.48
1:D:514:ASN:O	1:D:535:GLY:HA2	2.14	0.48
1:A:239:GLU:HG3	4:A:1176:HOH:O	2.14	0.47
1:B:316:LEU:HB3	1:B:321:HIS:CD2	2.49	0.47
1:A:315:ARG:NH1	1:A:467:GLU:HG3	2.29	0.47
1:C:10:TRP:CE3	1:C:57:LYS:HB2	2.49	0.47
1:B:411:THR:HA	1:B:430:VAL:O	2.14	0.47
1:A:109:THR:O	4:A:904:HOH:O	2.20	0.47
1:A:583:ASN:O	4:A:903:HOH:O	2.20	0.47
1:C:197:VAL:HG23	1:C:219:LEU:HD22	1.95	0.47
1:A:318:HIS:HA	1:A:339:GLU:OE2	2.14	0.47
1:B:68:ASP:HB2	1:B:151:LYS:N	2.29	0.47
1:B:46:ASN:HB2	1:B:522:ALA:HA	1.97	0.47
1:A:430:VAL:HG12	1:A:462:PRO:HG2	1.98	0.46
1:A:587:VAL:HG22	1:A:622:ALA:HB2	1.96	0.46
1:C:372:VAL:HG23	1:C:373:VAL:HG13	1.97	0.46
1:A:158:TYR:OH	3:A:802:FMN:H5'2	2.15	0.46
1:C:74:LEU:HD22	1:C:100:TRP:CZ2	2.51	0.46
1:A:591:VAL:HG21	1:A:610:VAL:HG22	1.97	0.46
1:B:33:PRO:HB3	1:B:141:TYR:O	2.15	0.46
1:A:171:LYS:HD3	1:A:171:LYS:HA	1.62	0.46
1:D:33:PRO:HB3	1:D:141:TYR:O	2.15	0.45
1:A:552:LYS:HG2	1:A:560:PHE:CE1	2.52	0.45
1:C:351:ARG:HB2	4:C:965:HOH:O	2.16	0.45
1:A:417:MET:SD	1:A:417:MET:N	2.87	0.45
1:B:71:TYR:HB2	1:B:147:ILE:HB	1.97	0.45
1:C:1:ARG:HB2	1:C:250:ASP:OD1	2.16	0.45
1:A:9:LYS:NZ	4:A:935:HOH:O	2.49	0.45
1:A:166:VAL:HG11	1:A:256:PHE:CG	2.52	0.45
1:C:38:GLU:HG3	1:C:39:ILE:HG13	1.98	0.45
1:A:60:LYS:HA	1:A:111:GLU:O	2.17	0.45
1:C:463:LEU:O	1:C:507:ILE:HA	2.16	0.45
1:B:303:ILE:HD13	1:B:328:LEU:HB3	1.98	0.45
1:D:151:LYS:O	1:D:153:HIS:ND1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:531:GLN:OE1	1:D:539:PHE:HB2	2.17	0.45
1:A:518:PHE:HA	1:A:539:PHE:CE1	2.52	0.44
1:B:230:LYS:NZ	1:B:307:CYS:SG	2.87	0.44
1:B:476:HIS:HA	1:B:488:GLU:OE2	2.18	0.44
1:A:13:THR:CG2	1:A:16:ALA:HB2	2.46	0.44
1:B:255:ARG:HH12	1:B:331:GLU:HG2	1.82	0.44
1:A:476:HIS:HA	1:A:488:GLU:OE2	2.17	0.44
1:D:457:GLU:HB3	1:D:458:PHE:CD2	2.52	0.44
1:D:513:TRP:HA	1:D:514:ASN:HA	1.59	0.44
1:B:11:ALA:O	1:B:55:TYR:HA	2.16	0.44
1:B:60:LYS:HA	1:B:111:GLU:O	2.17	0.44
1:D:300:ARG:O	1:D:303:ILE:HG22	2.17	0.44
1:A:540:ASP:O	1:A:541:ARG:HB2	2.18	0.44
1:B:528:GLU:O	1:B:533:HIS:HE1	2.01	0.44
1:C:5:ASN:HD21	1:C:7:ASN:HB2	1.82	0.44
1:A:264:ASP:O	4:A:905:HOH:O	2.21	0.43
1:A:328:LEU:O	1:A:332:ARG:HG2	2.18	0.43
1:C:214:GLU:OE2	1:C:216:LYS:HG2	2.18	0.43
1:C:2:GLU:CD	1:C:65:PRO:HG3	2.43	0.43
1:C:49:TYR:OH	4:C:904:HOH:O	2.20	0.43
1:D:81:ALA:HA	1:D:118:ALA:O	2.18	0.43
1:D:316:LEU:HD12	1:D:326:TYR:CE1	2.53	0.43
1:D:346:HIS:NE2	1:D:389:ASP:OD1	2.37	0.43
1:B:6:PHE:CZ	1:B:57:LYS:HD3	2.54	0.43
1:B:88:LYS:HE3	1:B:90:VAL:HG12	1.99	0.43
1:C:23:MET:HA	1:C:24:PRO:HD3	1.89	0.43
1:C:78:ASN:HA	1:C:79:ALA:HA	1.61	0.43
1:B:207:LYS:HG2	1:B:208:THR:N	2.33	0.43
1:B:315:ARG:HG3	1:B:337:TRP:CD1	2.53	0.43
1:D:320:GLN:HB2	1:D:361:LEU:HD13	2.00	0.43
1:A:11:ALA:HB1	1:A:27:TRP:HB2	2.00	0.43
1:C:15:GLU:OE2	1:C:15:GLU:O	2.36	0.43
1:C:434:ASN:HB3	1:C:467:GLU:HB2	2.00	0.43
1:B:330:ASP:CG	1:B:369:PRO:HD2	2.44	0.43
1:B:590:PHE:CE1	1:B:595:SER:HB2	2.54	0.43
1:C:36:TRP:CZ2	1:C:138:GLY:HA3	2.53	0.43
1:C:10:TRP:CE2	1:C:57:LYS:HD2	2.54	0.43
1:A:158:TYR:OH	3:A:802:FMN:H2'	2.19	0.42
1:B:74:LEU:HD21	1:B:117:ILE:HD13	2.00	0.42
1:C:513:TRP:HE3	4:C:905:HOH:O	2.02	0.42
1:A:78:ASN:HA	1:A:79:ALA:HA	1.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:347:MET:HE1	4:D:930:HOH:O	2.19	0.42
1:A:199:ASP:HB3	1:A:203:LYS:H	1.84	0.42
1:B:412:ILE:HG12	1:B:428:PRO:HG3	2.01	0.42
1:B:594:LYS:C	1:B:594:LYS:HE2	2.44	0.42
1:B:644:LYS:H	1:B:644:LYS:HD3	1.84	0.42
1:C:85:VAL:HG22	1:C:115:ILE:HD12	2.02	0.42
1:C:336:ILE:HG12	1:C:370:SER:HB2	2.02	0.42
1:D:411:THR:HG22	1:D:412:ILE:N	2.34	0.42
1:C:434:ASN:CB	1:C:467:GLU:HB2	2.49	0.42
1:B:351:ARG:NH2	1:B:389:ASP:OD1	2.49	0.42
1:D:413:ALA:HA	1:D:432:SER:O	2.20	0.42
1:C:316:LEU:HD23	1:C:316:LEU:HA	1.93	0.42
1:C:380:ILE:HD12	1:C:380:ILE:HA	1.86	0.42
1:D:446:MET:O	1:D:450:TRP:HB2	2.20	0.42
1:C:337:TRP:CD1	1:C:337:TRP:C	2.98	0.41
1:B:317:ALA:HB1	1:B:318:HIS:ND1	2.34	0.41
1:D:324:TYR:O	1:D:327:ASP:HB2	2.20	0.41
1:B:591:VAL:HG21	1:B:610:VAL:HG13	2.02	0.41
1:D:78:ASN:HA	1:D:79:ALA:HA	1.65	0.41
1:A:472:ALA:HB2	1:A:490:GLN:HB2	2.02	0.41
1:B:81:ALA:HA	1:B:118:ALA:O	2.21	0.41
1:C:1:ARG:HD2	1:C:250:ASP:OD1	2.21	0.41
1:A:158:TYR:CZ	3:A:802:FMN:H2'	2.55	0.41
1:A:545:LYS:O	1:A:548:PHE:HB3	2.21	0.41
1:C:494:HIS:ND1	1:C:551:TYR:OH	2.49	0.41
1:A:316:LEU:HD23	1:A:316:LEU:HA	1.87	0.41
1:A:317:ALA:HB1	1:A:318:HIS:CG	2.56	0.41
1:A:317:ALA:HB1	1:A:318:HIS:ND1	2.35	0.41
1:B:194:VAL:O	1:B:240:VAL:HA	2.20	0.41
1:D:74:LEU:HD21	1:D:117:ILE:HD13	2.03	0.41
1:C:34:HIS:H	1:C:141:TYR:HA	1.86	0.41
1:D:147:ILE:HG22	1:D:149:VAL:HG13	2.03	0.41
1:D:420:ILE:HD12	1:D:420:ILE:H	1.84	0.41
1:B:463:LEU:O	1:B:507:ILE:HA	2.21	0.41
1:D:420:ILE:HD11	1:D:450:TRP:CZ2	2.56	0.41
1:A:473:LEU:HD21	1:A:533:HIS:CE1	2.56	0.41
1:C:332:ARG:NH1	4:C:915:HOH:O	2.37	0.41
1:D:104:ILE:CD1	1:D:107:GLU:HB2	2.52	0.41
1:A:255:ARG:HH22	1:A:331:GLU:HG3	1.85	0.40
1:D:401:VAL:HG13	1:D:402:HIS:N	2.36	0.40
1:B:50:ARG:HA	1:B:121:ASN:OD1	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:287:ARG:HD3	4:B:1082:HOH:O	2.22	0.40
1:C:88:LYS:HE2	1:C:88:LYS:HB2	1.96	0.40
1:C:194:VAL:O	1:C:240:VAL:HA	2.21	0.40
1:C:473:LEU:HD21	1:C:533:HIS:CE1	2.56	0.40
1:A:452:ASP:O	4:A:906:HOH:O	2.22	0.40
1:A:566:LYS:HG2	1:A:629:SER:HB3	2.04	0.40
1:B:175:ALA:HB3	1:B:221:ILE:CG2	2.51	0.40
1:C:227:TRP:CZ2	1:C:333:GLY:HA2	2.57	0.40
3:C:802:FMN:H1'2	3:C:802:FMN:H9	1.42	0.40
1:D:58:GLN:HA	1:D:113:ASN:O	2.22	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	635/780 (81%)	614 (97%)	21 (3%)	0	100	100
1	B	634/780 (81%)	607 (96%)	27 (4%)	0	100	100
1	C	635/780 (81%)	609 (96%)	26 (4%)	0	100	100
1	D	287/780 (37%)	267 (93%)	19 (7%)	1 (0%)	36	50
All	All	2191/3120 (70%)	2097 (96%)	93 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	330	ASP

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	550/662 (83%)	535 (97%)	15 (3%)	39	62
1	B	549/662 (83%)	525 (96%)	24 (4%)	25	43
1	C	550/662 (83%)	534 (97%)	16 (3%)	37	60
1	D	265/662 (40%)	250 (94%)	15 (6%)	18	33
All	All	1914/2648 (72%)	1844 (96%)	70 (4%)	30	51

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	21	LYS
1	A	106	LYS
1	A	171	LYS
1	A	182	VAL
1	A	239	GLU
1	A	303	ILE
1	A	337	TRP
1	A	380	ILE
1	A	387	ASP
1	A	424	TYR
1	A	563	LEU
1	A	594	LYS
1	A	610	VAL
1	A	624	GLU
1	B	15	GLU
1	B	19	VAL
1	B	59	LEU
1	B	60	LYS
1	B	62	SER
1	B	66	GLU
1	B	109	THR
1	B	181	GLU
1	B	203	LYS

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Mol	Chain	Res	Type
1	B	221	ILE
1	B	274	GLU
1	B	294	LEU
1	B	337	TRP
1	B	403	GLU
1	B	424	TYR
1	B	431	ILE
1	B	481	LYS
1	B	482	GLN
1	B	572	VAL
1	B	594	LYS
1	B	599	LEU
1	B	610	VAL
1	B	624	GLU
1	B	632	ARG
1	C	17	THR
1	C	182	VAL
1	C	204	GLU
1	C	205	VAL
1	C	208	THR
1	C	303	ILE
1	C	328	LEU
1	C	337	TRP
1	C	393	ASN
1	C	417	MET
1	C	424	TYR
1	C	431	ILE
1	C	482	GLN
1	C	563	LEU
1	C	610	VAL
1	C	626	LYS
1	D	22	GLU
1	D	102	VAL
1	D	104	ILE
1	D	110	GLU
1	D	152	SER
1	D	297	GLU
1	D	351	ARG
1	D	356	SER
1	D	362	VAL
1	D	377	SER
1	D	424	TYR

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Mol	Chain	Res	Type
1	D	431	ILE
1	D	432	SER
1	D	516	PHE
1	D	538	THR

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	145	ASN
1	A	406	HIS
1	B	398	ASN
1	B	406	HIS
1	B	533	HIS
1	C	150	ASN
1	C	421	HIS
1	D	398	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	FMN	A	802	-	33,33,33	1.05	2 (6%)	48,50,50	1.30	7 (14%)
3	FMN	D	802	-	33,33,33	1.13	3 (9%)	48,50,50	1.75	10 (20%)
3	FMN	C	802	-	33,33,33	1.11	2 (6%)	48,50,50	1.63	12 (25%)
3	FMN	B	802	-	33,33,33	1.02	2 (6%)	48,50,50	1.26	10 (20%)

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	FMN	A	802	-	-	11/18/18/18	0/3/3/3
3	FMN	D	802	-	-	17/18/18/18	0/3/3/3
3	FMN	C	802	-	-	14/18/18/18	0/3/3/3
3	FMN	B	802	-	-	6/18/18/18	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	802	FMN	C4A-N5	3.46	1.38	1.30
3	A	802	FMN	C4A-N5	3.39	1.38	1.30
3	D	802	FMN	C4A-N5	3.37	1.38	1.30
3	B	802	FMN	C4A-N5	3.05	1.37	1.30
3	B	802	FMN	C10-N1	2.99	1.39	1.33
3	C	802	FMN	C10-N1	2.86	1.39	1.33
3	D	802	FMN	C10-N1	2.66	1.38	1.33
3	A	802	FMN	C10-N1	2.47	1.38	1.33
3	D	802	FMN	C1'-N10	2.41	1.53	1.47

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	802	FMN	C1'-N10-C9A	5.01	130.36	120.63
3	D	802	FMN	C4A-C10-N10	4.21	122.51	116.48
3	C	802	FMN	C5A-C9A-N10	3.89	121.49	117.97
3	D	802	FMN	C4-N3-C2	-3.68	119.11	125.64
3	D	802	FMN	C2'-C1'-N10	3.62	127.30	110.20
3	C	802	FMN	C4-N3-C2	-3.51	119.42	125.64
3	A	802	FMN	C4-N3-C2	-3.48	119.46	125.64
3	C	802	FMN	C1'-N10-C9A	-3.44	113.94	120.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	802	FMN	C2'-C1'-N10	3.30	125.78	110.20
3	A	802	FMN	C4A-C10-N10	3.19	121.04	116.48
3	B	802	FMN	C4A-C10-N10	2.86	120.58	116.48
3	D	802	FMN	C4A-C4-N3	2.81	120.40	113.25
3	B	802	FMN	C2'-C1'-N10	2.74	123.14	110.20
3	D	802	FMN	C10-C4A-N5	-2.69	119.31	124.81
3	C	802	FMN	C4A-C4-N3	2.66	120.02	113.25
3	D	802	FMN	C4A-C10-N1	-2.62	118.17	124.59
3	B	802	FMN	C4-N3-C2	-2.57	121.08	125.64
3	C	802	FMN	C9-C9A-N10	-2.57	118.40	121.85
3	C	802	FMN	O4-C4-C4A	-2.56	119.77	126.53
3	A	802	FMN	C4A-C10-N1	-2.56	118.32	124.59
3	A	802	FMN	C4-C4A-C10	2.49	121.21	116.93
3	C	802	FMN	C4A-C10-N1	-2.49	118.50	124.59
3	A	802	FMN	C4A-C4-N3	2.40	119.37	113.25
3	D	802	FMN	O4-C4-C4A	-2.36	120.30	126.53
3	B	802	FMN	C10-C4A-N5	-2.29	120.12	124.81
3	B	802	FMN	C4-C4A-C10	2.29	120.86	116.93
3	B	802	FMN	C5A-C9A-N10	2.26	120.01	117.97
3	A	802	FMN	O4-C4-C4A	-2.25	120.61	126.53
3	C	802	FMN	C9A-C5A-N5	-2.22	120.10	122.45
3	A	802	FMN	C10-C4A-N5	-2.20	120.32	124.81
3	B	802	FMN	O4-C4-C4A	-2.19	120.75	126.53
3	C	802	FMN	C4'-C3'-C2'	2.19	117.21	113.57
3	C	802	FMN	C4-C4A-C10	2.15	120.62	116.93
3	C	802	FMN	O2'-C2'-C3'	2.14	114.26	109.25
3	D	802	FMN	C10-N1-C2	2.13	121.47	116.85
3	B	802	FMN	C4A-C10-N1	-2.12	119.40	124.59
3	D	802	FMN	C4-C4A-C10	2.10	120.53	116.93
3	B	802	FMN	C9A-C5A-N5	-2.08	120.25	122.45
3	B	802	FMN	C4A-C4-N3	2.05	118.48	113.25

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	802	FMN	C1'-C2'-C3'-O3'
3	A	802	FMN	C1'-C2'-C3'-C4'
3	A	802	FMN	O2'-C2'-C3'-O3'
3	A	802	FMN	C3'-C4'-C5'-O5'
3	A	802	FMN	O4'-C4'-C5'-O5'
3	B	802	FMN	C2'-C3'-C4'-O4'

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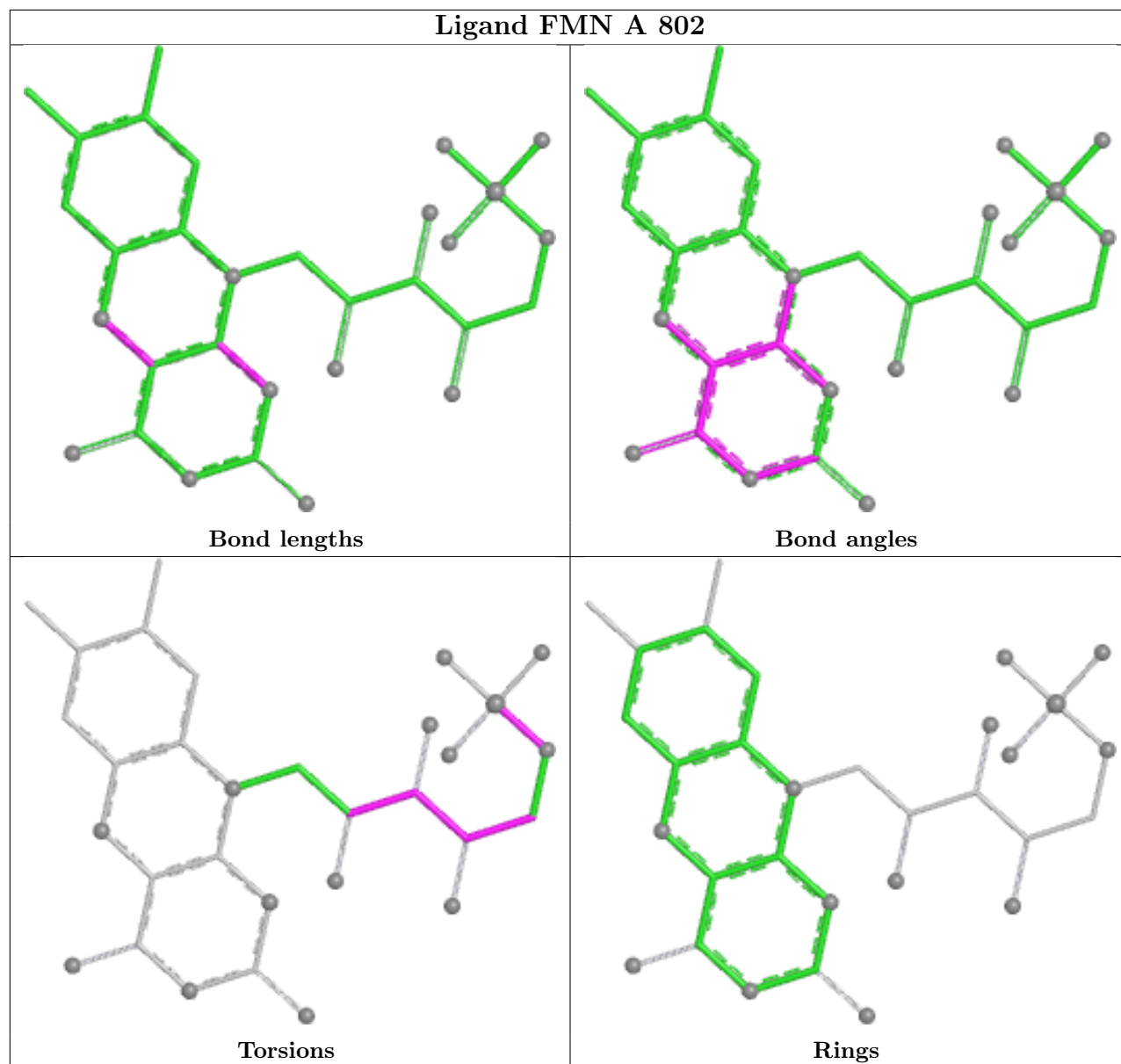
Mol	Chain	Res	Type	Atoms
3	B	802	FMN	C2'-C3'-C4'-C5'
3	B	802	FMN	O3'-C3'-C4'-O4'
3	B	802	FMN	O3'-C3'-C4'-C5'
3	B	802	FMN	C3'-C4'-C5'-O5'
3	B	802	FMN	O4'-C4'-C5'-O5'
3	C	802	FMN	C2'-C1'-N10-C10
3	C	802	FMN	N10-C1'-C2'-O2'
3	C	802	FMN	N10-C1'-C2'-C3'
3	C	802	FMN	C1'-C2'-C3'-O3'
3	C	802	FMN	C1'-C2'-C3'-C4'
3	C	802	FMN	O2'-C2'-C3'-O3'
3	C	802	FMN	O2'-C2'-C3'-C4'
3	C	802	FMN	C2'-C3'-C4'-O4'
3	C	802	FMN	O3'-C3'-C4'-O4'
3	C	802	FMN	O3'-C3'-C4'-C5'
3	C	802	FMN	C5'-O5'-P-O2P
3	C	802	FMN	C5'-O5'-P-O3P
3	D	802	FMN	C2'-C1'-N10-C9A
3	D	802	FMN	C2'-C1'-N10-C10
3	D	802	FMN	N10-C1'-C2'-O2'
3	D	802	FMN	N10-C1'-C2'-C3'
3	D	802	FMN	C1'-C2'-C3'-O3'
3	D	802	FMN	C1'-C2'-C3'-C4'
3	D	802	FMN	O2'-C2'-C3'-O3'
3	D	802	FMN	O2'-C2'-C3'-C4'
3	D	802	FMN	C2'-C3'-C4'-O4'
3	D	802	FMN	O3'-C3'-C4'-O4'
3	D	802	FMN	C3'-C4'-C5'-O5'
3	D	802	FMN	C5'-O5'-P-O1P
3	D	802	FMN	C5'-O5'-P-O2P
3	D	802	FMN	C5'-O5'-P-O3P
3	A	802	FMN	O2'-C2'-C3'-C4'
3	D	802	FMN	O3'-C3'-C4'-C5'
3	C	802	FMN	C2'-C3'-C4'-C5'
3	D	802	FMN	C2'-C3'-C4'-C5'
3	C	802	FMN	C5'-O5'-P-O1P
3	A	802	FMN	C2'-C3'-C4'-O4'
3	A	802	FMN	O3'-C3'-C4'-O4'
3	A	802	FMN	C2'-C3'-C4'-C5'
3	D	802	FMN	O4'-C4'-C5'-O5'
3	A	802	FMN	C5'-O5'-P-O1P
3	A	802	FMN	O3'-C3'-C4'-C5'

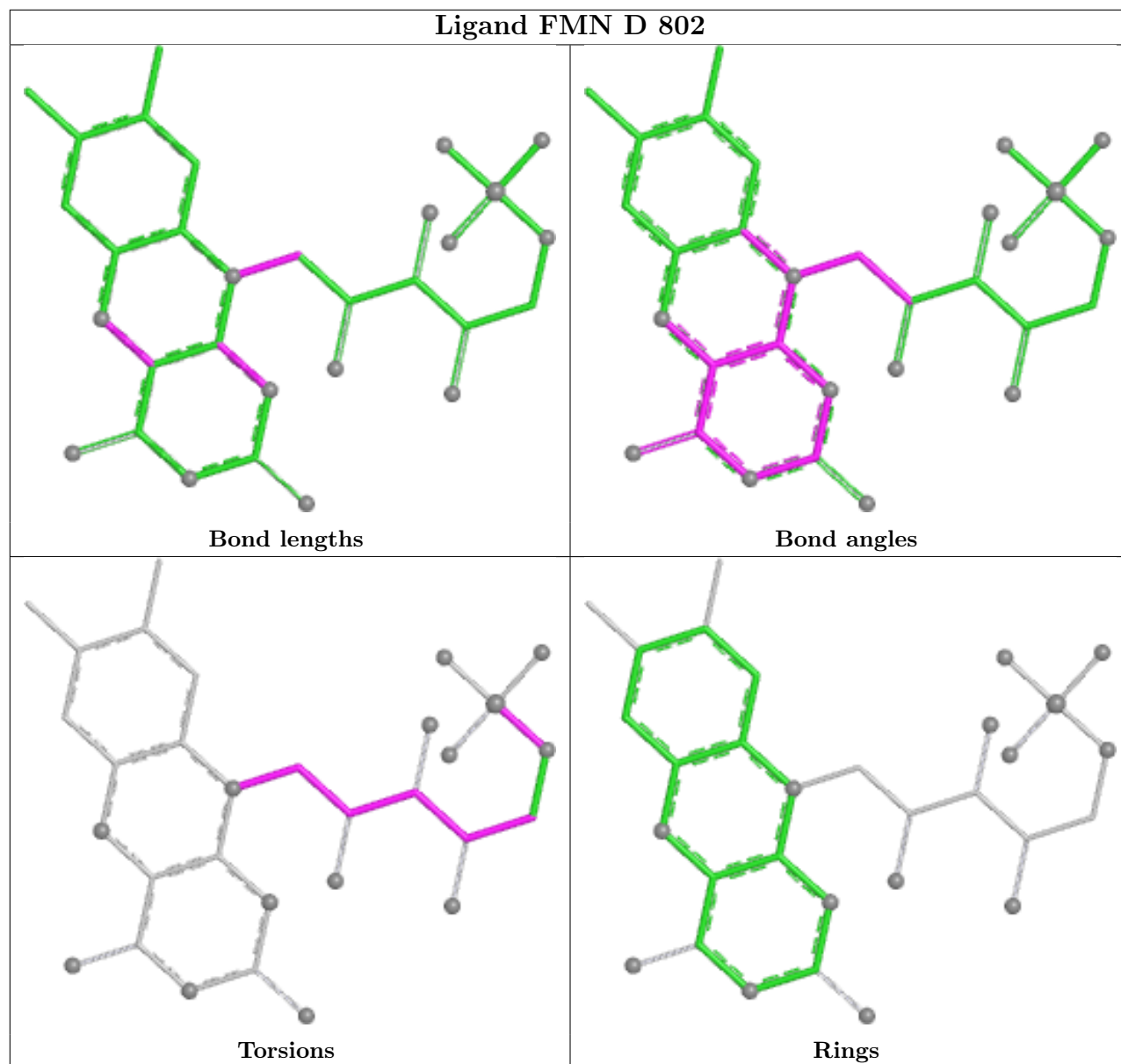
There are no ring outliers.

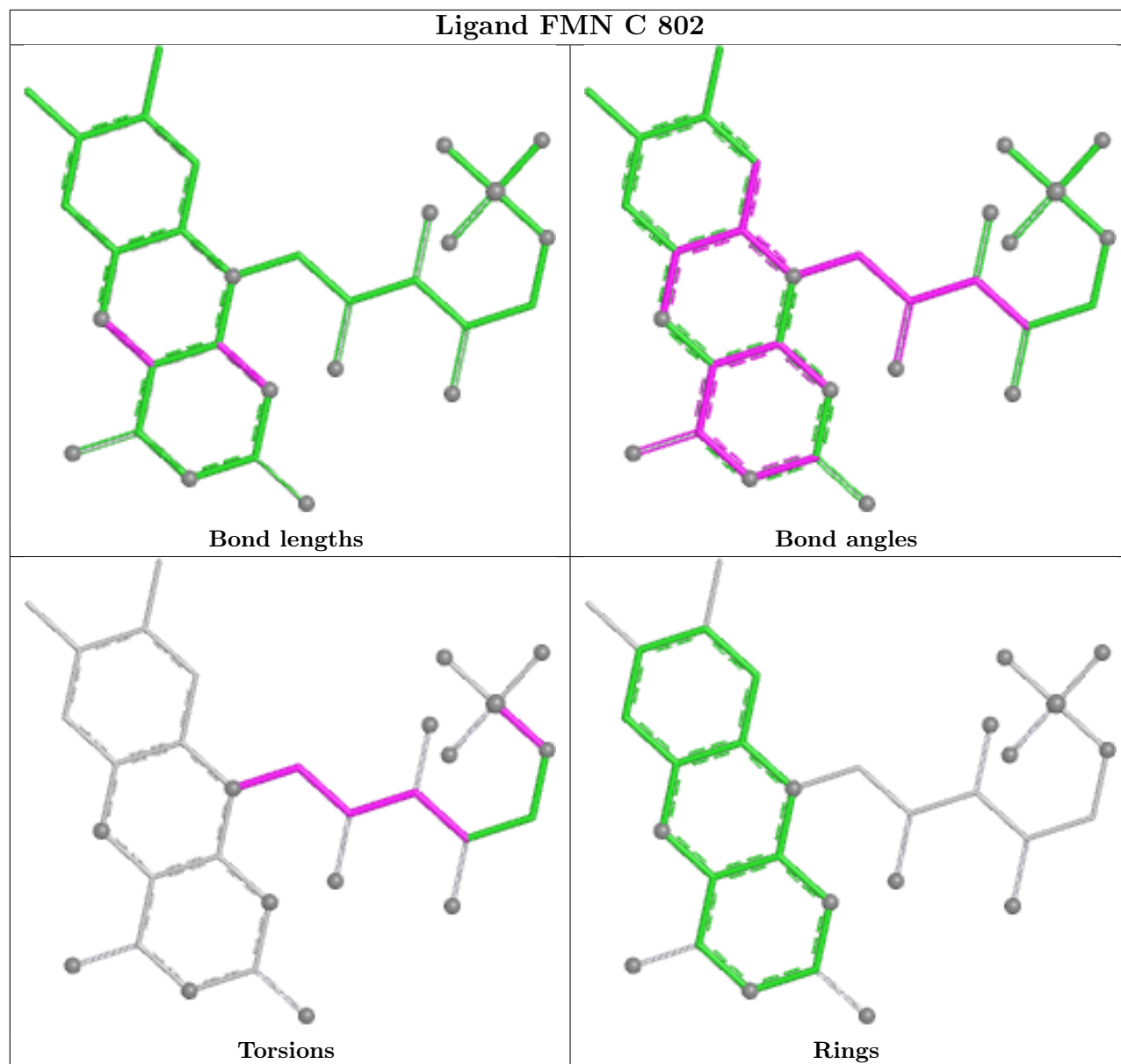
3 monomers are involved in 7 short contacts:

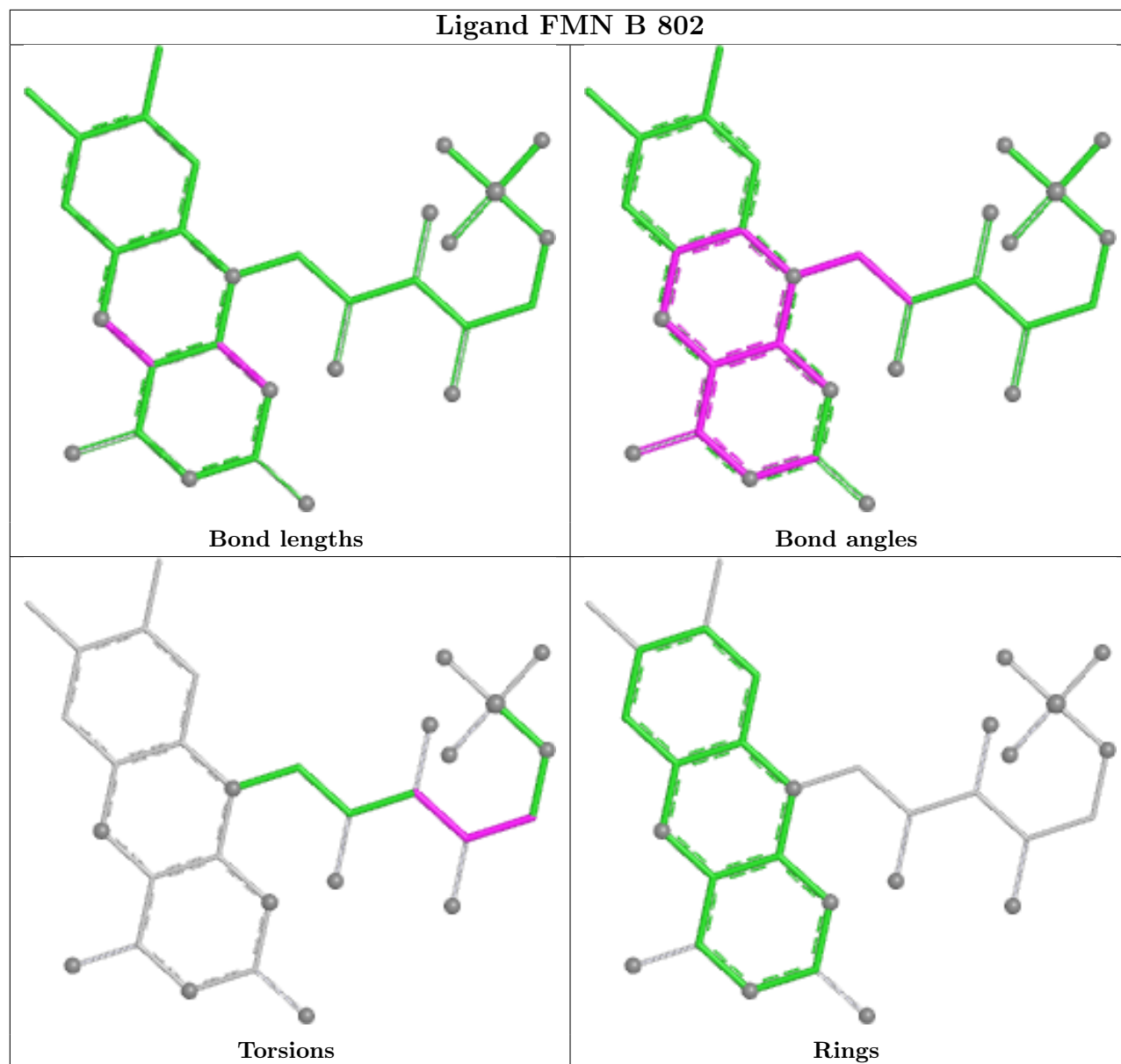
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	802	FMN	3	0
3	D	802	FMN	2	0
3	C	802	FMN	2	0

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	639/780 (81%)	-0.29	4 (0%) 85 83	19, 28, 46, 59	0
1	B	638/780 (81%)	-0.07	5 (0%) 82 80	22, 32, 56, 72	0
1	C	638/780 (81%)	-0.17	0 100 100	19, 32, 52, 62	1 (0%)
1	D	305/780 (39%)	0.76	38 (12%) 8 6	25, 46, 64, 80	0
All	All	2220/3120 (71%)	-0.05	47 (2%) 63 59	19, 32, 56, 80	1 (0%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	17	THR	4.1
1	D	375	GLY	4.0
1	A	16	ALA	3.9
1	D	162	PRO	3.7
1	D	154	PHE	3.6
1	D	149	VAL	3.4
1	D	451	MET	2.9
1	D	411	THR	2.9
1	B	189	ALA	2.9
1	D	317	ALA	2.7
1	D	304	ASP	2.7
1	D	445	SER	2.6
1	B	188	ALA	2.6
1	D	448	GLY	2.5
1	D	363	VAL	2.5
1	D	148	ALA	2.5
1	D	102	VAL	2.4
1	D	444	VAL	2.4
1	A	382	MET	2.4
1	D	403	GLU	2.4
1	D	156	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	440	TYR	2.4
1	D	159	TYR	2.4
1	D	362	VAL	2.4
1	A	246	GLU	2.4
1	D	382	MET	2.3
1	D	328	LEU	2.3
1	D	157	ASP	2.3
1	D	527	GLY	2.3
1	D	155	ASP	2.3
1	D	455	HIS	2.2
1	D	326	TYR	2.2
1	D	463	LEU	2.2
1	D	158	TYR	2.1
1	D	151	LYS	2.1
1	B	186	ASN	2.1
1	D	434	ASN	2.1
1	D	161	GLY	2.1
1	D	536	LEU	2.1
1	D	20	PRO	2.1
1	D	365	ASN	2.0
1	B	440	TYR	2.0
1	D	462	PRO	2.0
1	D	402	HIS	2.0
1	D	453	ASN	2.0
1	D	420	ILE	2.0
1	B	244	SER	2.0

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

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

6.3 Carbohydrates [i](#)

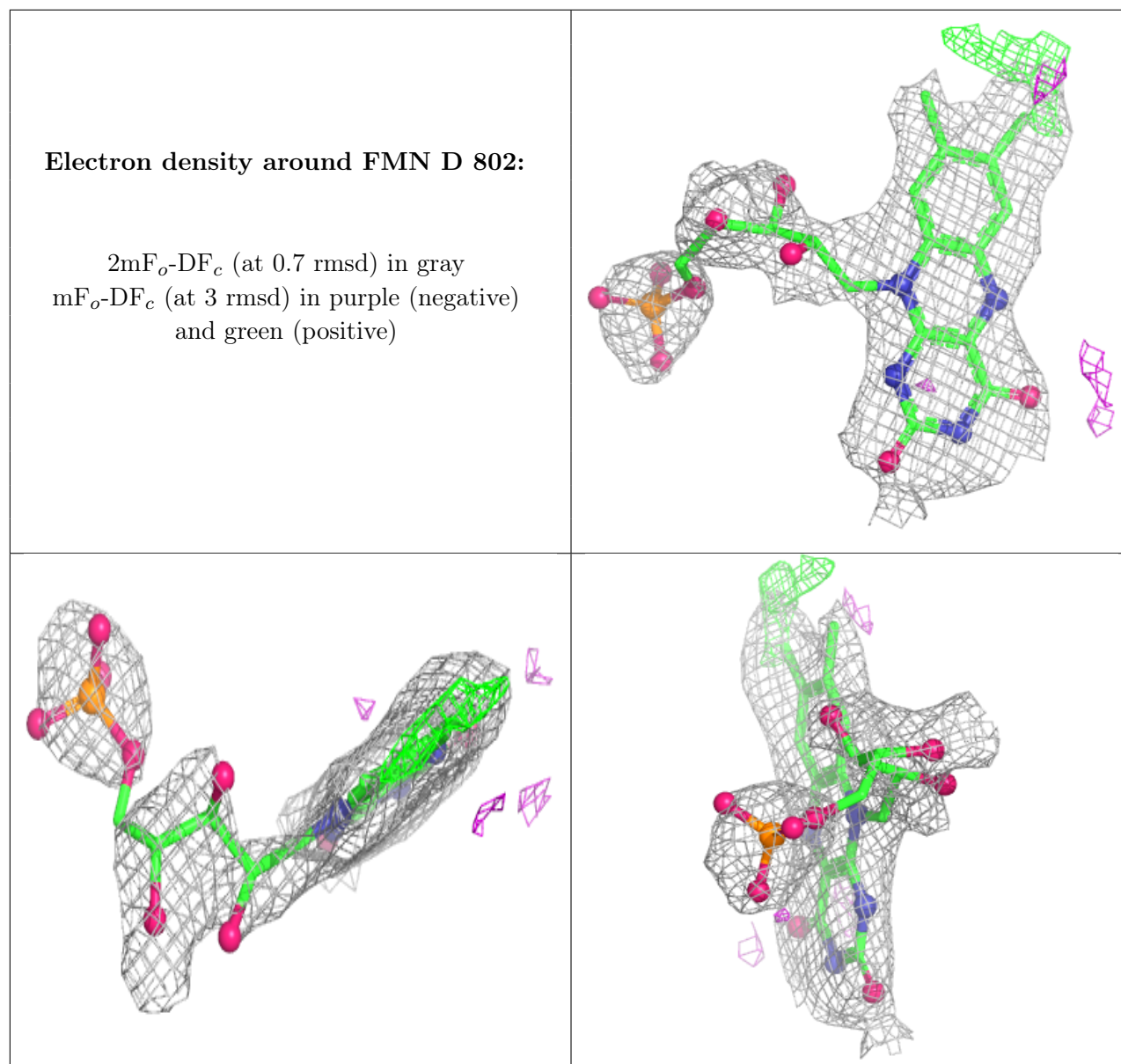
There are no oligosaccharides in this entry.

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.

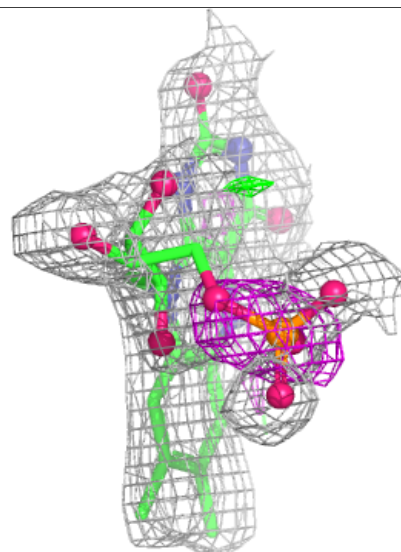
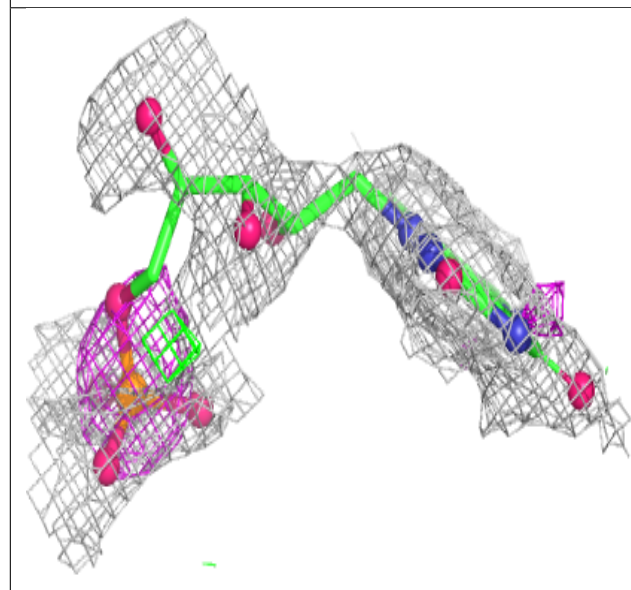
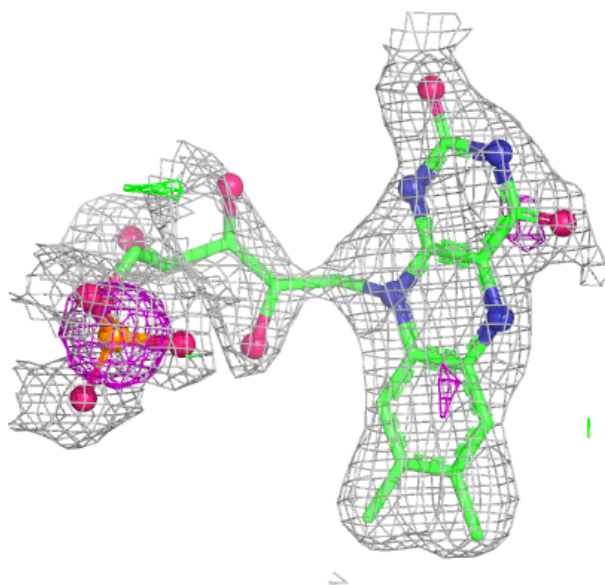
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	FMN	D	802	31/31	0.64	0.17	42,61,85,90	0
3	FMN	A	802	31/31	0.71	0.15	29,41,67,88	0
3	FMN	C	802	31/31	0.77	0.13	39,49,76,100	0
3	FMN	B	802	31/31	0.82	0.12	36,43,86,88	0
2	CA	A	801	1/1	0.97	0.02	26,26,26,26	0
2	CA	C	801	1/1	0.99	0.02	26,26,26,26	0
2	CA	D	801	1/1	0.99	0.02	33,33,33,33	0
2	CA	B	801	1/1	0.99	0.01	23,23,23,23	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.



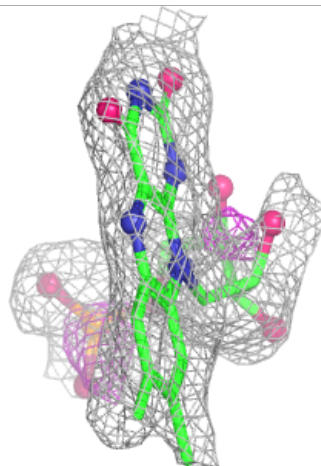
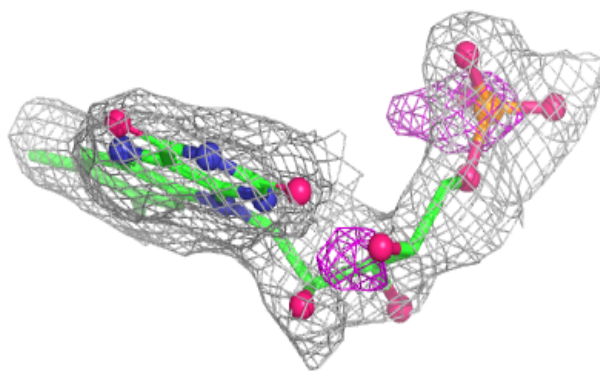
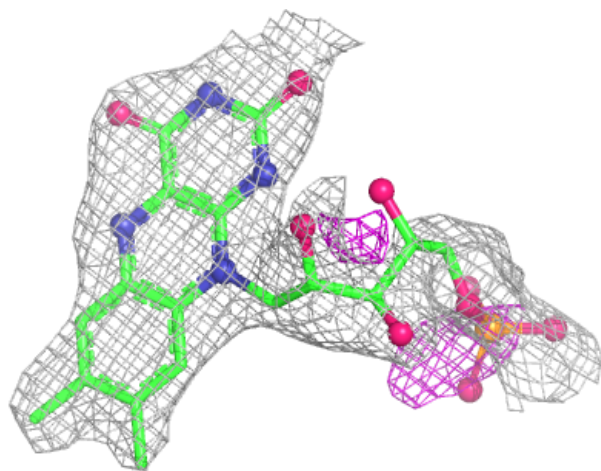
Electron density around FMN A 802:

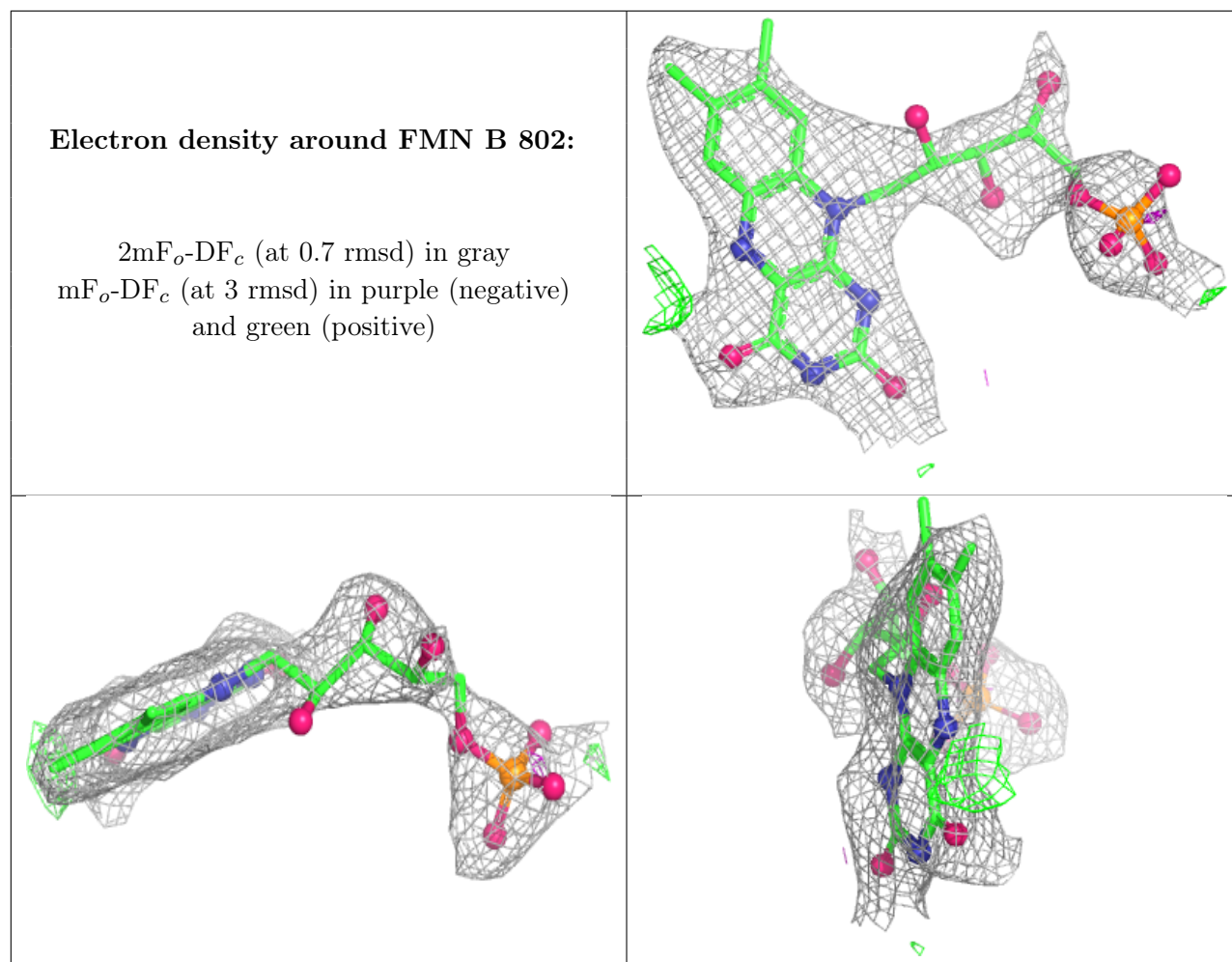
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FMN C 802:

$2mF_o-DF_c$ (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.