Full wwPDB X-ray Structure Validation Report (i)

Mar 5, 2026 – 11:38 AM UTC

PDB ID	:	7T6H / pdb_00007t6h
Title	:	E. coli dihydroorotate dehydrogenase
Authors	:	Horwitz, S.M.; Ambarian, J.A.; Davis, K.M.
Deposited on	:	2021-12-13
Resolution	:	2.42 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 (i) 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 (i)) were used in the production of this report:

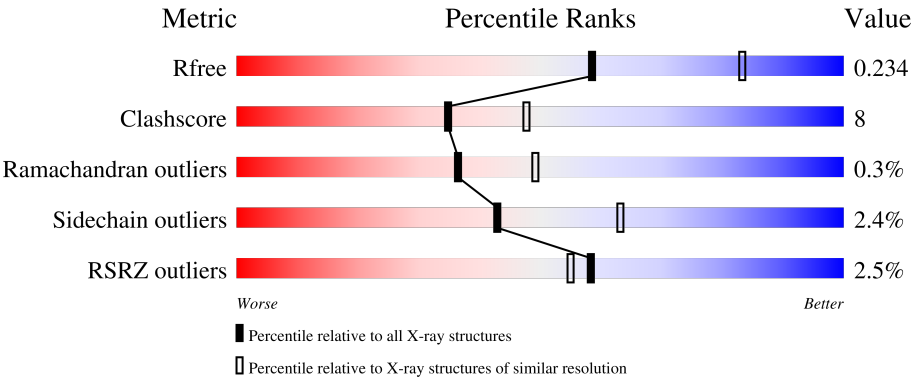
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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 <=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	368	3% 79% 11% 9%
1	B	368	% 73% 18% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ORO	B	401	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5344 atoms, of which 1 is hydrogen 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 Dihydroorotate dehydrogenase (quinone).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	1	0	0
			2553	1617	450	476	10			
1	B	335	Total	C	N	O	S	0	0	0
			2572	1633	451	478	10			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-31	HIS	-	expression tag	UNP P0A7E1
A	-30	HIS	-	expression tag	UNP P0A7E1
A	-29	HIS	-	expression tag	UNP P0A7E1
A	-28	HIS	-	expression tag	UNP P0A7E1
A	-27	HIS	-	expression tag	UNP P0A7E1
A	-26	HIS	-	expression tag	UNP P0A7E1
A	-25	SER	-	expression tag	UNP P0A7E1
A	-24	SER	-	expression tag	UNP P0A7E1
A	-23	GLY	-	expression tag	UNP P0A7E1
A	-22	LEU	-	expression tag	UNP P0A7E1
A	-21	VAL	-	expression tag	UNP P0A7E1
A	-20	PRO	-	expression tag	UNP P0A7E1
A	-19	ARG	-	expression tag	UNP P0A7E1
A	-18	GLY	-	expression tag	UNP P0A7E1
A	-17	SER	-	expression tag	UNP P0A7E1
A	-16	HIS	-	expression tag	UNP P0A7E1
A	-15	MET	-	expression tag	UNP P0A7E1
A	-14	ALA	-	expression tag	UNP P0A7E1
A	-13	SER	-	expression tag	UNP P0A7E1
A	-12	MET	-	expression tag	UNP P0A7E1
A	-11	THR	-	expression tag	UNP P0A7E1
A	-10	GLY	-	expression tag	UNP P0A7E1
A	-9	GLY	-	expression tag	UNP P0A7E1
A	-8	GLN	-	expression tag	UNP P0A7E1
A	-7	GLN	-	expression tag	UNP P0A7E1

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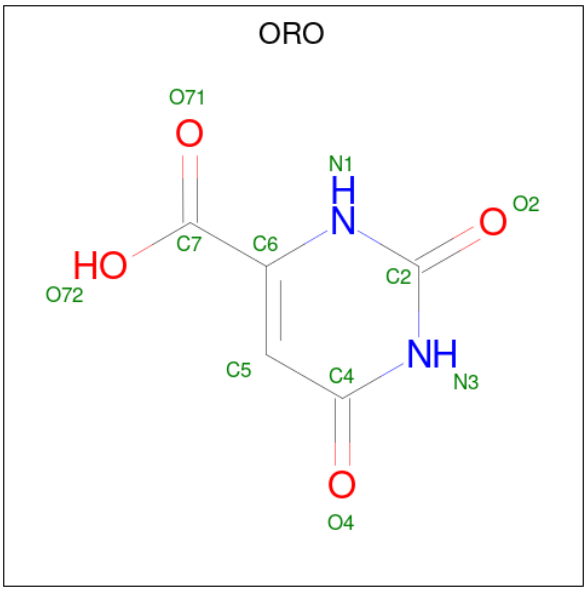
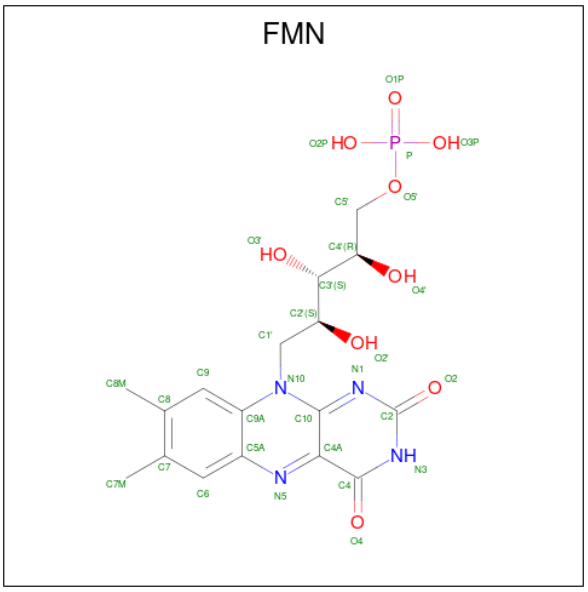
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	expression tag	UNP P0A7E1
A	-5	GLY	-	expression tag	UNP P0A7E1
A	-4	ARG	-	expression tag	UNP P0A7E1
A	-3	GLY	-	expression tag	UNP P0A7E1
A	-2	SER	-	expression tag	UNP P0A7E1
A	-1	GLU	-	expression tag	UNP P0A7E1
A	0	PHE	-	expression tag	UNP P0A7E1
B	-31	HIS	-	expression tag	UNP P0A7E1
B	-30	HIS	-	expression tag	UNP P0A7E1
B	-29	HIS	-	expression tag	UNP P0A7E1
B	-28	HIS	-	expression tag	UNP P0A7E1
B	-27	HIS	-	expression tag	UNP P0A7E1
B	-26	HIS	-	expression tag	UNP P0A7E1
B	-25	SER	-	expression tag	UNP P0A7E1
B	-24	SER	-	expression tag	UNP P0A7E1
B	-23	GLY	-	expression tag	UNP P0A7E1
B	-22	LEU	-	expression tag	UNP P0A7E1
B	-21	VAL	-	expression tag	UNP P0A7E1
B	-20	PRO	-	expression tag	UNP P0A7E1
B	-19	ARG	-	expression tag	UNP P0A7E1
B	-18	GLY	-	expression tag	UNP P0A7E1
B	-17	SER	-	expression tag	UNP P0A7E1
B	-16	HIS	-	expression tag	UNP P0A7E1
B	-15	MET	-	expression tag	UNP P0A7E1
B	-14	ALA	-	expression tag	UNP P0A7E1
B	-13	SER	-	expression tag	UNP P0A7E1
B	-12	MET	-	expression tag	UNP P0A7E1
B	-11	THR	-	expression tag	UNP P0A7E1
B	-10	GLY	-	expression tag	UNP P0A7E1
B	-9	GLY	-	expression tag	UNP P0A7E1
B	-8	GLN	-	expression tag	UNP P0A7E1
B	-7	GLN	-	expression tag	UNP P0A7E1
B	-6	MET	-	expression tag	UNP P0A7E1
B	-5	GLY	-	expression tag	UNP P0A7E1
B	-4	ARG	-	expression tag	UNP P0A7E1
B	-3	GLY	-	expression tag	UNP P0A7E1
B	-2	SER	-	expression tag	UNP P0A7E1
B	-1	GLU	-	expression tag	UNP P0A7E1
B	0	PHE	-	expression tag	UNP P0A7E1

- Molecule 2 is OROTIC ACID (CCD ID: ORO) (formula: $C_5H_4N_2O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 11 5 2 4	0	0
2	B	1	Total C N O 11 5 2 4	0	0

- Molecule 3 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
3	A	1	Total C N O P 31 17 4 9 1	0	0

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	H	N	O	P	0	0
			32	17	1	4	9	1		

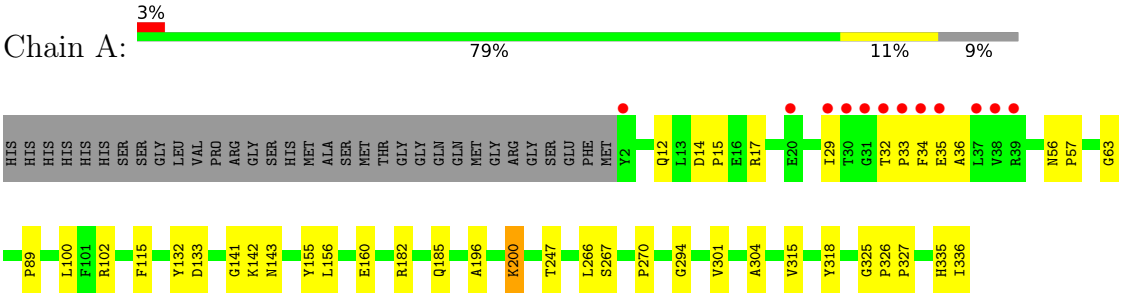
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	72	Total	O	0	0
			72	72		
4	B	62	Total	O	0	0
			62	62		

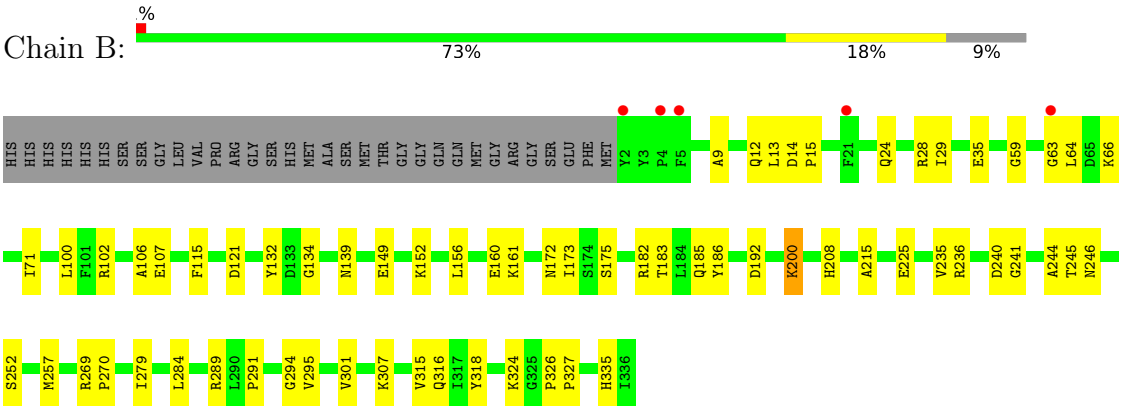
3 Residue-property plots (i)

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: Dihydroorotate dehydrogenase (quinone)



• Molecule 1: Dihydroorotate dehydrogenase (quinone)



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	104.53Å 169.50Å 129.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.70 – 2.42 49.70 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.70-2.42) 99.2 (49.70-2.42)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.33 (at 2.42Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.200 , 0.235 0.201 , 0.234	Depositor DCC
R_{free} test set	1617 reflections (2.96%)	wwPDB-VP
Wilson B-factor (Å^2)	57.1	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 33.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5344	wwPDB-VP
Average B, all atoms (Å^2)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ORO, 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	0.51	0/2596	0.73	1/3514 (0.0%)
1	B	0.49	0/2613	0.71	0/3533
All	All	0.50	0/5209	0.72	1/7047 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	12	GLN	N-CA-C	-5.22	106.17	112.54

There are no chirality outliers.

There are no planarity outliers.

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	2553	0	2603	27	0
1	B	2572	0	2630	52	0
2	A	11	0	3	2	0
2	B	11	0	3	10	0
3	A	31	0	19	1	0
3	B	31	1	19	0	0
4	A	72	0	0	0	0
4	B	62	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5343	1	5277	79	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 8.

All (79) 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:318:TYR:CD2	1:B:318:TYR:CD1	2.37	1.04
1:B:115:PHE:HE1	2:B:401:ORO:H5	1.25	0.98
1:B:115:PHE:CE1	2:B:401:ORO:H5	2.01	0.94
1:B:318:TYR:CD1	1:B:318:TYR:CB	2.54	0.90
1:B:318:TYR:CD2	1:B:318:TYR:CB	2.54	0.90
1:B:63:GLY:CA	1:B:64:LEU:N	2.45	0.80
1:B:175:SER:HB2	2:B:401:ORO:C4	2.16	0.74
1:B:63:GLY:CA	1:B:63:GLY:O	2.37	0.73
1:A:115:PHE:HE1	2:A:401:ORO:H5	1.54	0.72
1:B:63:GLY:O	1:B:64:LEU:N	2.24	0.70
1:A:15:PRO:HG3	1:A:102:ARG:HG3	1.74	0.68
1:B:246:ASN:ND2	2:B:401:ORO:O4	2.25	0.68
1:A:115:PHE:CE1	2:A:401:ORO:H5	2.29	0.67
1:B:173:ILE:HD12	1:B:185:GLN:HB3	1.78	0.65
1:B:301:VAL:HG21	1:B:335:HIS:HB2	1.81	0.62
1:B:246:ASN:HD21	2:B:401:ORO:C4	2.13	0.61
1:A:182:ARG:O	1:A:185:GLN:HG2	2.01	0.61
1:B:172:ASN:ND2	2:B:401:ORO:O4	2.30	0.57
1:B:139:ASN:OD1	1:B:172:ASN:HB2	2.05	0.57
1:A:34:PHE:C	1:A:36:ALA:H	2.12	0.56
1:B:29:ILE:HD12	1:B:35:GLU:HA	1.88	0.54
1:A:56:ASN:HB2	1:A:57:PRO:CD	2.39	0.53
1:B:240:ASP:O	1:B:291:PRO:HD2	2.10	0.52
1:A:247:THR:HG22	1:A:267:SER:HB3	1.91	0.51
1:A:34:PHE:C	1:A:36:ALA:N	2.66	0.51
1:B:175:SER:HB2	2:B:401:ORO:C5	2.40	0.51
1:B:215:ALA:HA	1:B:241:GLY:O	2.11	0.51
1:B:121:ASP:OD1	1:B:161:LYS:HE2	2.12	0.50
1:B:175:SER:CB	2:B:401:ORO:C4	2.88	0.50
1:B:132:TYR:CZ	1:B:134:GLY:HA3	2.47	0.49
1:A:32:THR:C	1:A:34:PHE:H	2.21	0.49
1:B:326:PRO:N	1:B:327:PRO:CD	2.76	0.49
1:A:318:TYR:CD1	1:A:318:TYR:C	2.90	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:TYR:C	1:A:318:TYR:HD1	2.21	0.48
1:B:235:VAL:HG11	1:B:289:ARG:NH1	2.28	0.48
1:B:71:ILE:HG22	1:B:132:TYR:HB2	1.96	0.48
1:A:301:VAL:HG11	1:A:335:HIS:HB2	1.95	0.48
1:A:304:ALA:HB1	1:A:336:ILE:HD12	1.96	0.47
1:A:156:LEU:HG	1:A:200:LYS:HG2	1.97	0.47
1:B:208:HIS:HE1	4:B:542:HOH:O	1.98	0.47
1:B:24:GLN:O	1:B:28:ARG:HG3	2.14	0.46
1:B:208:HIS:HD2	4:B:560:HOH:O	1.99	0.46
1:B:284:LEU:HD23	1:B:284:LEU:HA	1.69	0.46
1:B:318:TYR:CD1	1:B:318:TYR:CA	2.99	0.46
1:A:141:GLY:HA2	1:A:155:TYR:CE1	2.51	0.46
1:B:152:LYS:O	1:B:156:LEU:HD13	2.16	0.46
1:B:15:PRO:HG3	1:B:102:ARG:HG3	1.98	0.46
1:B:107:GLU:OE1	1:B:324:LYS:NZ	2.45	0.45
1:B:244:ALA:HA	1:B:245:THR:HA	1.66	0.45
1:B:318:TYR:CD1	1:B:318:TYR:C	2.95	0.45
1:B:183:THR:HA	1:B:186:TYR:CD2	2.52	0.45
1:B:246:ASN:CG	2:B:401:ORO:HN3	2.24	0.45
1:A:156:LEU:O	1:A:160:GLU:HG3	2.17	0.44
1:A:326:PRO:N	1:A:327:PRO:CD	2.81	0.44
1:B:182:ARG:O	1:B:185:GLN:HG2	2.17	0.44
1:B:149:GLU:H	1:B:149:GLU:CD	2.26	0.44
1:B:252:SER:O	4:B:501:HOH:O	2.20	0.44
1:A:133:ASP:OD1	1:A:133:ASP:N	2.51	0.44
1:B:200:LYS:HA	1:B:200:LYS:HD3	1.72	0.43
1:B:59:GLY:O	1:B:316:GLN:HA	2.19	0.42
1:A:89:PRO:HD2	1:A:143:ASN:OD1	2.20	0.42
1:A:266:LEU:HD21	1:A:270:PRO:HB2	2.00	0.42
1:B:175:SER:HB2	2:B:401:ORO:O4	2.20	0.42
1:B:208:HIS:CE1	4:B:542:HOH:O	2.73	0.42
1:B:294:GLY:O	1:B:315:VAL:HA	2.19	0.42
1:A:294:GLY:O	1:A:315:VAL:HA	2.18	0.42
1:B:9:ALA:O	1:B:12:GLN:HB2	2.20	0.42
1:A:156:LEU:HD21	1:A:196:ALA:HB1	2.02	0.42
1:A:63:GLY:HA3	3:A:402:FMN:N5	2.35	0.41
1:B:160:GLU:OE2	1:B:200:LYS:HE3	2.20	0.41
1:A:15:PRO:HB2	1:A:100:LEU:HG	2.01	0.41
1:A:142:LYS:HG2	1:A:143:ASN:O	2.21	0.41
1:A:325:GLY:C	1:A:327:PRO:HD2	2.46	0.41
1:B:106:ALA:H	1:B:257:MET:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:PRO:HB2	1:B:100:LEU:HG	2.03	0.41
1:B:269:ARG:N	1:B:270:PRO:CD	2.84	0.41
1:B:279:ILE:CD1	1:B:307:LYS:HD3	2.51	0.40
1:A:200:LYS:HA	1:A:200:LYS:HD3	1.56	0.40
1:B:63:GLY:O	1:B:64:LEU:CA	2.70	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	333/368 (90%)	322 (97%)	9 (3%)	2 (1%)	21 30
1	B	331/368 (90%)	320 (97%)	11 (3%)	0	100 100
All	All	664/736 (90%)	642 (97%)	20 (3%)	2 (0%)	36 49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	TYR
1	A	33	PRO

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	272/304 (90%)	267 (98%)	5 (2%)	51 71
1	B	276/304 (91%)	268 (97%)	8 (3%)	37 57
All	All	548/608 (90%)	535 (98%)	13 (2%)	43 63

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ASP
1	A	17	ARG
1	A	29	ILE
1	A	35	GLU
1	A	200	LYS
1	B	13	LEU
1	B	14	ASP
1	B	66	LYS
1	B	192	ASP
1	B	200	LYS
1	B	225	GLU
1	B	236	ARG
1	B	295	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	208	HIS
1	A	335	HIS
1	B	12	GLN
1	B	177	ASN
1	B	255	GLN
1	B	335	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no oligosaccharides 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	$\#\|Z\| > 2$
2	ORO	A	401	-	11,11,11	1.52	2 (18%)	14,15,15	1.31	1 (7%)
3	FMN	B	402	-	33,33,33	1.06	2 (6%)	48,50,50	1.45	11 (22%)
3	FMN	A	402	-	33,33,33	1.22	3 (9%)	48,50,50	1.38	6 (12%)
2	ORO	B	401	-	11,11,11	1.10	1 (9%)	14,15,15	0.76	1 (7%)

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	ORO	A	401	-	-	2/4/4/4	0/1/1/1
3	FMN	B	402	-	-	3/18/18/18	0/3/3/3
3	FMN	A	402	-	-	1/18/18/18	0/3/3/3
2	ORO	B	401	-	-	4/4/4/4	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	FMN	C4A-N5	3.65	1.38	1.30
2	A	401	ORO	O72-C7	-3.50	1.21	1.30
3	A	402	FMN	C4A-N5	3.45	1.38	1.30
2	B	401	ORO	O72-C7	-3.19	1.21	1.30
3	A	402	FMN	C10-N1	2.61	1.38	1.33
2	A	401	ORO	C6-C7	-2.31	1.43	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	FMN	C5'-C4'	2.05	1.54	1.51
3	A	402	FMN	C4A-C10	-2.02	1.38	1.44

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	402	FMN	C4'-C3'-C2'	4.47	121.02	113.57
2	A	401	ORO	O71-C7-C6	-3.48	113.96	121.01
3	B	402	FMN	C5A-C9A-N10	3.27	120.92	117.97
3	B	402	FMN	C4A-C10-N10	2.99	120.76	116.48
3	B	402	FMN	C4-N3-C2	-2.72	120.81	125.64
3	B	402	FMN	O4-C4-C4A	-2.56	119.77	126.53
3	A	402	FMN	O4-C4-C4A	-2.54	119.83	126.53
3	A	402	FMN	C4A-C4-N3	2.51	119.65	113.25
3	A	402	FMN	C9A-C5A-N5	-2.39	119.92	122.45
3	A	402	FMN	C5A-C9A-N10	2.32	120.07	117.97
3	B	402	FMN	C10-C4A-N5	-2.29	120.14	124.81
3	B	402	FMN	C9A-C5A-N5	-2.28	120.03	122.45
3	B	402	FMN	C4A-C4-N3	2.27	119.04	113.25
3	B	402	FMN	C9-C9A-N10	-2.26	118.81	121.85
3	A	402	FMN	C4-N3-C2	-2.21	121.72	125.64
3	B	402	FMN	C10-N1-C2	2.17	121.55	116.85
3	B	402	FMN	C4'-C3'-C2'	2.08	117.04	113.57
3	B	402	FMN	C4A-C10-N1	-2.06	119.54	124.59
2	B	401	ORO	O71-C7-C6	-2.04	116.87	121.01

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ORO	N1-C6-C7-O71
2	A	401	ORO	N1-C6-C7-O72
2	B	401	ORO	N1-C6-C7-O71
2	B	401	ORO	N1-C6-C7-O72
2	B	401	ORO	C5-C6-C7-O72
2	B	401	ORO	C5-C6-C7-O71
3	B	402	FMN	C4'-C5'-O5'-P
3	B	402	FMN	C2'-C3'-C4'-O4'
3	A	402	FMN	C4'-C5'-O5'-P
3	B	402	FMN	O3'-C3'-C4'-O4'

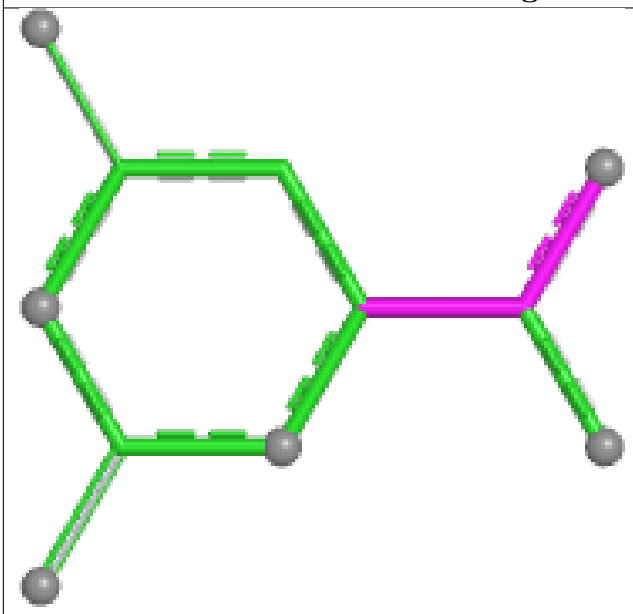
There are no ring outliers.

3 monomers are involved in 13 short contacts:

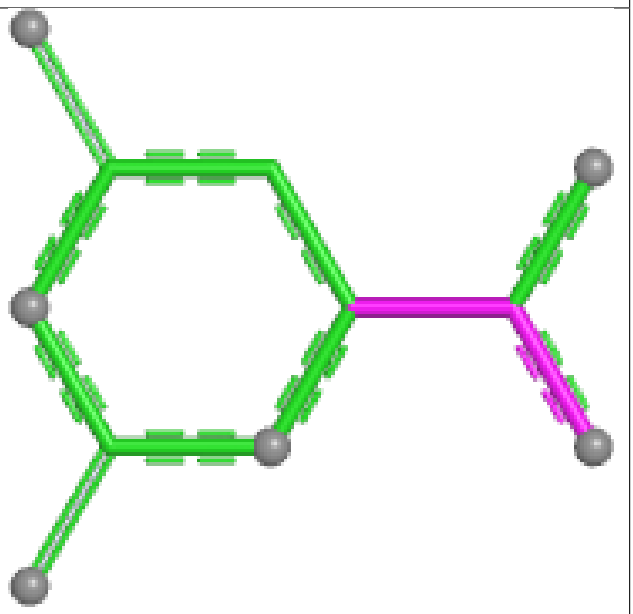
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ORO	2	0
3	A	402	FMN	1	0
2	B	401	ORO	10	0

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

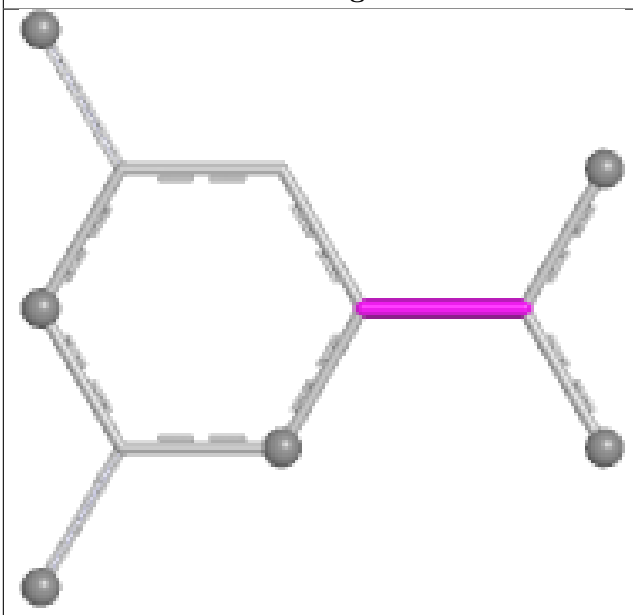
Ligand ORO A 401



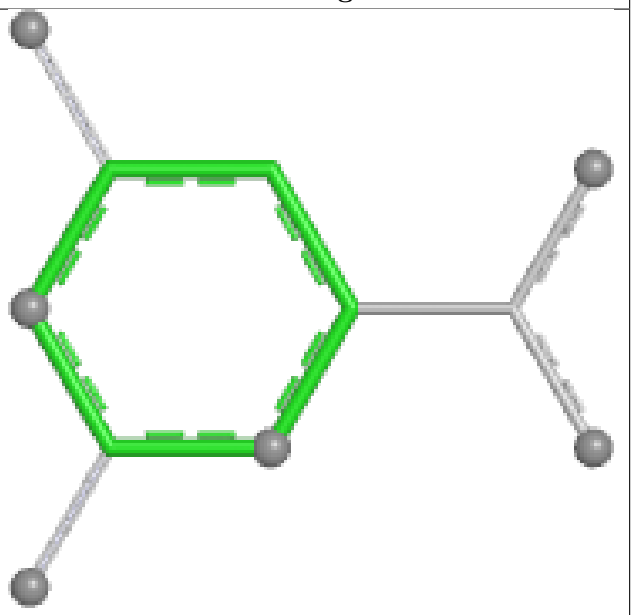
Bond lengths



Bond angles

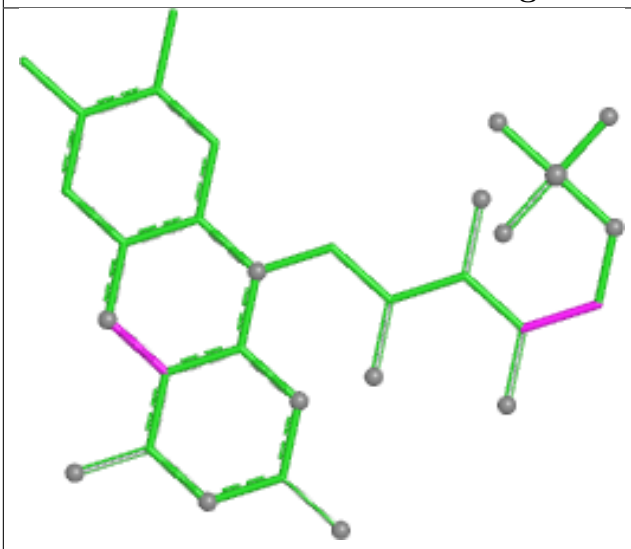


Torsions

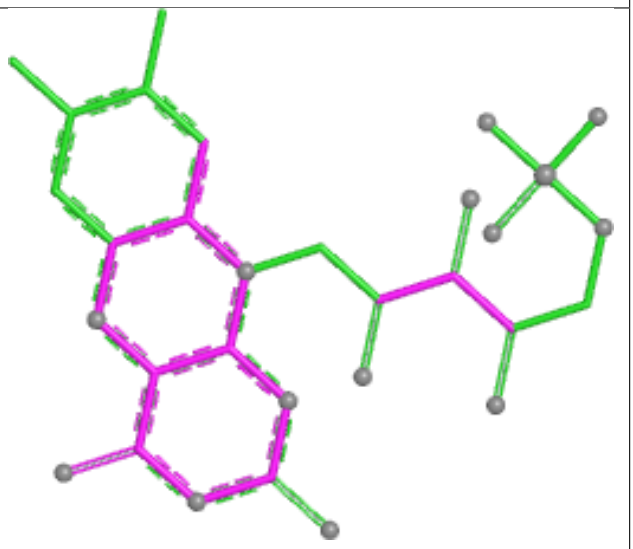


Rings

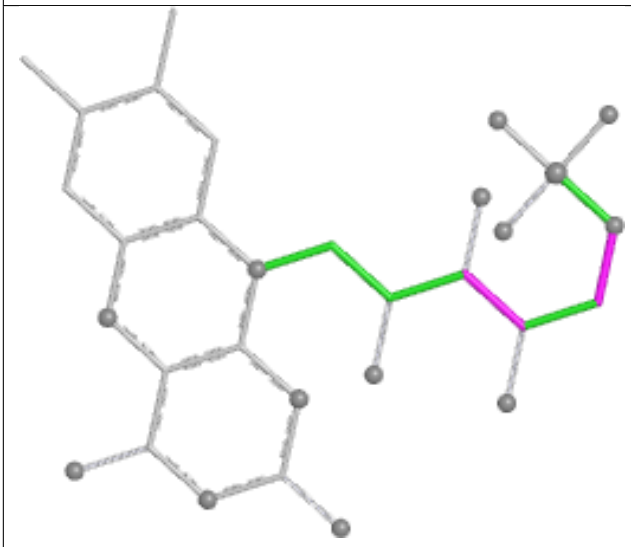
Ligand FMN B 402



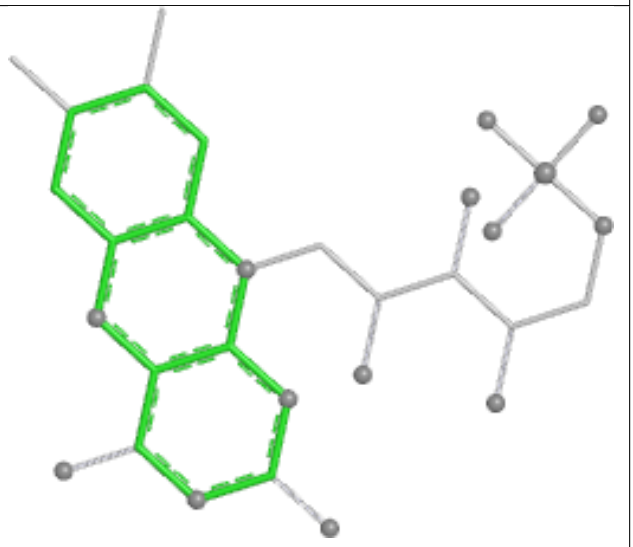
Bond lengths



Bond angles

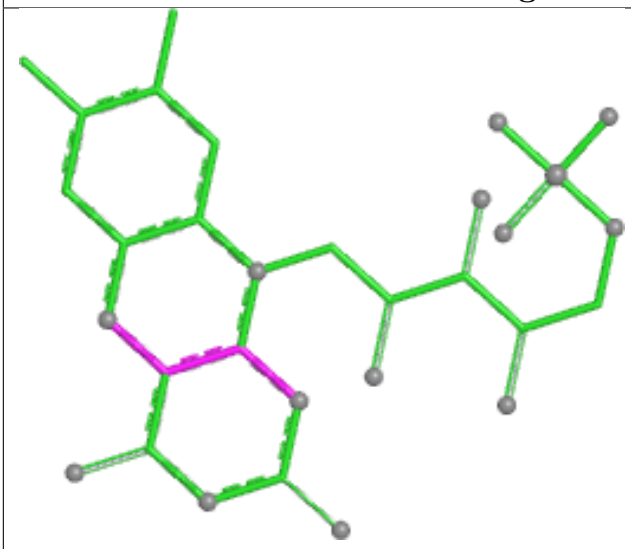


Torsions

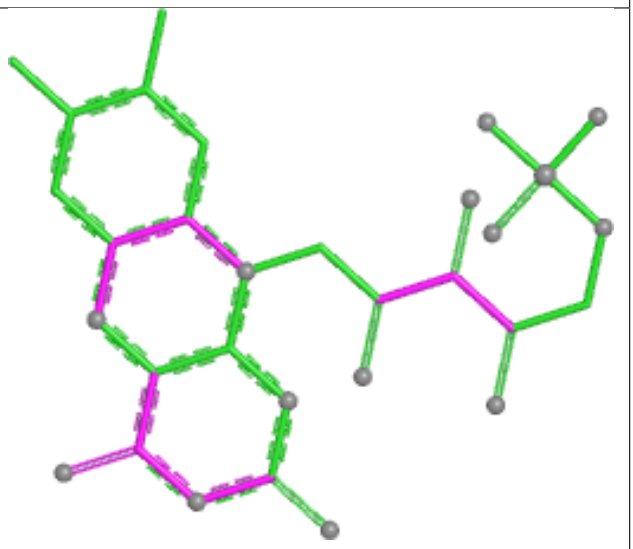


Rings

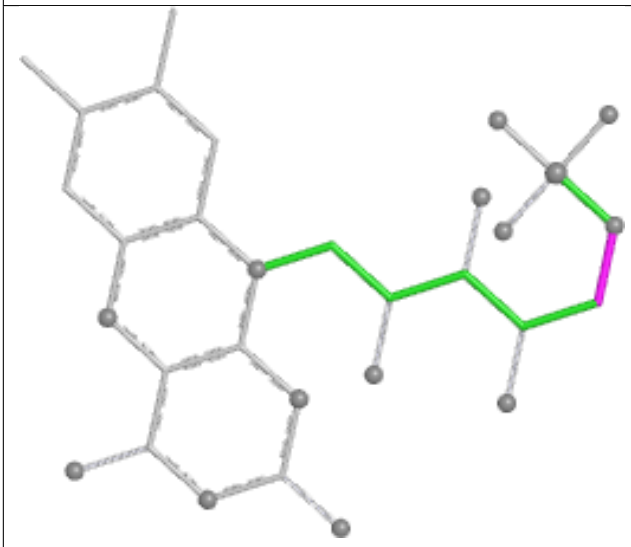
Ligand FMN A 402



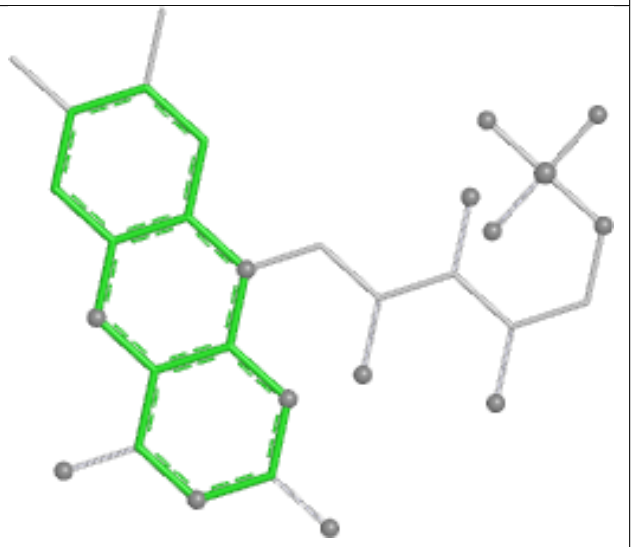
Bond lengths



Bond angles

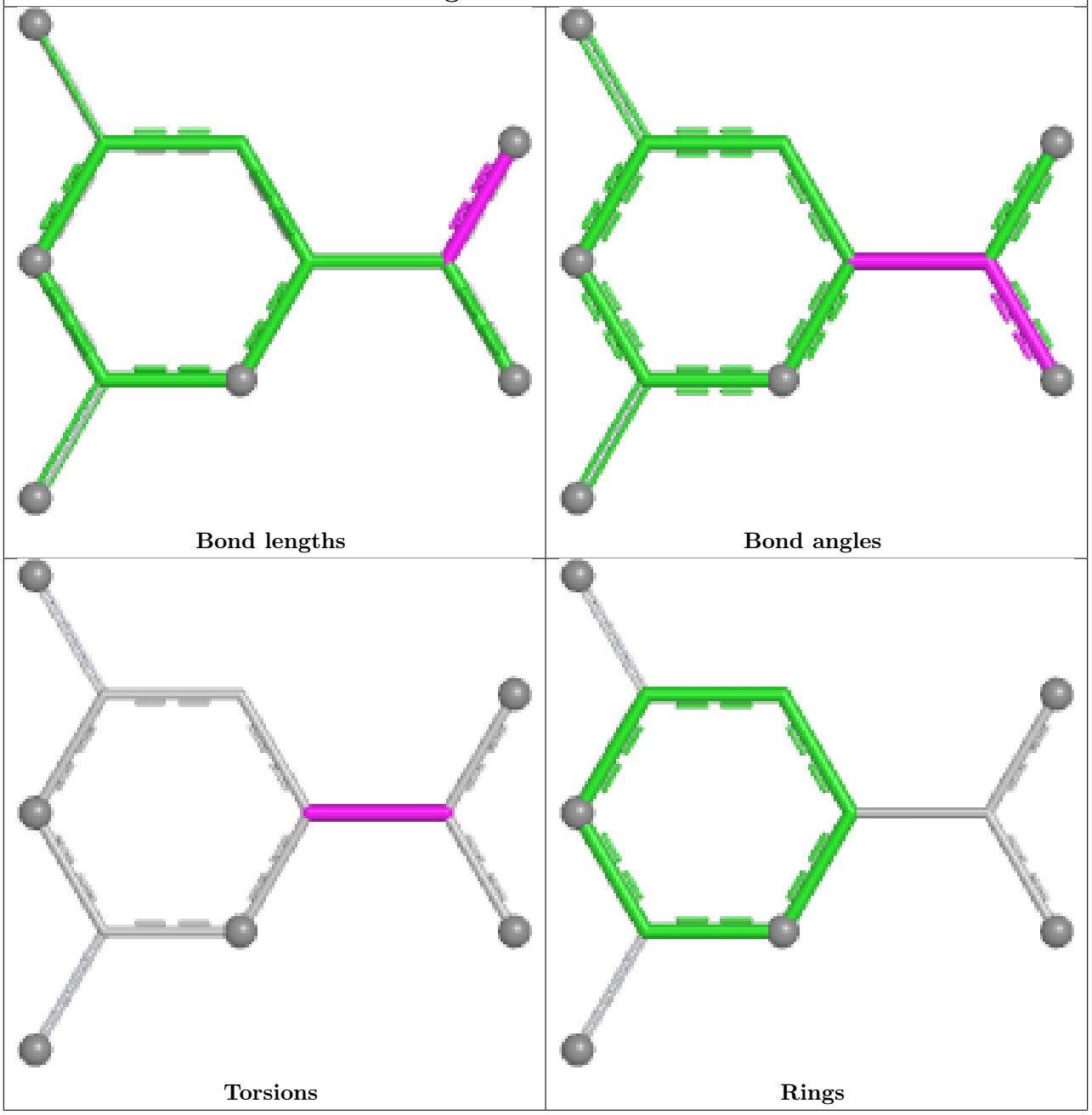


Torsions



Rings

Ligand ORO B 401



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	335/368 (91%)	0.07	12 (3%) 46 42	30, 52, 70, 116	2 (0%)
1	B	335/368 (91%)	0.12	5 (1%) 72 69	42, 58, 75, 89	0
All	All	670/736 (91%)	0.09	17 (2%) 58 55	30, 55, 74, 116	2 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	34	PHE	9.3
1	A	33	PRO	6.2
1	A	32	THR	5.7
1	A	37	LEU	5.0
1	B	2	TYR	4.6
1	A	31	GLY	3.1
1	B	63	GLY	3.0
1	A	2	TYR	2.8
1	B	5	PHE	2.7
1	A	35	GLU	2.6
1	A	30	THR	2.4
1	A	29	ILE	2.2
1	A	20	GLU	2.1
1	A	39	ARG	2.1
1	A	38	VAL	2.1
1	B	21	PHE	2.0
1	B	4	PRO	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)

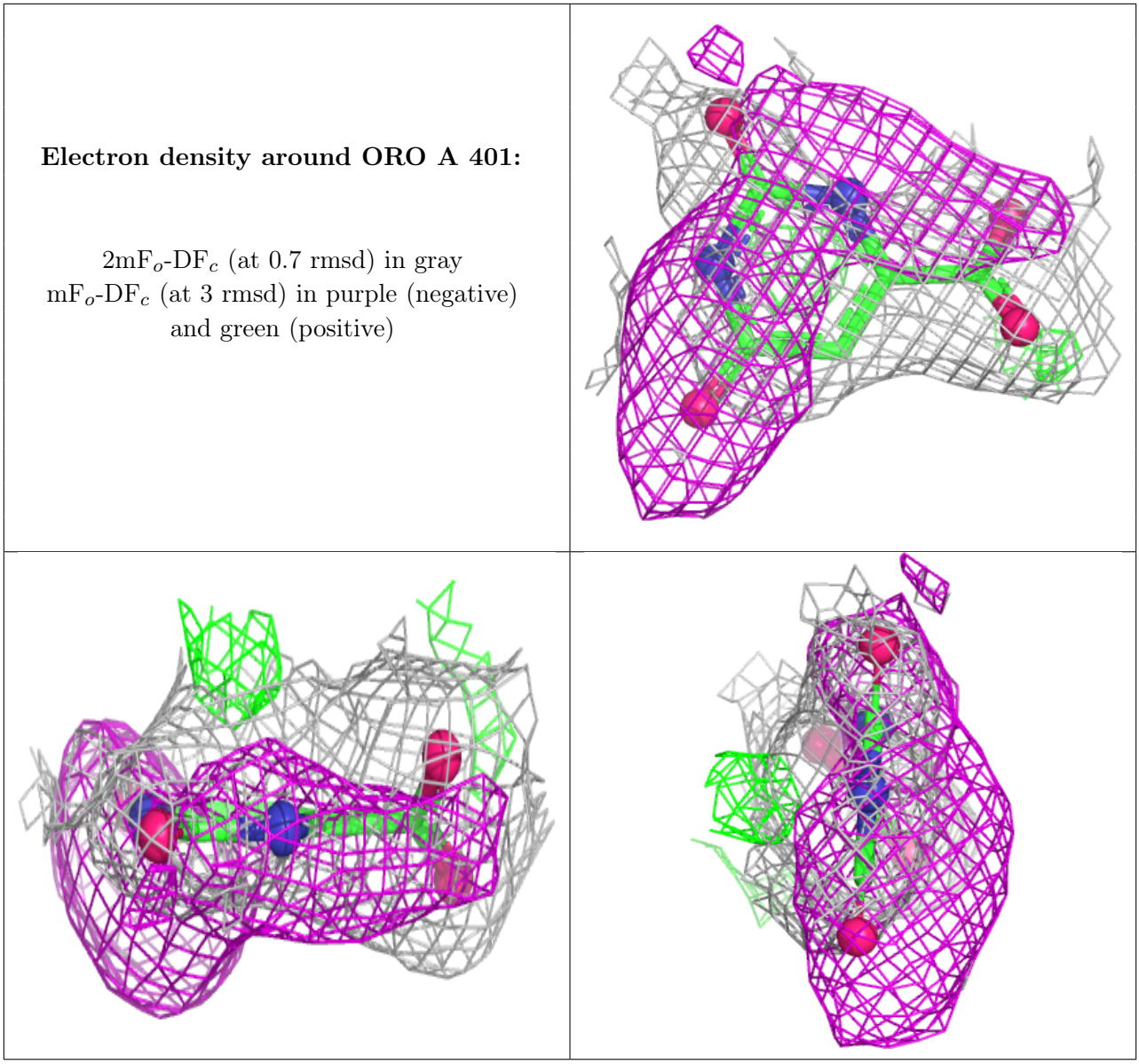
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ORO	A	401	11/11	0.78	0.17	57,59,64,66	0
2	ORO	B	401	11/11	0.86	0.12	65,67,70,74	0
3	FMN	B	402	31/31	0.96	0.08	42,54,59,61	0
3	FMN	A	402	31/31	0.97	0.07	39,47,51,52	0

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

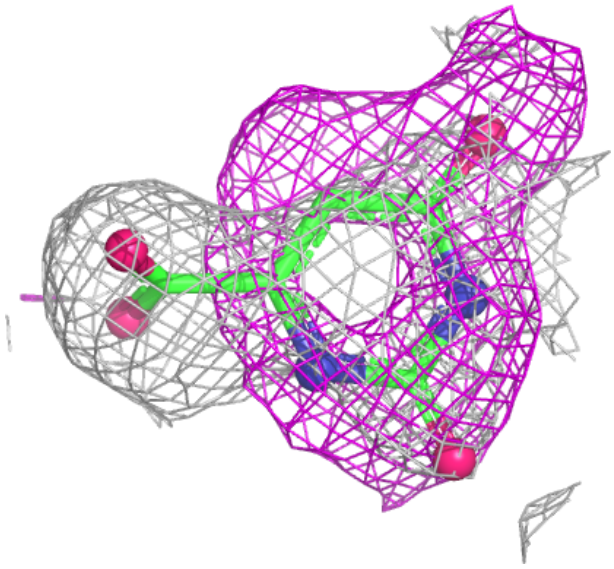
Electron density around ORO A 401:

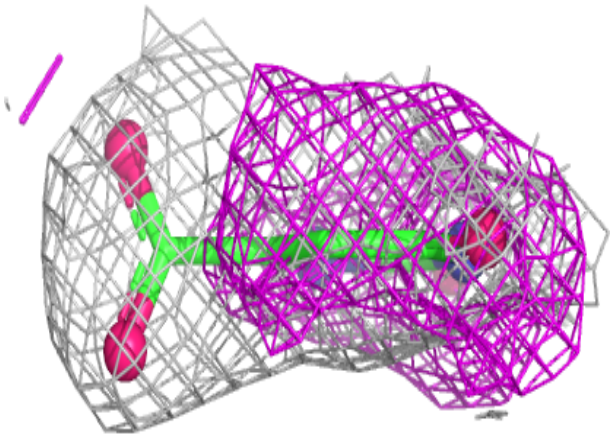
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

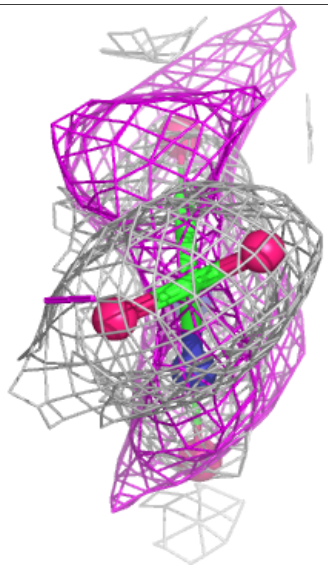


Electron density around ORO B 401:

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

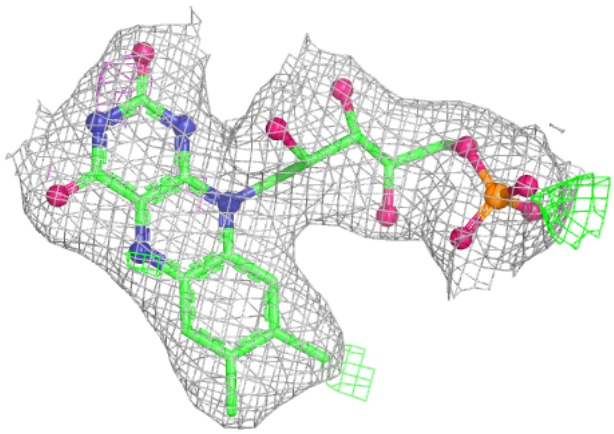


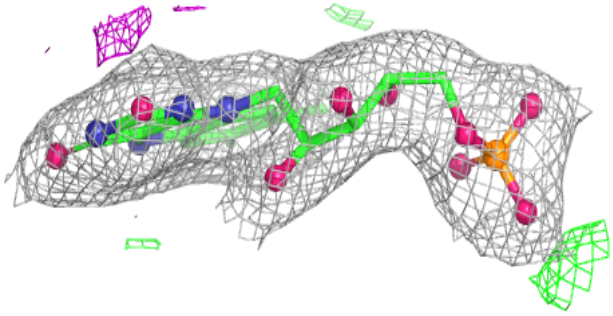


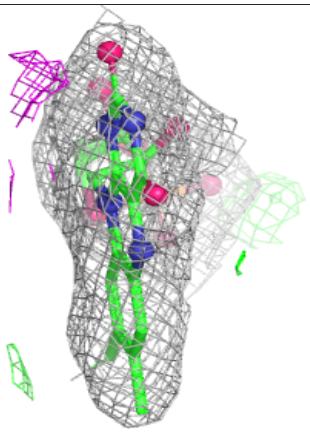


Electron density around FMN B 402:

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

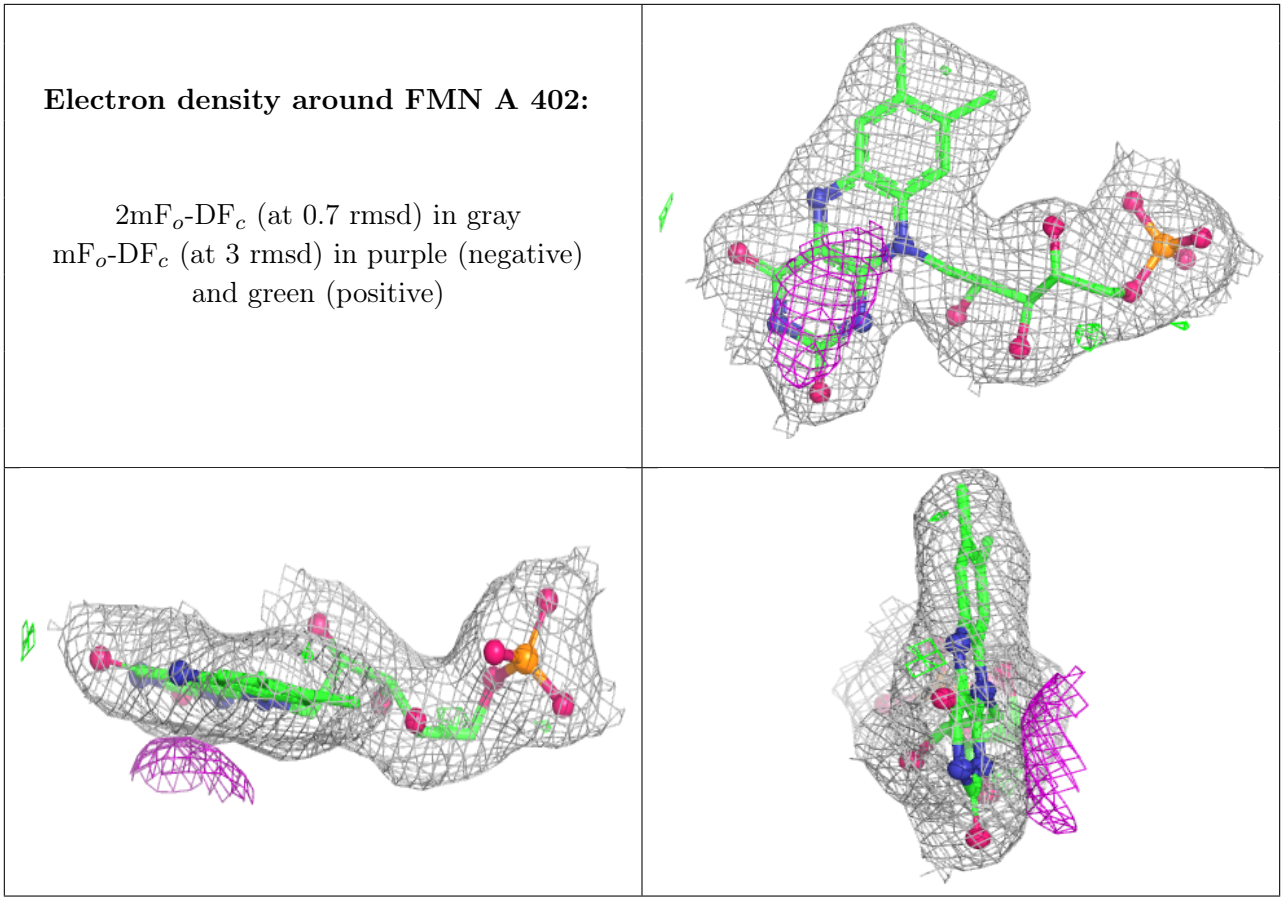






Electron density around FMN A 402:

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



6.5 Other polymers

There are no such residues in this entry.