



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:09 PM UTC

PDB ID : 3AH5 / pdb_00003ah5
Title : Crystal Structure of flavin dependent thymidylate synthase ThyX from helicobacter pylori complexed with FAD and dUMP
Authors : Zhang, X.; Zhang, J.; Hu, Y.; Zou, Q.; Wang, D.
Deposited on : 2010-04-14
Resolution : 2.50 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:

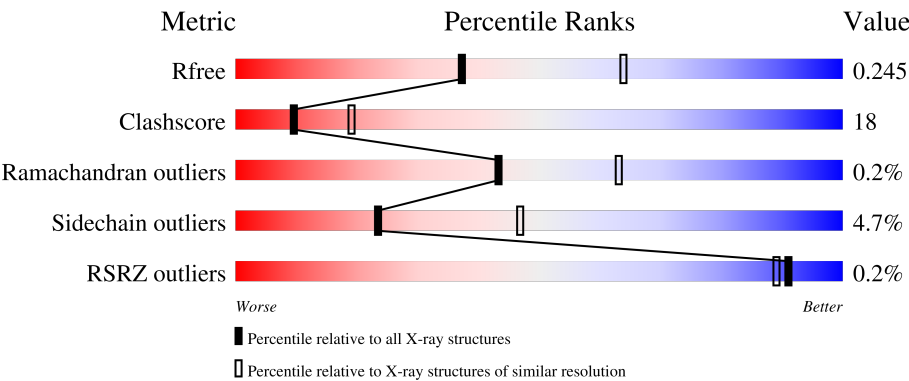
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal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

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

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	216	66%30%.
1	B	216	65%30%..
1	C	216	66%27%..
1	D	216	% 65%28%..
1	E	216	64%30%..

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Mol	Chain	Length	Quality of chain
1	F	216	 64% 28% . .

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	UMP	C	241	-	-	X	-
4	SO4	E	2	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11206 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 Thymidylate synthase thyX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1760	1113	313	326	8			
1	B	207	Total	C	N	O	S	0	0	0
			1683	1066	293	316	8			
1	C	207	Total	C	N	O	S	0	0	0
			1683	1066	293	316	8			
1	D	207	Total	C	N	O	S	0	0	0
			1683	1066	293	316	8			
1	E	207	Total	C	N	O	S	0	0	0
			1683	1066	293	316	8			
1	F	207	Total	C	N	O	S	0	0	0
			1683	1066	293	316	8			

There are 48 discrepancies between the modelled and reference sequences:

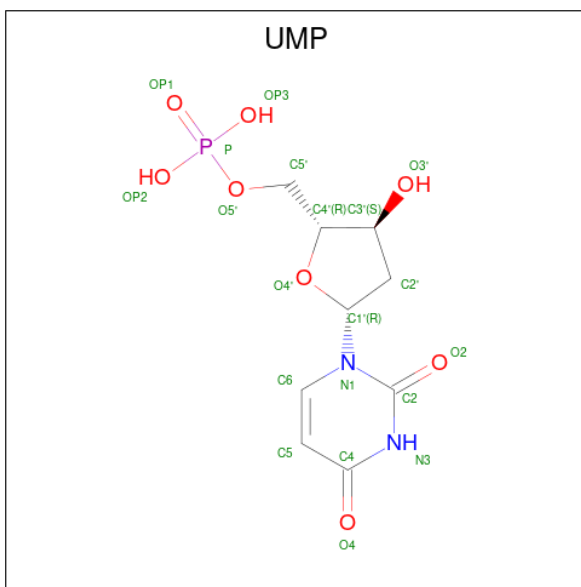
Chain	Residue	Modelled	Actual	Comment	Reference
A	232	LEU	-	expression tag	UNP O26061
A	233	GLU	-	expression tag	UNP O26061
A	234	HIS	-	expression tag	UNP O26061
A	235	HIS	-	expression tag	UNP O26061
A	236	HIS	-	expression tag	UNP O26061
A	237	HIS	-	expression tag	UNP O26061
A	238	HIS	-	expression tag	UNP O26061
A	239	HIS	-	expression tag	UNP O26061
B	232	LEU	-	expression tag	UNP O26061
B	233	GLU	-	expression tag	UNP O26061
B	234	HIS	-	expression tag	UNP O26061
B	235	HIS	-	expression tag	UNP O26061
B	236	HIS	-	expression tag	UNP O26061
B	237	HIS	-	expression tag	UNP O26061
B	238	HIS	-	expression tag	UNP O26061
B	239	HIS	-	expression tag	UNP O26061
C	232	LEU	-	expression tag	UNP O26061

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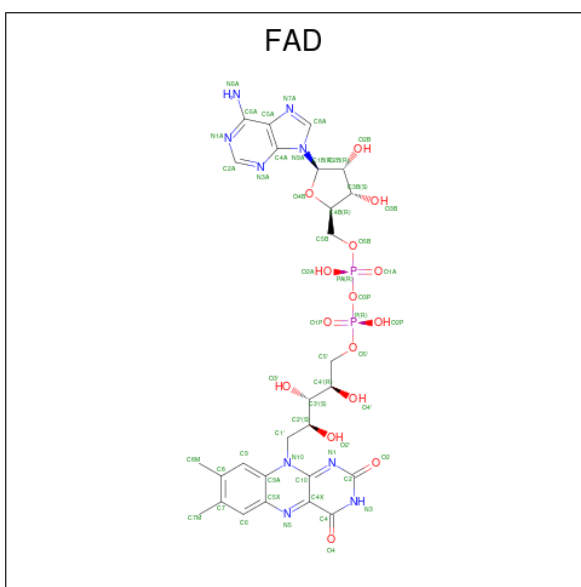
Chain	Residue	Modelled	Actual	Comment	Reference
C	233	GLU	-	expression tag	UNP O26061
C	234	HIS	-	expression tag	UNP O26061
C	235	HIS	-	expression tag	UNP O26061
C	236	HIS	-	expression tag	UNP O26061
C	237	HIS	-	expression tag	UNP O26061
C	238	HIS	-	expression tag	UNP O26061
C	239	HIS	-	expression tag	UNP O26061
D	232	LEU	-	expression tag	UNP O26061
D	233	GLU	-	expression tag	UNP O26061
D	234	HIS	-	expression tag	UNP O26061
D	235	HIS	-	expression tag	UNP O26061
D	236	HIS	-	expression tag	UNP O26061
D	237	HIS	-	expression tag	UNP O26061
D	238	HIS	-	expression tag	UNP O26061
D	239	HIS	-	expression tag	UNP O26061
E	232	LEU	-	expression tag	UNP O26061
E	233	GLU	-	expression tag	UNP O26061
E	234	HIS	-	expression tag	UNP O26061
E	235	HIS	-	expression tag	UNP O26061
E	236	HIS	-	expression tag	UNP O26061
E	237	HIS	-	expression tag	UNP O26061
E	238	HIS	-	expression tag	UNP O26061
E	239	HIS	-	expression tag	UNP O26061
F	232	LEU	-	expression tag	UNP O26061
F	233	GLU	-	expression tag	UNP O26061
F	234	HIS	-	expression tag	UNP O26061
F	235	HIS	-	expression tag	UNP O26061
F	236	HIS	-	expression tag	UNP O26061
F	237	HIS	-	expression tag	UNP O26061
F	238	HIS	-	expression tag	UNP O26061
F	239	HIS	-	expression tag	UNP O26061

- Molecule 2 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (CCD ID: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	B	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	C	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	D	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	E	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	F	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	B	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	C	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	D	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	E	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	F	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	98	Total	O	0	0
			98	98		

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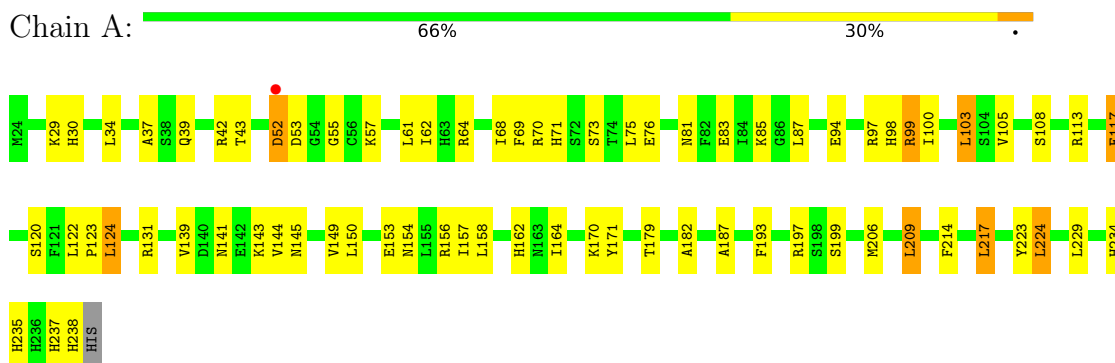
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	98	Total 98	O 98	0	0
5	C	112	Total 112	O 112	0	0
5	D	84	Total 84	O 84	0	0
5	E	74	Total 74	O 74	0	0
5	F	72	Total 72	O 72	0	0

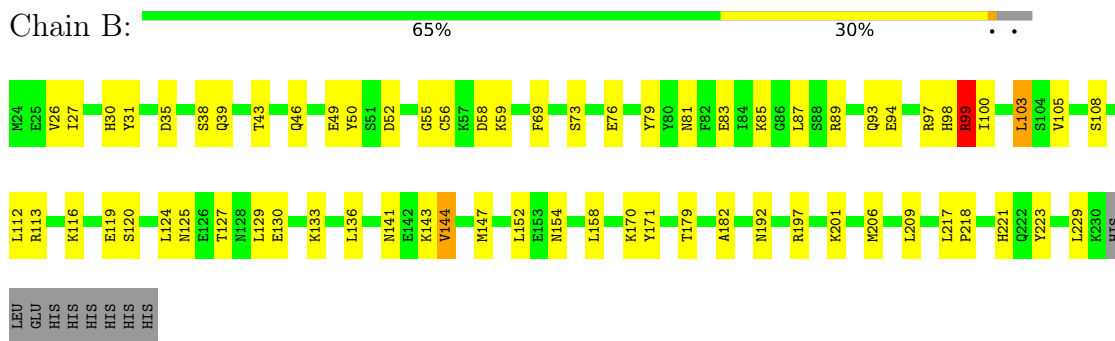
3 Residue-property plots [i](#)

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.

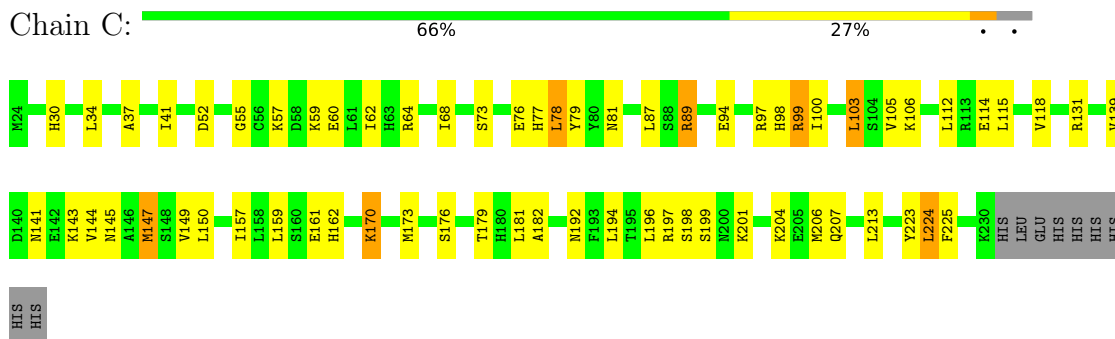
• Molecule 1: Thymidylate synthase thyX



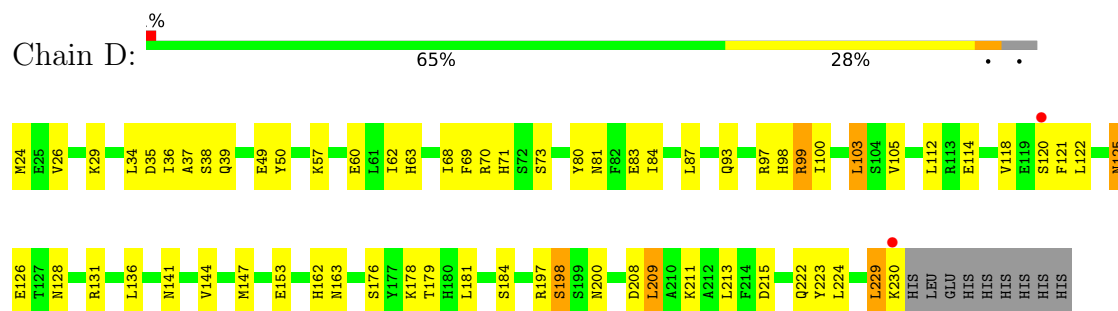
• Molecule 1: Thymidylate synthase thyX



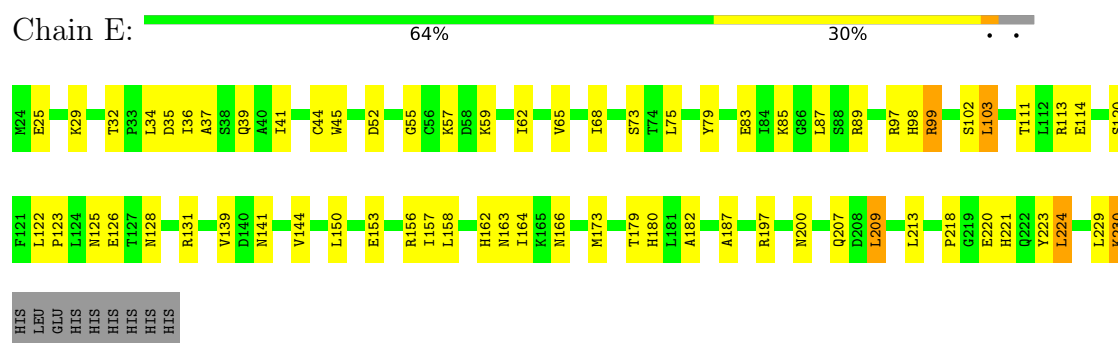
• Molecule 1: Thymidylate synthase thyX



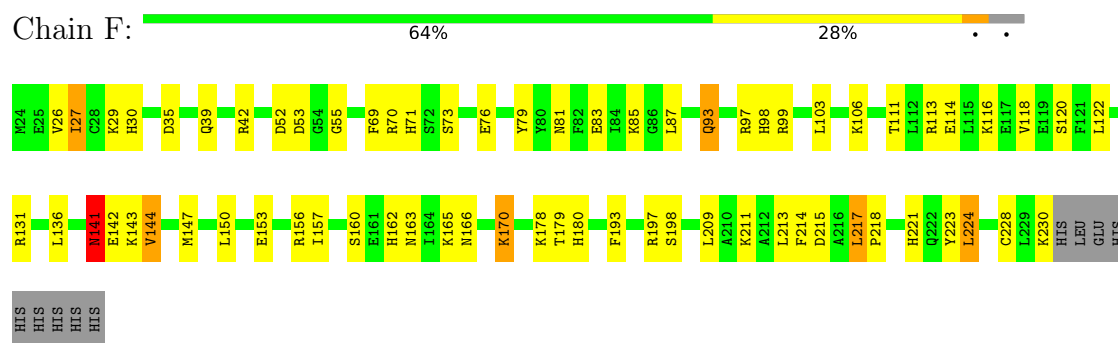
- Molecule 1: Thymidylate synthase thyX



- Molecule 1: Thymidylate synthase thyX



- Molecule 1: Thymidylate synthase thyX



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	221.92Å 49.43Å 143.02Å 90.00° 98.84° 90.00°	Depositor
Resolution (Å)	55.63 – 2.50 55.63 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (55.63-2.50) 99.7 (55.63-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.50Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.194 , 0.246 0.194 , 0.245	Depositor DCC
R_{free} test set	5207 reflections (9.68%)	wwPDB-VP
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11206	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, FAD, UMP

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	0.37	0/1797	0.89	3/2420 (0.1%)
1	B	0.38	0/1714	0.89	4/2307 (0.2%)
1	C	0.39	0/1714	0.89	5/2307 (0.2%)
1	D	0.36	0/1714	0.90	7/2307 (0.3%)
1	E	0.36	0/1714	0.86	5/2307 (0.2%)
1	F	0.35	0/1714	0.86	2/2307 (0.1%)
All	All	0.37	0/10367	0.88	26/13955 (0.2%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	126	GLU	N-CA-C	-7.98	102.45	112.90
1	D	105	VAL	N-CA-C	6.53	117.99	108.53
1	D	198	SER	N-CA-C	-6.43	105.40	113.18
1	D	36	ILE	N-CA-C	-6.40	104.26	110.72
1	C	99	ARG	N-CA-C	6.39	119.24	111.82
1	B	99	ARG	N-CA-C	6.33	118.99	111.71
1	B	69	PHE	N-CA-C	6.10	118.71	111.33
1	F	99	ARG	N-CA-C	5.98	118.58	111.71
1	A	235	HIS	N-CA-C	5.86	119.96	112.34
1	C	198	SER	N-CA-C	-5.69	106.33	113.28
1	A	105	VAL	N-CA-C	5.64	116.71	108.53
1	B	182	ALA	N-CA-C	-5.61	100.10	109.24
1	C	182	ALA	N-CA-C	-5.58	100.62	109.76
1	E	99	ARG	N-CA-C	5.56	119.16	112.38
1	D	121	PHE	N-CA-C	-5.49	105.80	112.88
1	A	182	ALA	N-CA-C	-5.42	100.41	109.24
1	D	34	LEU	N-CA-C	5.41	117.61	111.11
1	C	105	VAL	N-CA-C	5.41	117.11	108.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	207	GLN	N-CA-C	-5.38	105.49	111.36
1	C	68	ILE	N-CA-C	-5.27	100.03	107.98
1	F	69	PHE	N-CA-C	5.21	119.97	111.37
1	E	68	ILE	N-CA-C	-5.20	100.28	108.23
1	D	229	LEU	N-CA-C	5.19	118.20	110.52
1	E	182	ALA	N-CA-C	-5.17	101.27	109.76
1	D	222	GLN	N-CA-C	5.16	117.64	111.71
1	B	105	VAL	N-CA-C	5.15	116.00	108.53

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	1760	0	1740	88	0
1	B	1683	0	1681	61	0
1	C	1683	0	1681	75	0
1	D	1683	0	1681	67	0
1	E	1683	0	1681	56	0
1	F	1683	0	1681	62	0
2	A	20	0	11	6	0
2	B	20	0	11	6	0
2	C	20	0	11	7	0
2	D	20	0	11	6	0
2	E	20	0	11	4	0
2	F	20	0	11	6	0
3	A	53	0	31	8	0
3	B	53	0	31	7	0
3	C	53	0	31	7	0
3	D	53	0	31	8	0
3	E	53	0	31	2	0
3	F	53	0	31	3	0
4	A	15	0	0	0	0
4	C	15	0	0	0	0
4	D	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	10	0	0	4	0
4	F	5	0	0	0	0
5	A	98	0	0	7	0
5	B	98	0	0	7	0
5	C	112	0	0	1	0
5	D	84	0	0	3	0
5	E	74	0	0	1	1
5	F	72	0	0	2	1
All	All	11206	0	10397	375	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 18.

All (375) 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:E:97:ARG:HG3	2:E:241:UMP:H2''	1.38	1.02
2:A:241:UMP:H2''	1:C:97:ARG:HG3	1.43	1.01
1:B:97:ARG:HG3	2:D:241:UMP:H2''	1.41	1.01
1:A:197:ARG:HH12	2:C:241:UMP:HN3	0.97	0.97
1:B:197:ARG:HH12	2:D:241:UMP:HN3	1.09	0.95
1:D:125:ASN:ND2	1:D:128:ASN:H	1.67	0.93
2:B:241:UMP:H2''	1:D:97:ARG:HG3	1.51	0.93
1:F:197:ARG:HH12	2:F:241:UMP:HN3	1.12	0.92
1:E:197:ARG:HH12	2:E:241:UMP:HN3	1.12	0.91
1:A:30:HIS:HB3	1:A:81:ASN:HD22	1.34	0.91
1:A:97:ARG:HG3	2:C:241:UMP:H2''	1.51	0.90
2:B:241:UMP:HN3	1:D:197:ARG:HH12	1.20	0.89
1:C:99:ARG:HG3	1:C:99:ARG:HH11	1.37	0.88
1:F:97:ARG:HG3	2:F:241:UMP:H2''	1.53	0.87
1:F:98:HIS:HE1	1:F:197:ARG:HE	1.18	0.87
1:A:154:ASN:HB3	1:C:147:MET:HG3	1.58	0.86
1:F:29:LYS:HD2	1:F:83:GLU:HB2	1.58	0.85
1:B:98:HIS:HE1	1:B:197:ARG:HE	1.22	0.85
1:C:98:HIS:HE1	1:C:197:ARG:HE	1.26	0.84
2:A:241:UMP:HN3	1:C:197:ARG:HH12	1.27	0.83
1:F:141:ASN:HD21	1:F:143:LYS:HB2	1.44	0.82
1:A:69:PHE:HZ	1:B:192:ASN:HD21	1.29	0.79
3:F:240:FAD:H51A	3:F:240:FAD:H8A	1.66	0.78
3:B:240:FAD:H51A	3:B:240:FAD:H8A	1.64	0.78
1:B:30:HIS:HB3	1:B:81:ASN:HD22	1.48	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:HIS:HE1	1:D:197:ARG:HE	1.32	0.78
3:D:240:FAD:H51A	3:D:240:FAD:H8A	1.66	0.77
3:E:240:FAD:H51A	3:E:240:FAD:H8A	1.67	0.77
5:B:535:HOH:O	1:C:77:HIS:HA	1.84	0.77
1:D:125:ASN:C	1:D:125:ASN:HD22	1.94	0.76
3:C:240:FAD:H8A	3:C:240:FAD:H51A	1.67	0.76
1:B:46:GLN:HG3	1:C:106:LYS:HE2	1.67	0.75
1:E:29:LYS:HE2	1:E:83:GLU:HB2	1.69	0.75
1:A:99:ARG:O	1:C:99:ARG:HD2	1.87	0.74
1:F:27:ILE:CD1	1:F:83:GLU:HB3	2.18	0.73
1:A:141:ASN:O	1:A:144:VAL:HG22	1.89	0.73
1:A:197:ARG:NH1	2:C:241:UMP:HN3	1.80	0.73
1:E:35:ASP:O	1:E:39:GLN:HG3	1.88	0.73
1:F:27:ILE:HD11	1:F:83:GLU:HB3	1.71	0.72
1:A:97:ARG:CG	2:C:241:UMP:H2''	2.19	0.72
1:A:237:HIS:O	1:A:238:HIS:HB2	1.88	0.72
1:B:99:ARG:HB3	1:D:99:ARG:HB2	1.71	0.72
1:F:98:HIS:CE1	1:F:197:ARG:HE	2.05	0.72
1:A:29:LYS:HE3	1:D:39:GLN:OE1	1.90	0.71
1:C:34:LEU:HD12	1:C:59:LYS:HE3	1.73	0.70
1:D:100:ILE:HD12	3:D:240:FAD:H52A	1.72	0.70
1:E:123:PRO:HA	4:E:2:SO4:O1	1.91	0.70
3:A:240:FAD:H51A	3:A:240:FAD:H8A	1.72	0.70
1:E:98:HIS:HE1	1:E:197:ARG:HE	1.38	0.70
1:C:147:MET:HE3	1:C:147:MET:HA	1.72	0.70
1:D:99:ARG:HG2	3:D:240:FAD:O2A	1.91	0.70
1:A:154:ASN:HB3	1:C:147:MET:CG	2.23	0.69
1:B:100:ILE:HD12	3:B:240:FAD:H52A	1.75	0.69
1:C:99:ARG:HG3	1:C:99:ARG:NH1	2.08	0.68
1:E:120:SER:HA	1:E:156:ARG:HE	1.58	0.68
1:D:125:ASN:HD21	1:D:128:ASN:H	1.41	0.68
1:A:83:GLU:OE2	1:A:85:LYS:HE3	1.94	0.68
1:A:158:LEU:HD13	1:C:144:VAL:HG12	1.75	0.68
1:C:98:HIS:CE1	1:C:197:ARG:HE	2.10	0.68
1:F:223:TYR:CE1	1:F:224:LEU:HD13	2.29	0.68
1:A:39:GLN:NE2	1:D:29:LYS:HE3	2.09	0.67
1:A:98:HIS:HE1	1:A:197:ARG:HE	1.40	0.67
1:D:100:ILE:CD1	3:D:240:FAD:H52A	2.24	0.67
1:B:49:GLU:HG3	1:B:50:TYR:CD1	2.29	0.67
1:E:83:GLU:OE1	1:E:85:LYS:HE3	1.93	0.67
1:A:29:LYS:HD2	1:A:83:GLU:HB2	1.74	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:HIS:CE1	1:B:197:ARG:HE	2.10	0.66
2:A:241:UMP:H2''	1:C:97:ARG:CG	2.23	0.66
1:A:214:PHE:O	1:A:217:LEU:HB2	1.95	0.66
1:A:69:PHE:HZ	1:B:192:ASN:ND2	1.94	0.66
1:D:38:SER:N	1:D:62:ILE:HD11	2.11	0.65
1:A:100:ILE:HD12	3:A:240:FAD:H52A	1.79	0.65
1:A:103:LEU:N	1:A:103:LEU:HD23	2.12	0.65
1:B:100:ILE:CD1	3:B:240:FAD:H52A	2.27	0.65
1:A:52:ASP:HB3	1:A:55:GLY:H	1.61	0.64
1:C:37:ALA:HB2	1:C:78:LEU:HD13	1.78	0.64
1:C:100:ILE:O	1:C:100:ILE:HG22	1.97	0.64
2:B:241:UMP:H2''	1:D:97:ARG:CG	2.27	0.64
1:D:29:LYS:HD2	1:D:83:GLU:HB2	1.79	0.64
1:D:98:HIS:CE1	1:D:197:ARG:HE	2.16	0.63
1:F:35:ASP:O	1:F:39:GLN:HG3	1.98	0.63
1:B:35:ASP:O	1:B:39:GLN:HG3	1.99	0.63
1:E:220:GLU:HG3	1:E:221:HIS:HD2	1.63	0.63
1:F:166:ASN:O	1:F:170:LYS:HG3	1.99	0.63
1:C:76:GLU:OE1	1:D:99:ARG:NH2	2.32	0.62
1:B:97:ARG:CG	2:D:241:UMP:H2''	2.23	0.62
1:F:153:GLU:O	1:F:157:ILE:HG12	1.99	0.62
1:A:99:ARG:HG2	3:A:240:FAD:O2A	2.00	0.61
1:E:32:THR:HG23	1:E:36:ILE:HD12	1.81	0.61
1:C:100:ILE:HD13	3:C:240:FAD:H52A	1.82	0.61
1:C:73:SER:HB3	3:D:240:FAD:O4'	1.99	0.61
1:E:120:SER:HB2	1:E:156:ARG:HH21	1.66	0.61
1:A:120:SER:HA	1:A:156:ARG:HE	1.66	0.61
1:B:83:GLU:OE2	1:B:85:LYS:HE3	2.01	0.61
1:A:154:ASN:CG	1:C:147:MET:HE2	2.25	0.60
1:A:99:ARG:HD3	5:A:6:HOH:O	2.00	0.60
1:B:49:GLU:HG3	1:B:50:TYR:HD1	1.65	0.60
1:E:97:ARG:CG	2:E:241:UMP:H2''	2.23	0.60
1:A:139:VAL:CG2	1:A:144:VAL:HG21	2.31	0.60
1:A:87:LEU:O	1:A:179:THR:HG22	2.02	0.59
1:F:141:ASN:C	1:F:141:ASN:HD22	2.10	0.59
1:A:76:GLU:OE1	1:B:99:ARG:NH2	2.33	0.59
1:E:230:LYS:HE3	1:E:230:LYS:N	2.18	0.59
1:C:89:ARG:HH22	2:C:241:UMP:H5''	1.67	0.59
1:D:24:MET:HE1	1:D:208:ASP:C	2.27	0.59
1:C:89:ARG:NH2	2:C:241:UMP:OP3	2.35	0.59
1:A:113:ARG:HG3	5:A:412:HOH:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:241:UMP:C2'	1:C:97:ARG:HG3	2.26	0.59
1:F:106:LYS:HG2	1:F:180:HIS:H	1.68	0.59
1:C:141:ASN:O	1:C:144:VAL:HG22	2.03	0.58
3:A:240:FAD:O4'	1:B:73:SER:HB3	2.03	0.58
1:F:98:HIS:HE1	1:F:197:ARG:NE	1.96	0.58
1:A:98:HIS:CE1	1:A:197:ARG:HE	2.21	0.58
1:A:113:ARG:HB2	5:A:254:HOH:O	2.03	0.58
1:E:122:LEU:HB3	4:E:2:SO4:O1	2.04	0.58
1:B:100:ILE:CG1	3:B:240:FAD:H52A	2.34	0.57
1:C:201:LYS:HE3	5:D:459:HOH:O	2.04	0.57
1:B:97:ARG:HG3	2:D:241:UMP:C2'	2.24	0.57
1:A:52:ASP:OD2	1:A:57:LYS:HB2	2.05	0.57
1:C:118:VAL:HG22	1:C:131:ARG:HH12	1.70	0.57
1:F:214:PHE:O	1:F:217:LEU:HB2	2.04	0.57
1:A:158:LEU:HD22	1:A:164:ILE:HD12	1.87	0.56
1:D:100:ILE:CG1	3:D:240:FAD:H52A	2.34	0.56
1:D:122:LEU:CD1	1:D:153:GLU:HG2	2.35	0.56
1:C:223:TYR:CZ	1:C:224:LEU:HD13	2.40	0.56
1:D:229:LEU:O	1:D:230:LYS:HB2	2.04	0.56
1:F:97:ARG:CG	2:F:241:UMP:H2''	2.30	0.56
1:E:98:HIS:CE1	1:E:197:ARG:HE	2.20	0.56
1:E:103:LEU:HD23	1:E:103:LEU:N	2.19	0.56
1:F:73:SER:HB3	3:F:240:FAD:O4'	2.06	0.56
1:A:162:HIS:CD2	1:C:143:LYS:HE2	2.40	0.56
1:F:122:LEU:HD11	1:F:153:GLU:HG2	1.89	0.55
1:A:122:LEU:HD23	1:A:153:GLU:HG2	1.87	0.55
1:B:38:SER:OG	1:B:58:ASP:HB3	2.07	0.55
1:D:141:ASN:O	1:D:144:VAL:HG22	2.07	0.55
1:C:98:HIS:HE1	1:C:197:ARG:NE	2.00	0.55
1:B:100:ILE:HG13	3:B:240:FAD:H52A	1.89	0.54
1:E:218:PRO:HA	4:E:1:SO4:O1	2.08	0.54
1:C:76:GLU:CD	1:D:99:ARG:HH22	2.15	0.54
3:C:240:FAD:O4'	1:D:73:SER:HB3	2.08	0.54
1:E:141:ASN:HB3	1:E:144:VAL:HG22	1.90	0.54
1:E:111:THR:O	1:E:114:GLU:HG2	2.08	0.54
1:E:218:PRO:HG2	1:E:221:HIS:HB2	1.90	0.53
1:D:100:ILE:HG13	3:D:240:FAD:H52A	1.90	0.53
1:A:143:LYS:HD3	1:C:162:HIS:CE1	2.44	0.53
5:B:535:HOH:O	1:C:77:HIS:CA	2.52	0.53
1:F:113:ARG:HH21	1:F:116:LYS:NZ	2.06	0.53
1:A:153:GLU:O	1:A:157:ILE:HG13	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:LYS:HD3	1:C:162:HIS:ND1	2.23	0.53
1:C:30:HIS:HB3	1:C:81:ASN:HD22	1.74	0.53
1:E:153:GLU:O	1:E:157:ILE:HG13	2.08	0.53
1:F:27:ILE:HD13	1:F:83:GLU:HB3	1.90	0.53
1:A:97:ARG:HG3	2:C:241:UMP:C2'	2.34	0.53
1:A:100:ILE:CD1	3:A:240:FAD:H52A	2.38	0.53
1:C:145:ASN:O	1:C:149:VAL:HG23	2.08	0.53
1:D:211:LYS:HE3	1:D:215:ASP:OD2	2.09	0.53
1:F:70:ARG:NH1	1:F:71:HIS:CE1	2.77	0.52
1:A:144:VAL:HG23	1:A:145:ASN:N	2.24	0.52
1:C:141:ASN:HB3	1:C:144:VAL:HG22	1.92	0.52
1:F:116:LYS:HA	1:F:160:SER:OG	2.10	0.52
1:B:87:LEU:O	1:B:179:THR:HG22	2.09	0.52
1:E:158:LEU:HD22	1:E:164:ILE:HD12	1.91	0.52
2:B:241:UMP:C2'	1:D:97:ARG:HG3	2.32	0.52
1:A:39:GLN:HE22	1:D:29:LYS:HE3	1.73	0.52
1:C:192:ASN:HD21	1:D:69:PHE:HZ	1.57	0.52
1:D:87:LEU:O	1:D:179:THR:HG22	2.09	0.52
1:A:100:ILE:CG1	3:A:240:FAD:H52A	2.40	0.52
1:D:70:ARG:NH1	1:D:71:HIS:HE1	2.08	0.52
1:E:120:SER:O	1:E:131:ARG:NH2	2.43	0.52
1:A:113:ARG:HB3	5:A:280:HOH:O	2.10	0.52
1:A:71:HIS:HE1	5:A:269:HOH:O	1.93	0.51
1:B:56:CYS:HB3	5:B:514:HOH:O	2.10	0.51
1:B:127:THR:HG22	5:B:428:HOH:O	2.10	0.51
1:E:52:ASP:OD1	1:E:57:LYS:HB2	2.10	0.51
1:F:118:VAL:HG23	1:F:156:ARG:HD2	1.92	0.51
1:A:122:LEU:HD13	1:A:124:LEU:HD13	1.93	0.51
1:D:103:LEU:N	1:D:103:LEU:HD23	2.25	0.51
1:E:32:THR:CG2	1:E:36:ILE:HD12	2.40	0.51
1:C:196:LEU:HD21	1:D:69:PHE:CE2	2.46	0.51
1:F:141:ASN:HB3	1:F:144:VAL:HG13	1.93	0.51
1:D:125:ASN:HD22	1:D:128:ASN:H	1.54	0.51
1:C:89:ARG:HD2	1:C:173:MET:HE3	1.93	0.51
2:B:241:UMP:H2''	1:D:97:ARG:HH11	1.74	0.51
1:F:113:ARG:HH21	1:F:116:LYS:HZ3	1.57	0.51
1:B:129:LEU:O	1:B:133:LYS:HG3	2.11	0.51
1:B:143:LYS:HE2	1:D:162:HIS:CD2	2.45	0.51
1:A:69:PHE:CZ	1:B:192:ASN:ND2	2.72	0.50
1:A:70:ARG:NH2	1:B:201:LYS:HE3	2.25	0.50
1:B:43:THR:HG21	1:C:181:LEU:HA	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:LEU:CD1	1:A:124:LEU:HD13	2.42	0.50
1:B:108:SER:N	2:B:241:UMP:OP2	2.45	0.50
1:D:125:ASN:ND2	1:D:125:ASN:C	2.66	0.50
1:E:102:SER:OG	1:F:76:GLU:HB3	2.12	0.49
1:E:55:GLY:O	1:E:59:LYS:HG3	2.11	0.49
1:C:100:ILE:HD13	3:C:240:FAD:C5B	2.43	0.49
3:B:240:FAD:O3P	3:B:240:FAD:H3B	2.12	0.49
1:A:52:ASP:HB3	1:A:55:GLY:N	2.26	0.49
1:F:76:GLU:HG2	3:F:240:FAD:N1A	2.27	0.49
1:D:49:GLU:HG2	5:D:274:HOH:O	2.12	0.49
1:D:209:LEU:HD22	1:D:213:LEU:HG	1.94	0.49
1:F:26:VAL:HG11	1:F:213:LEU:HD23	1.94	0.49
1:B:229:LEU:HD12	1:B:229:LEU:N	2.28	0.48
1:C:55:GLY:O	1:C:59:LYS:HD3	2.13	0.48
1:E:87:LEU:O	1:E:179:THR:HG22	2.13	0.48
1:F:141:ASN:ND2	1:F:143:LYS:H	2.11	0.48
1:B:103:LEU:N	1:B:103:LEU:HD23	2.28	0.48
1:B:154:ASN:HB3	1:D:147:MET:SD	2.54	0.48
1:C:161:GLU:HG3	1:C:162:HIS:CD2	2.48	0.48
1:A:234:HIS:HD2	5:A:536:HOH:O	1.96	0.48
1:A:99:ARG:NH2	1:B:76:GLU:OE1	2.47	0.47
1:A:120:SER:O	1:A:131:ARG:NH2	2.48	0.47
1:B:100:ILE:HG22	1:B:100:ILE:O	2.13	0.47
1:C:94:GLU:OE2	1:C:206:MET:HG3	2.15	0.47
1:C:103:LEU:HD23	1:C:103:LEU:N	2.28	0.47
1:D:50:TYR:O	1:D:57:LYS:HE3	2.13	0.47
1:E:36:ILE:HD11	1:F:30:HIS:HB3	1.94	0.47
1:F:97:ARG:HG3	2:F:241:UMP:C2'	2.37	0.47
1:F:113:ARG:HG3	5:F:243:HOH:O	2.12	0.47
1:D:35:ASP:HA	1:D:38:SER:OG	2.13	0.47
1:B:35:ASP:HA	1:B:38:SER:HB3	1.95	0.47
1:C:157:ILE:HG23	1:C:161:GLU:CD	2.39	0.47
1:D:229:LEU:O	1:D:230:LYS:CB	2.61	0.47
1:B:26:VAL:O	1:B:27:ILE:HD13	2.15	0.47
1:A:73:SER:HB3	3:B:240:FAD:O4'	2.14	0.47
1:F:211:LYS:HG2	1:F:215:ASP:OD2	2.14	0.47
1:C:112:LEU:O	1:C:159:LEU:HD13	2.14	0.47
1:E:123:PRO:HA	4:E:2:SO4:S	2.55	0.47
1:C:144:VAL:HG23	1:C:145:ASN:N	2.29	0.46
1:D:26:VAL:HG13	1:D:84:ILE:HD13	1.96	0.46
1:B:171:TYR:HB3	1:D:176:SER:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:LEU:HD11	1:B:152:LEU:HD22	1.97	0.46
1:B:158:LEU:HD21	1:D:147:MET:HG3	1.98	0.46
1:D:38:SER:CA	1:D:62:ILE:HD11	2.46	0.46
1:E:209:LEU:HD22	1:E:213:LEU:HG	1.97	0.46
1:F:52:ASP:OD2	1:F:55:GLY:HA3	2.14	0.46
1:F:122:LEU:CD1	1:F:153:GLU:HG2	2.45	0.46
1:A:117:GLU:CD	1:A:117:GLU:H	2.23	0.46
1:B:113:ARG:O	1:B:116:LYS:HB2	2.15	0.46
1:E:229:LEU:O	1:E:230:LYS:HB2	2.16	0.46
1:A:103:LEU:N	1:A:103:LEU:CD2	2.77	0.46
1:B:89:ARG:NH2	1:D:93:GLN:HG2	2.31	0.46
1:C:98:HIS:CE1	1:C:197:ARG:HH21	2.34	0.46
1:E:125:ASN:ND2	1:E:128:ASN:ND2	2.64	0.46
1:F:197:ARG:NH1	2:F:241:UMP:HN3	1.95	0.46
1:F:198:SER:HB2	1:F:230:LYS:H	1.81	0.46
1:A:122:LEU:HA	1:A:123:PRO:C	2.41	0.46
1:F:87:LEU:O	1:F:179:THR:HG22	2.16	0.46
1:C:87:LEU:O	1:C:179:THR:HG22	2.15	0.46
1:D:136:LEU:O	1:D:178:LYS:HE3	2.16	0.46
2:A:241:UMP:H2"	1:C:97:ARG:HH11	1.81	0.45
1:B:143:LYS:O	1:B:147:MET:HG2	2.16	0.45
1:F:218:PRO:HG2	1:F:221:HIS:HB2	1.97	0.45
1:E:85:LYS:HG2	1:E:180:HIS:CD2	2.52	0.45
1:D:125:ASN:O	1:D:126:GLU:C	2.59	0.45
1:C:115:LEU:O	1:C:118:VAL:HG12	2.16	0.45
1:A:100:ILE:O	1:A:100:ILE:HG22	2.17	0.45
1:B:79:TYR:CD1	1:C:79:TYR:HB2	2.52	0.45
1:C:199:SER:OG	1:C:201:LYS:HB2	2.16	0.45
1:E:25:GLU:CG	1:E:85:LYS:HB2	2.46	0.45
1:E:139:VAL:CG2	1:E:144:VAL:HG21	2.47	0.45
1:F:42:ARG:HG2	1:F:42:ARG:HH11	1.82	0.45
1:B:113:ARG:HD3	5:B:319:HOH:O	2.17	0.44
1:A:113:ARG:O	1:A:113:ARG:HG2	2.17	0.44
1:B:31:TYR:HA	1:B:79:TYR:O	2.18	0.44
5:C:288:HOH:O	1:D:71:HIS:HD2	2.00	0.44
1:E:223:TYR:CE1	1:E:224:LEU:HD13	2.52	0.44
1:D:112:LEU:C	1:D:114:GLU:H	2.24	0.44
1:C:37:ALA:O	1:C:41:ILE:HG13	2.18	0.44
1:D:141:ASN:OD1	1:D:144:VAL:HG13	2.17	0.44
1:A:94:GLU:OE2	1:A:206:MET:HG3	2.16	0.44
1:A:154:ASN:ND2	1:C:147:MET:HE2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:GLU:OE1	1:A:117:GLU:N	2.50	0.44
1:A:145:ASN:O	1:A:149:VAL:HG23	2.18	0.44
1:B:130:GLU:HG3	5:B:431:HOH:O	2.17	0.44
1:D:98:HIS:HE1	1:D:197:ARG:NE	2.08	0.44
1:B:97:ARG:HH11	2:D:241:UMP:H2'	1.82	0.44
1:E:44:CYS:HB2	1:E:45:TRP:CE3	2.52	0.44
1:F:165:LYS:HE2	5:F:504:HOH:O	2.17	0.44
5:B:379:HOH:O	1:C:106:LYS:HE3	2.17	0.43
3:C:240:FAD:H2'	3:C:240:FAD:H9	2.00	0.43
1:E:36:ILE:HA	1:E:39:GLN:HE21	1.83	0.43
1:F:143:LYS:O	1:F:147:MET:HG2	2.18	0.43
1:C:89:ARG:HD2	1:C:173:MET:CE	2.49	0.43
1:C:62:ILE:HG21	1:C:224:LEU:HD21	2.00	0.43
1:E:79:TYR:CD2	1:F:79:TYR:HB2	2.53	0.43
1:C:60:GLU:HG3	1:C:64:ARG:HH12	1.83	0.43
1:E:73:SER:HB3	3:E:240:FAD:O4'	2.17	0.43
1:B:98:HIS:HE1	1:B:197:ARG:NE	2.02	0.43
1:C:139:VAL:CG2	1:C:144:VAL:HG21	2.48	0.43
1:A:199:SER:OG	1:A:237:HIS:CD2	2.71	0.43
1:C:52:ASP:CG	1:C:57:LYS:HB2	2.42	0.43
1:E:141:ASN:O	1:E:144:VAL:HG22	2.19	0.43
1:A:43:THR:HG21	1:D:181:LEU:HA	2.01	0.43
1:A:64:ARG:O	1:A:68:ILE:HG13	2.19	0.43
1:D:68:ILE:HG22	1:D:70:ARG:HG2	2.00	0.43
1:E:36:ILE:HD11	1:F:30:HIS:CB	2.49	0.43
1:A:42:ARG:CZ	1:A:61:LEU:HD22	2.49	0.42
1:B:99:ARG:HB3	1:D:99:ARG:CB	2.45	0.42
1:D:120:SER:O	1:D:131:ARG:NH2	2.52	0.42
1:D:230:LYS:HD3	1:D:230:LYS:HA	1.88	0.42
1:E:200:ASN:ND2	5:E:493:HOH:O	2.52	0.42
1:F:141:ASN:HD22	1:F:143:LYS:N	2.17	0.42
1:A:34:LEU:HD22	1:A:62:ILE:HD12	2.01	0.42
1:A:223:TYR:CZ	1:A:224:LEU:HD13	2.54	0.42
1:E:89:ARG:NE	1:E:173:MET:HE3	2.34	0.42
1:F:103:LEU:N	1:F:103:LEU:HD23	2.34	0.42
1:C:99:ARG:HH12	3:C:240:FAD:C5'	2.32	0.42
1:C:141:ASN:HB3	1:C:144:VAL:CG2	2.49	0.42
1:B:49:GLU:OE1	1:B:50:TYR:HE1	2.02	0.42
1:B:112:LEU:HD23	1:B:112:LEU:HA	1.80	0.42
1:E:32:THR:O	1:E:221:HIS:HE1	2.02	0.42
1:E:39:GLN:OE1	1:F:29:LYS:HE3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ALA:HB3	1:A:62:ILE:HD13	2.02	0.42
1:A:99:ARG:HB2	1:C:99:ARG:HB3	2.02	0.42
1:C:34:LEU:CD1	1:C:59:LYS:HE3	2.47	0.42
1:B:59:LYS:HE3	1:B:59:LYS:HB2	1.81	0.42
1:D:37:ALA:CB	1:D:62:ILE:HD12	2.49	0.42
1:B:93:GLN:HG3	2:D:241:UMP:H3'	2.01	0.42
1:E:41:ILE:HD13	1:E:65:VAL:HB	2.02	0.42
1:A:70:ARG:HH11	1:A:71:HIS:CE1	2.37	0.42
1:A:100:ILE:HG13	3:A:240:FAD:H52A	2.02	0.41
1:C:194:LEU:CD1	1:C:225:PHE:HB3	2.50	0.41
1:D:26:VAL:HG11	1:D:213:LEU:HD23	2.02	0.41
1:E:75:LEU:O	1:E:187:ALA:HB3	2.20	0.41
1:B:94:GLU:OE2	1:B:206:MET:HG3	2.20	0.41
1:D:198:SER:HB3	1:D:229:LEU:HD12	2.01	0.41
1:A:52:ASP:O	1:A:53:ASP:HB2	2.21	0.41
1:B:218:PRO:HG2	1:B:221:HIS:CD2	2.54	0.41
1:C:112:LEU:C	1:C:114:GLU:H	2.29	0.41
1:C:170:LYS:HB2	1:C:170:LYS:NZ	2.35	0.41
1:D:118:VAL:HG13	1:D:131:ARG:HH12	1.86	0.41
1:F:111:THR:O	1:F:114:GLU:HG2	2.19	0.41
1:B:136:LEU:CD1	1:B:152:LEU:HD22	2.51	0.41
1:E:97:ARG:HH11	2:E:241:UMP:H2''	1.85	0.41
1:F:97:ARG:HH11	2:F:241:UMP:H2''	1.85	0.41
1:A:52:ASP:HB3	1:A:55:GLY:CA	2.51	0.41
1:A:144:VAL:CG2	1:A:145:ASN:N	2.83	0.41
1:A:193:PHE:CZ	1:A:209:LEU:HD13	2.56	0.41
1:C:73:SER:CB	3:D:240:FAD:O4'	2.67	0.41
1:E:34:LEU:HG	1:E:221:HIS:NE2	2.35	0.41
1:F:93:GLN:HE21	1:F:93:GLN:HB2	1.59	0.41
1:F:193:PHE:CZ	1:F:209:LEU:HD12	2.55	0.41
1:F:83:GLU:OE2	1:F:85:LYS:HE3	2.20	0.41
1:A:75:LEU:O	1:A:187:ALA:HB3	2.21	0.41
1:B:125:ASN:OD1	1:B:127:THR:HG23	2.21	0.41
1:B:141:ASN:HB3	1:B:144:VAL:HG13	2.03	0.41
1:C:197:ARG:O	1:C:207:GLN:HA	2.20	0.41
1:D:80:TYR:O	1:D:184:SER:HA	2.21	0.41
1:F:113:ARG:NH2	1:F:116:LYS:HZ3	2.18	0.41
1:A:234:HIS:CD2	5:A:536:HOH:O	2.73	0.41
1:D:81:ASN:ND2	5:D:271:HOH:O	2.54	0.41
1:E:113:ARG:HG2	1:E:166:ASN:HD21	1.86	0.41
1:E:158:LEU:HD22	1:E:164:ILE:CD1	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:118:VAL:HG23	1:F:156:ARG:CD	2.51	0.41
1:F:141:ASN:ND2	1:F:143:LYS:HB2	2.24	0.41
1:C:98:HIS:HA	3:C:240:FAD:O1A	2.21	0.40
1:D:63:HIS:HB2	1:D:223:TYR:CE1	2.56	0.40
1:E:36:ILE:HD13	1:F:81:ASN:CG	2.46	0.40
1:E:37:ALA:HB3	1:E:62:ILE:HD13	2.02	0.40
1:A:29:LYS:N	1:A:81:ASN:O	2.51	0.40
1:A:97:ARG:HD2	1:A:97:ARG:HA	1.89	0.40
1:A:108:SER:N	2:A:241:UMP:OP2	2.53	0.40
1:A:171:TYR:O	1:C:176:SER:HB3	2.21	0.40
1:B:52:ASP:OD1	1:B:55:GLY:N	2.54	0.40
1:F:120:SER:O	1:F:131:ARG:NH2	2.55	0.40
1:F:141:ASN:C	1:F:141:ASN:ND2	2.74	0.40
1:F:141:ASN:ND2	1:F:143:LYS:N	2.68	0.40
1:F:162:HIS:O	1:F:163:ASN:C	2.64	0.40
1:A:193:PHE:HZ	1:A:209:LEU:HD13	1.86	0.40
1:C:196:LEU:HD21	1:D:69:PHE:CD2	2.56	0.40
1:A:98:HIS:HA	3:A:240:FAD:O1A	2.21	0.40
1:A:103:LEU:HD23	1:A:103:LEU:H	1.87	0.40
1:A:223:TYR:CE1	1:A:224:LEU:HD13	2.57	0.40
1:E:162:HIS:O	1:E:163:ASN:C	2.65	0.40
1:F:136:LEU:O	1:F:178:LYS:HE3	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:257:HOH:O	5:F:503:HOH:O[2_555]	2.15	0.05

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	213/216 (99%)	203 (95%)	10 (5%)	0	100	100
1	B	205/216 (95%)	191 (93%)	13 (6%)	1 (0%)	24	43
1	C	205/216 (95%)	197 (96%)	8 (4%)	0	100	100
1	D	205/216 (95%)	196 (96%)	9 (4%)	0	100	100
1	E	205/216 (95%)	196 (96%)	9 (4%)	0	100	100
1	F	205/216 (95%)	193 (94%)	10 (5%)	2 (1%)	12	24
All	All	1238/1296 (96%)	1176 (95%)	59 (5%)	3 (0%)	43	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	120	SER
1	F	53	ASP
1	F	141	ASN

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	194/195 (100%)	183 (94%)	11 (6%)	18	39
1	B	186/195 (95%)	177 (95%)	9 (5%)	23	46
1	C	186/195 (95%)	177 (95%)	9 (5%)	23	46
1	D	186/195 (95%)	178 (96%)	8 (4%)	26	51
1	E	186/195 (95%)	180 (97%)	6 (3%)	34	62
1	F	186/195 (95%)	176 (95%)	10 (5%)	20	41
All	All	1124/1170 (96%)	1071 (95%)	53 (5%)	23	47

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	ASP
1	A	99	ARG

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Mol	Chain	Res	Type
1	A	103	LEU
1	A	117	GLU
1	A	124	LEU
1	A	150	LEU
1	A	170	LYS
1	A	209	LEU
1	A	217	LEU
1	A	224	LEU
1	A	229	LEU
1	B	99	ARG
1	B	103	LEU
1	B	119	GLU
1	B	124	LEU
1	B	144	VAL
1	B	170	LYS
1	B	209	LEU
1	B	217	LEU
1	B	223	TYR
1	C	78	LEU
1	C	89	ARG
1	C	103	LEU
1	C	147	MET
1	C	150	LEU
1	C	170	LYS
1	C	204	LYS
1	C	213	LEU
1	C	224	LEU
1	D	60	GLU
1	D	99	ARG
1	D	103	LEU
1	D	125	ASN
1	D	163	ASN
1	D	200	ASN
1	D	209	LEU
1	D	224	LEU
1	E	99	ARG
1	E	103	LEU
1	E	150	LEU
1	E	209	LEU
1	E	224	LEU
1	E	230	LYS
1	F	27	ILE

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Mol	Chain	Res	Type
1	F	93	GLN
1	F	141	ASN
1	F	142	GLU
1	F	144	VAL
1	F	150	LEU
1	F	170	LYS
1	F	217	LEU
1	F	224	LEU
1	F	228	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	71	HIS
1	A	81	ASN
1	A	98	HIS
1	A	162	HIS
1	A	163	ASN
1	A	166	ASN
1	A	180	HIS
1	A	192	ASN
1	A	231	HIS
1	A	237	HIS
1	A	238	HIS
1	B	67	ASN
1	B	81	ASN
1	B	93	GLN
1	B	98	HIS
1	B	128	ASN
1	B	162	HIS
1	B	163	ASN
1	B	180	HIS
1	B	192	ASN
1	C	39	GLN
1	C	46	GLN
1	C	71	HIS
1	C	81	ASN
1	C	98	HIS
1	C	162	HIS
1	C	192	ASN
1	C	207	GLN

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Mol	Chain	Res	Type
1	D	46	GLN
1	D	71	HIS
1	D	81	ASN
1	D	98	HIS
1	D	125	ASN
1	D	162	HIS
1	D	180	HIS
1	D	192	ASN
1	D	200	ASN
1	E	46	GLN
1	E	81	ASN
1	E	98	HIS
1	E	125	ASN
1	E	180	HIS
1	E	200	ASN
1	E	222	GLN
1	F	71	HIS
1	F	81	ASN
1	F	93	GLN
1	F	98	HIS
1	F	141	ASN
1	F	200	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](#)

23 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	$\# Z > 2$
4	SO4	C	8	-	4,4,4	0.37	0	6,6,6	0.06	0
3	FAD	E	240	-	58,58,58	1.58	10 (17%)	85,89,89	1.20	5 (5%)
2	UMP	D	241	-	21,21,21	1.83	1 (4%)	30,31,31	1.85	8 (26%)
4	SO4	A	10	-	4,4,4	0.42	0	6,6,6	0.12	0
4	SO4	E	2	-	4,4,4	0.32	0	6,6,6	0.25	0
4	SO4	A	11	-	4,4,4	0.43	0	6,6,6	0.06	0
4	SO4	C	4	-	4,4,4	0.37	0	6,6,6	0.09	0
3	FAD	D	240	-	58,58,58	1.62	10 (17%)	85,89,89	1.23	7 (8%)
2	UMP	A	241	-	21,21,21	1.80	1 (4%)	30,31,31	1.88	9 (30%)
2	UMP	B	241	-	21,21,21	1.86	1 (4%)	30,31,31	1.89	9 (30%)
4	SO4	A	5	-	4,4,4	0.38	0	6,6,6	0.07	0
4	SO4	E	1	-	4,4,4	0.35	0	6,6,6	0.06	0
4	SO4	F	7	-	4,4,4	0.39	0	6,6,6	0.09	0
3	FAD	C	240	-	58,58,58	1.58	9 (15%)	85,89,89	1.21	7 (8%)
3	FAD	F	240	-	58,58,58	1.56	8 (13%)	85,89,89	1.16	6 (7%)
2	UMP	E	241	-	21,21,21	1.85	1 (4%)	30,31,31	1.86	9 (30%)
3	FAD	B	240	-	58,58,58	1.61	9 (15%)	85,89,89	1.16	5 (5%)
2	UMP	C	241	-	21,21,21	1.84	1 (4%)	30,31,31	1.88	8 (26%)
4	SO4	C	9	-	4,4,4	0.40	0	6,6,6	0.09	0
4	SO4	D	3	-	4,4,4	0.36	0	6,6,6	0.19	0
2	UMP	F	241	-	21,21,21	1.79	1 (4%)	30,31,31	1.84	7 (23%)
4	SO4	D	6	-	4,4,4	0.36	0	6,6,6	0.07	0
3	FAD	A	240	-	58,58,58	1.58	9 (15%)	85,89,89	1.13	7 (8%)

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	UMP	A	241	-	-	2/10/22/22	0/2/2/2
2	UMP	B	241	-	-	2/10/22/22	0/2/2/2
2	UMP	F	241	-	-	2/10/22/22	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	UMP	E	241	-	-	2/10/22/22	0/2/2/2
3	FAD	B	240	-	-	4/34/50/50	0/6/6/6
2	UMP	D	241	-	-	2/10/22/22	0/2/2/2
3	FAD	C	240	-	-	6/34/50/50	0/6/6/6
3	FAD	E	240	-	-	4/34/50/50	0/6/6/6
3	FAD	A	240	-	-	6/34/50/50	0/6/6/6
2	UMP	C	241	-	-	2/10/22/22	0/2/2/2
3	FAD	F	240	-	-	5/34/50/50	0/6/6/6
3	FAD	D	240	-	-	4/34/50/50	0/6/6/6

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	241	UMP	C6-C5	7.26	1.51	1.35
2	C	241	UMP	C6-C5	7.13	1.51	1.35
2	D	241	UMP	C6-C5	7.10	1.51	1.35
2	B	241	UMP	C6-C5	7.03	1.51	1.35
2	F	241	UMP	C6-C5	6.96	1.51	1.35
2	A	241	UMP	C6-C5	6.87	1.51	1.35
3	B	240	FAD	C4X-N5	5.92	1.43	1.30
3	C	240	FAD	C4X-N5	5.55	1.42	1.30
3	F	240	FAD	C4X-N5	5.48	1.42	1.30
3	E	240	FAD	C4X-N5	5.28	1.42	1.30
3	D	240	FAD	C4X-N5	5.24	1.42	1.30
3	A	240	FAD	C4X-N5	5.18	1.41	1.30
3	D	240	FAD	C9-C8	4.65	1.45	1.39
3	A	240	FAD	C9-C8	4.44	1.45	1.39
3	E	240	FAD	C9-C8	4.28	1.45	1.39
3	F	240	FAD	C9-C8	4.11	1.45	1.39
3	C	240	FAD	C9A-N10	4.00	1.48	1.41
3	D	240	FAD	C9A-N10	3.97	1.48	1.41
3	C	240	FAD	C9-C8	3.90	1.44	1.39
3	F	240	FAD	C9A-N10	3.80	1.47	1.41
3	B	240	FAD	C9-C8	3.74	1.44	1.39
3	A	240	FAD	C9A-N10	3.71	1.47	1.41
3	A	240	FAD	C4A-N3A	3.62	1.41	1.34
3	B	240	FAD	C9A-N10	3.58	1.47	1.41
3	A	240	FAD	C9A-C5X	3.42	1.46	1.41
3	C	240	FAD	C9A-C5X	3.36	1.46	1.41
3	E	240	FAD	C6-C5X	3.35	1.45	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	240	FAD	C9A-C5X	3.34	1.46	1.41
3	E	240	FAD	C9A-N10	3.34	1.46	1.41
3	F	240	FAD	C6-C5X	3.32	1.45	1.40
3	B	240	FAD	C6-C5X	3.30	1.45	1.40
3	C	240	FAD	C4A-N3A	3.27	1.40	1.34
3	B	240	FAD	C9A-C5X	3.25	1.46	1.41
3	F	240	FAD	C4A-N3A	3.24	1.40	1.34
3	B	240	FAD	C4A-N3A	3.23	1.40	1.34
3	F	240	FAD	C9A-C5X	3.16	1.46	1.41
3	D	240	FAD	C4A-N3A	3.05	1.40	1.34
3	E	240	FAD	C4A-N3A	2.98	1.40	1.34
3	C	240	FAD	C6-C5X	2.94	1.44	1.40
3	E	240	FAD	C9A-C5X	2.93	1.45	1.41
3	D	240	FAD	PA-O5B	-2.88	1.48	1.59
3	D	240	FAD	C6-C5X	2.84	1.44	1.40
3	E	240	FAD	PA-O5B	-2.74	1.48	1.59
3	C	240	FAD	PA-O5B	-2.71	1.48	1.59
3	A	240	FAD	C10-N1	2.65	1.38	1.33
3	F	240	FAD	PA-O5B	-2.63	1.49	1.59
3	B	240	FAD	C10-N1	2.63	1.38	1.33
3	E	240	FAD	PA-O3P	-2.62	1.56	1.59
3	A	240	FAD	C6-C5X	2.61	1.43	1.40
3	D	240	FAD	C1'-C2'	-2.59	1.49	1.52
3	B	240	FAD	PA-O5B	-2.58	1.49	1.59
3	A	240	FAD	PA-O5B	-2.54	1.49	1.59
3	C	240	FAD	C1'-C2'	-2.49	1.49	1.52
3	D	240	FAD	PA-O3P	-2.42	1.56	1.59
3	C	240	FAD	C10-N1	2.35	1.38	1.33
3	B	240	FAD	C4'-C3'	-2.34	1.49	1.53
3	F	240	FAD	C10-N1	2.34	1.37	1.33
3	E	240	FAD	C10-N1	2.32	1.37	1.33
3	E	240	FAD	C4'-C3'	-2.14	1.49	1.53
3	A	240	FAD	C2A-N1A	2.13	1.37	1.33
3	D	240	FAD	C10-N1	2.03	1.37	1.33

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	241	UMP	N3-C2-N1	5.03	121.43	114.89
2	E	241	UMP	N3-C2-N1	4.83	121.17	114.89
2	D	241	UMP	N3-C2-N1	4.81	121.16	114.89
2	C	241	UMP	N3-C2-N1	4.81	121.15	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	241	UMP	N3-C2-N1	4.60	120.88	114.89
3	E	240	FAD	O5'-C5'-C4'	-4.59	97.12	109.36
2	F	241	UMP	N3-C2-N1	4.58	120.85	114.89
3	B	240	FAD	O5'-C5'-C4'	-4.43	97.53	109.36
3	D	240	FAD	O5'-C5'-C4'	-4.33	97.79	109.36
3	F	240	FAD	O5'-C5'-C4'	-4.09	98.44	109.36
3	A	240	FAD	O5'-C5'-C4'	-4.08	98.48	109.36
3	C	240	FAD	O5'-C5'-C4'	-3.91	98.92	109.36
2	A	241	UMP	C5-C4-N3	3.84	120.18	114.80
2	F	241	UMP	C5-C4-N3	3.70	119.99	114.80
3	E	240	FAD	C4'-C3'-C2'	-3.68	107.45	113.57
2	B	241	UMP	C5-C4-N3	3.64	119.90	114.80
2	D	241	UMP	C5-C4-N3	3.64	119.90	114.80
2	C	241	UMP	C5-C4-N3	3.59	119.82	114.80
2	E	241	UMP	C5-C4-N3	3.58	119.82	114.80
3	C	240	FAD	O4B-C4B-C5B	-3.42	98.38	109.33
3	B	240	FAD	C4'-C3'-C2'	-3.41	107.89	113.57
2	F	241	UMP	O4-C4-C5	-3.40	119.30	125.16
3	D	240	FAD	C4'-C3'-C2'	-3.36	107.98	113.57
3	C	240	FAD	C4'-C3'-C2'	-3.33	108.03	113.57
2	A	241	UMP	O4-C4-C5	-3.32	119.43	125.16
3	F	240	FAD	C4'-C3'-C2'	-3.27	108.12	113.57
2	C	241	UMP	O4-C4-C5	-3.23	119.58	125.16
3	C	240	FAD	C4X-C10-N10	3.20	121.06	116.48
3	D	240	FAD	O4B-C4B-C5B	-3.19	99.13	109.33
3	E	240	FAD	O4B-C4B-C5B	-3.11	99.36	109.33
2	E	241	UMP	O4-C4-C5	-3.10	119.81	125.16
2	D	241	UMP	O4-C4-C5	-3.07	119.86	125.16
3	D	240	FAD	C4X-C10-N10	3.06	120.87	116.48
3	F	240	FAD	O4B-C4B-C5B	-3.04	99.60	109.33
3	A	240	FAD	O4B-C4B-C5B	-3.03	99.64	109.33
3	E	240	FAD	C4X-C10-N10	2.97	120.73	116.48
3	F	240	FAD	C4X-C10-N10	2.97	120.73	116.48
3	B	240	FAD	O4B-C4B-C5B	-2.97	99.83	109.33
2	D	241	UMP	O4'-C1'-N1	2.92	113.04	107.86
2	B	241	UMP	O4'-C1'-N1	2.92	113.04	107.86
2	B	241	UMP	O4-C4-C5	-2.91	120.14	125.16
3	A	240	FAD	C4X-C10-N10	2.83	120.53	116.48
3	C	240	FAD	C1'-N10-C9A	2.82	126.11	120.63
2	A	241	UMP	O4'-C1'-N1	2.82	112.87	107.86
2	E	241	UMP	C4-N3-C2	-2.82	123.11	126.61
2	C	241	UMP	C4-N3-C2	-2.81	123.12	126.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	240	FAD	C4'-C3'-C2'	-2.81	108.90	113.57
2	F	241	UMP	C4-N3-C2	-2.80	123.13	126.61
2	C	241	UMP	O4'-C1'-N1	2.73	112.71	107.86
2	A	241	UMP	C6-C5-C4	-2.71	116.07	119.53
2	A	241	UMP	C4-N3-C2	-2.71	123.25	126.61
2	E	241	UMP	O4'-C1'-N1	2.70	112.65	107.86
3	B	240	FAD	C4X-C10-N10	2.65	120.28	116.48
2	F	241	UMP	O4'-C1'-N1	2.61	112.49	107.86
2	D	241	UMP	C6-C5-C4	-2.61	116.20	119.53
3	D	240	FAD	C4A-N9A-C8A	-2.59	103.02	105.74
2	B	241	UMP	C4-N3-C2	-2.57	123.42	126.61
2	B	241	UMP	C6-C5-C4	-2.56	116.26	119.53
2	D	241	UMP	C4-N3-C2	-2.56	123.44	126.61
2	E	241	UMP	C2'-C1'-N1	2.55	120.19	113.81
2	C	241	UMP	C6-C5-C4	-2.55	116.28	119.53
2	E	241	UMP	C6-C5-C4	-2.47	116.38	119.53
2	B	241	UMP	C2'-C1'-N1	2.44	119.91	113.81
3	E	240	FAD	C4A-N9A-C8A	-2.42	103.20	105.74
2	C	241	UMP	C2'-C1'-N1	2.37	119.73	113.81
2	F	241	UMP	C1'-N1-C2	2.36	122.28	117.66
3	D	240	FAD	C1'-N10-C9A	2.36	125.23	120.63
2	B	241	UMP	C1'-N1-C2	2.31	122.17	117.66
2	F	241	UMP	C6-C5-C4	-2.30	116.60	119.53
3	D	240	FAD	C9-C9A-N10	2.28	124.92	121.85
2	A	241	UMP	C1'-N1-C2	2.27	122.11	117.66
2	B	241	UMP	O2-C2-N3	-2.27	117.31	121.49
2	A	241	UMP	C2'-C1'-N1	2.26	119.47	113.81
3	C	240	FAD	C9-C9A-N10	2.22	124.83	121.85
2	D	241	UMP	C1'-N1-C2	2.21	121.99	117.66
2	C	241	UMP	C1'-N1-C2	2.20	121.97	117.66
2	E	241	UMP	C1'-N1-C2	2.19	121.94	117.66
2	D	241	UMP	C2'-C1'-N1	2.15	119.19	113.81
3	C	240	FAD	C4A-N9A-C8A	-2.15	103.48	105.74
3	B	240	FAD	C4A-N9A-C8A	-2.13	103.50	105.74
3	F	240	FAD	C4A-N9A-C8A	-2.10	103.53	105.74
3	A	240	FAD	C1'-N10-C9A	2.07	124.65	120.63
3	A	240	FAD	C9-C9A-N10	2.05	124.61	121.85
3	A	240	FAD	C4A-N9A-C8A	-2.04	103.60	105.74
3	F	240	FAD	C1'-N10-C9A	2.03	124.58	120.63
2	A	241	UMP	O5'-P-OP1	2.02	111.90	106.44
2	E	241	UMP	O5'-P-OP1	2.01	111.87	106.44

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	240	FAD	C2'-C1'-N10-C10
3	D	240	FAD	C2'-C1'-N10-C10
2	A	241	UMP	C2'-C1'-N1-C2
2	B	241	UMP	C2'-C1'-N1-C2
2	D	241	UMP	C2'-C1'-N1-C2
2	E	241	UMP	C2'-C1'-N1-C2
2	F	241	UMP	C2'-C1'-N1-C2
2	C	241	UMP	C2'-C1'-N1-C2
2	A	241	UMP	C2'-C1'-N1-C6
2	B	241	UMP	C2'-C1'-N1-C6
2	D	241	UMP	C2'-C1'-N1-C6
2	E	241	UMP	C2'-C1'-N1-C6
2	F	241	UMP	C2'-C1'-N1-C6
3	A	240	FAD	P-O3P-PA-O1A
3	B	240	FAD	P-O3P-PA-O1A
3	C	240	FAD	P-O3P-PA-O1A
3	D	240	FAD	P-O3P-PA-O1A
3	E	240	FAD	P-O3P-PA-O1A
3	F	240	FAD	P-O3P-PA-O1A
2	C	241	UMP	C2'-C1'-N1-C6
3	C	240	FAD	C4'-C5'-O5'-P
3	B	240	FAD	PA-O3P-P-O2P
3	F	240	FAD	PA-O3P-P-O2P
3	D	240	FAD	C4'-C5'-O5'-P
3	F	240	FAD	C4'-C5'-O5'-P
3	A	240	FAD	PA-O3P-P-O2P
3	A	240	FAD	C4'-C5'-O5'-P
3	B	240	FAD	C4B-C5B-O5B-PA
3	E	240	FAD	C4'-C5'-O5'-P
3	A	240	FAD	PA-O3P-P-O1P
3	B	240	FAD	PA-O3P-P-O1P
3	C	240	FAD	PA-O3P-P-O1P
3	D	240	FAD	PA-O3P-P-O1P
3	E	240	FAD	PA-O3P-P-O1P
3	F	240	FAD	O4B-C4B-C5B-O5B
3	A	240	FAD	C4B-C5B-O5B-PA
3	C	240	FAD	C4B-C5B-O5B-PA
3	F	240	FAD	C4B-C5B-O5B-PA
3	E	240	FAD	PA-O3P-P-O2P
3	A	240	FAD	O4B-C4B-C5B-O5B
3	C	240	FAD	O4B-C4B-C5B-O5B

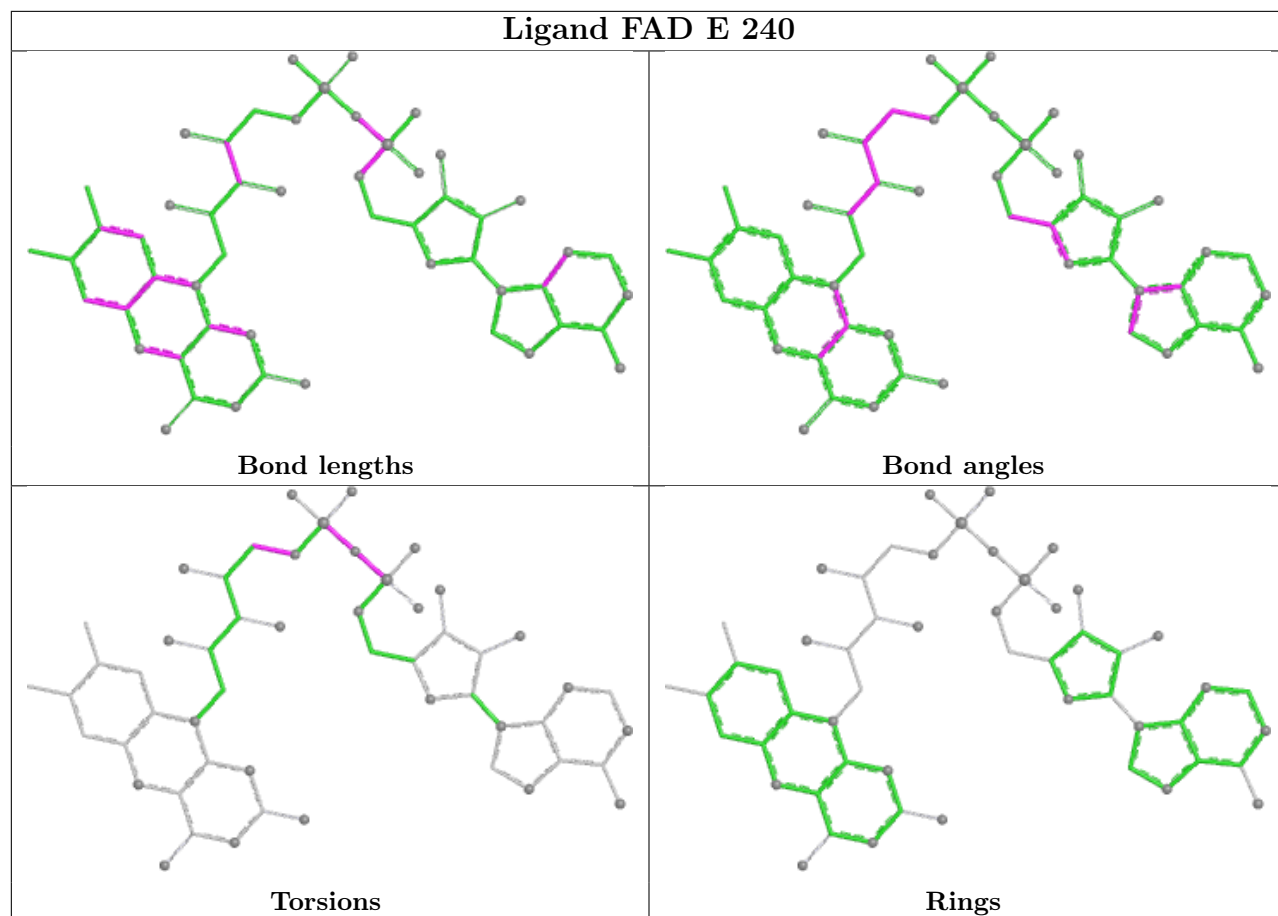
There are no ring outliers.

14 monomers are involved in 74 short contacts:

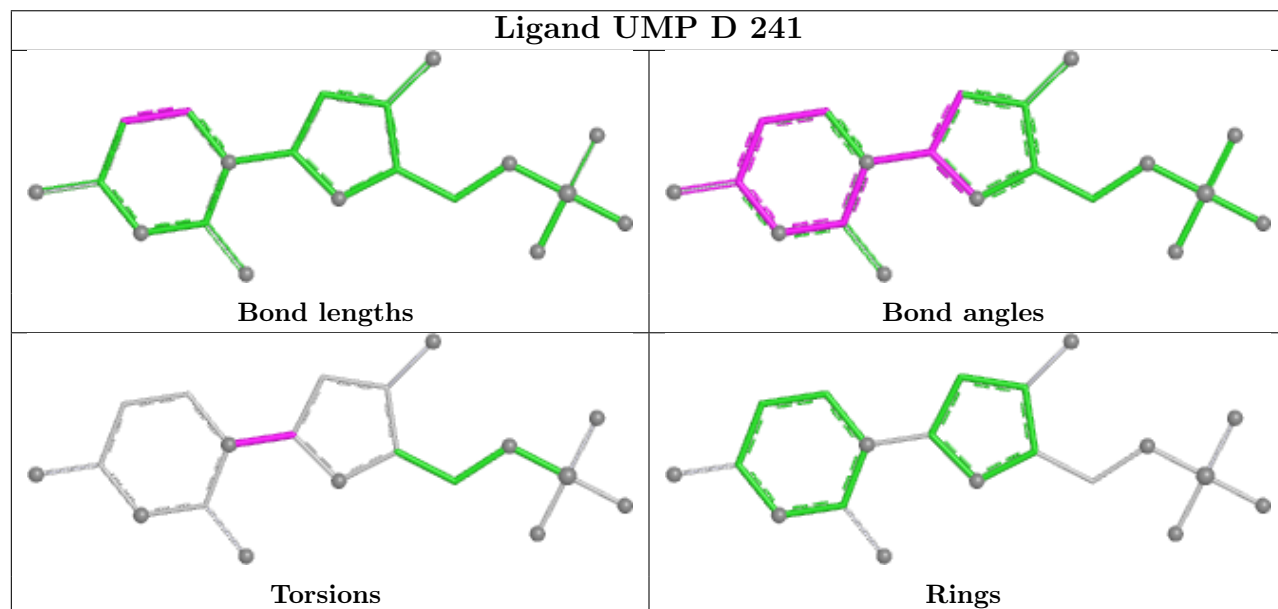
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	240	FAD	2	0
2	D	241	UMP	6	0
4	E	2	SO4	3	0
3	D	240	FAD	8	0
2	A	241	UMP	6	0
2	B	241	UMP	6	0
4	E	1	SO4	1	0
3	C	240	FAD	7	0
3	F	240	FAD	3	0
2	E	241	UMP	4	0
3	B	240	FAD	7	0
2	C	241	UMP	7	0
2	F	241	UMP	6	0
3	A	240	FAD	8	0

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.

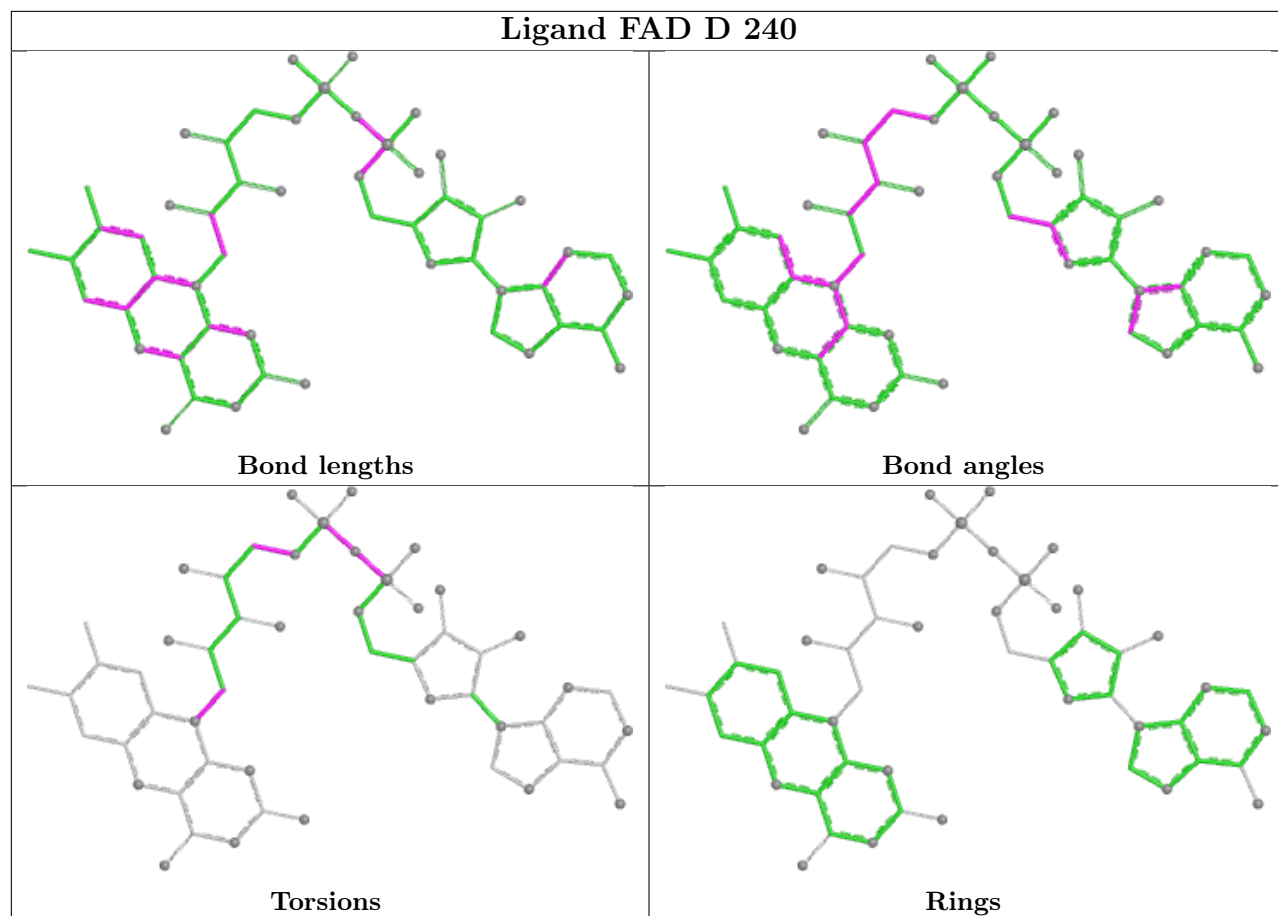
Ligand FAD E 240



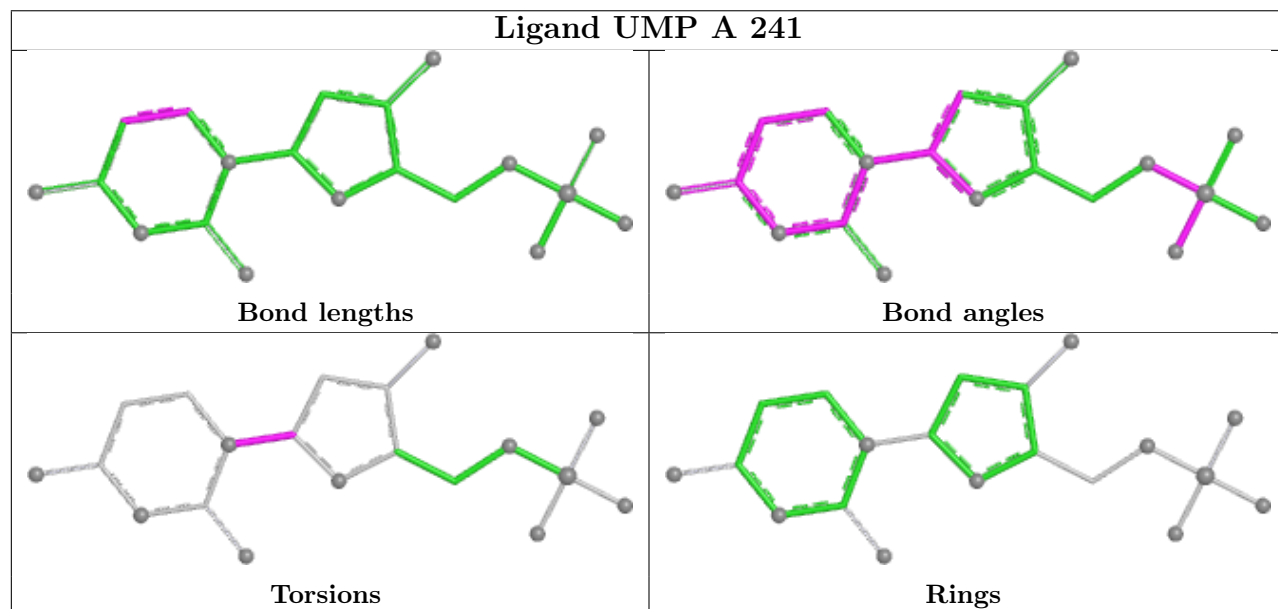
Ligand UMP D 241

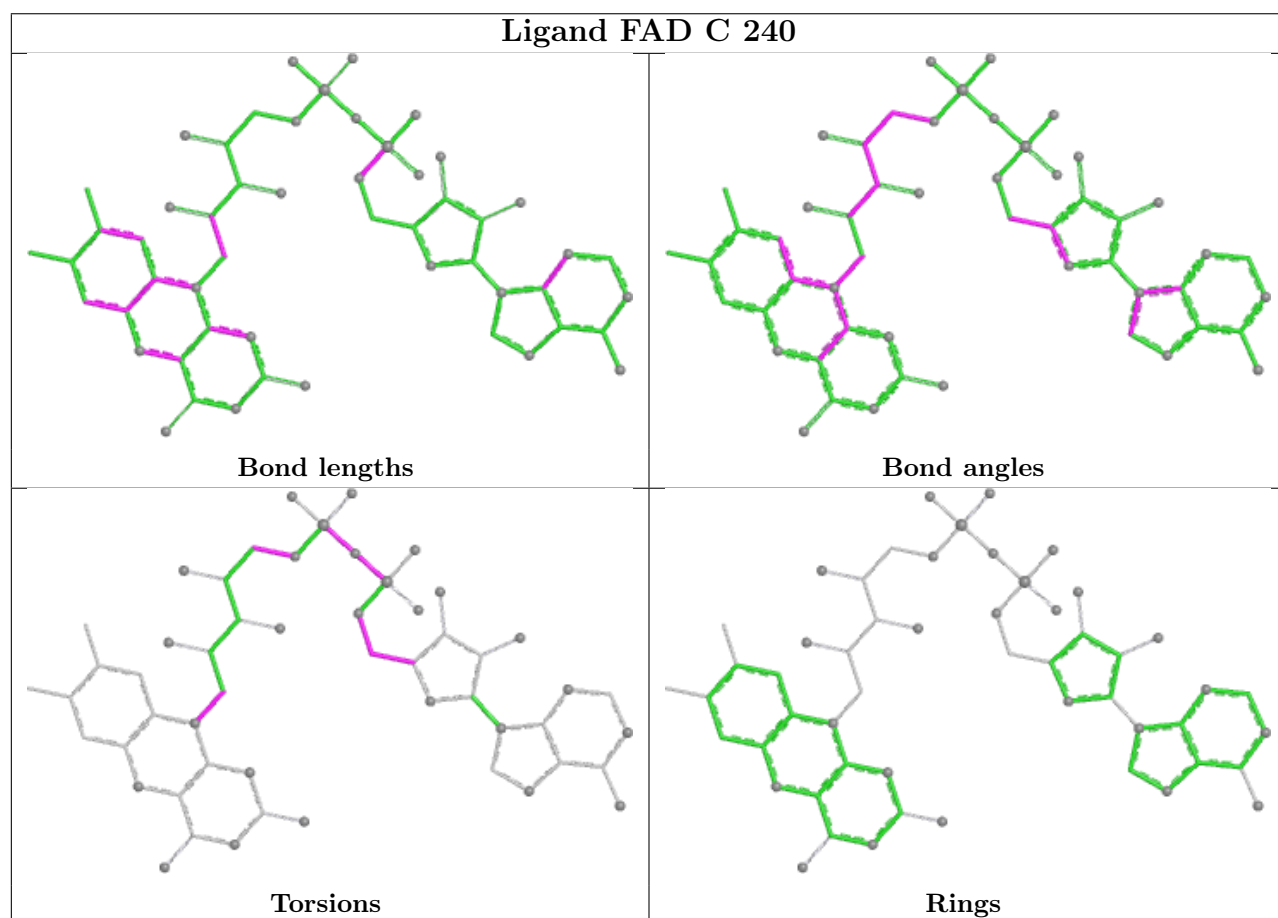
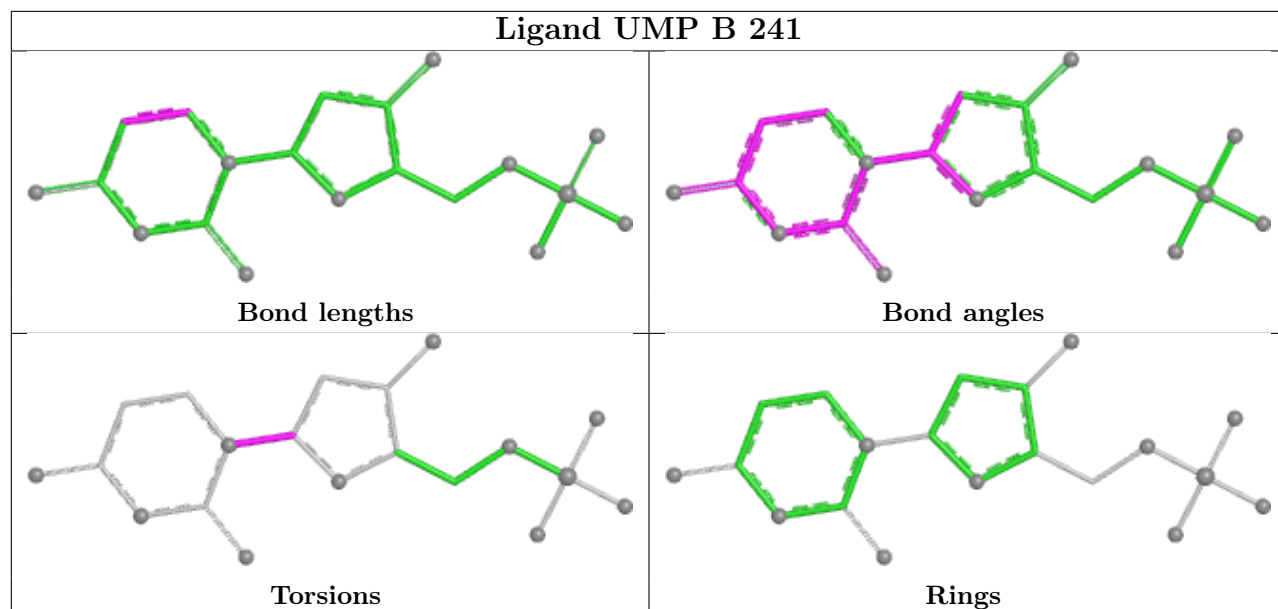


Ligand FAD D 240

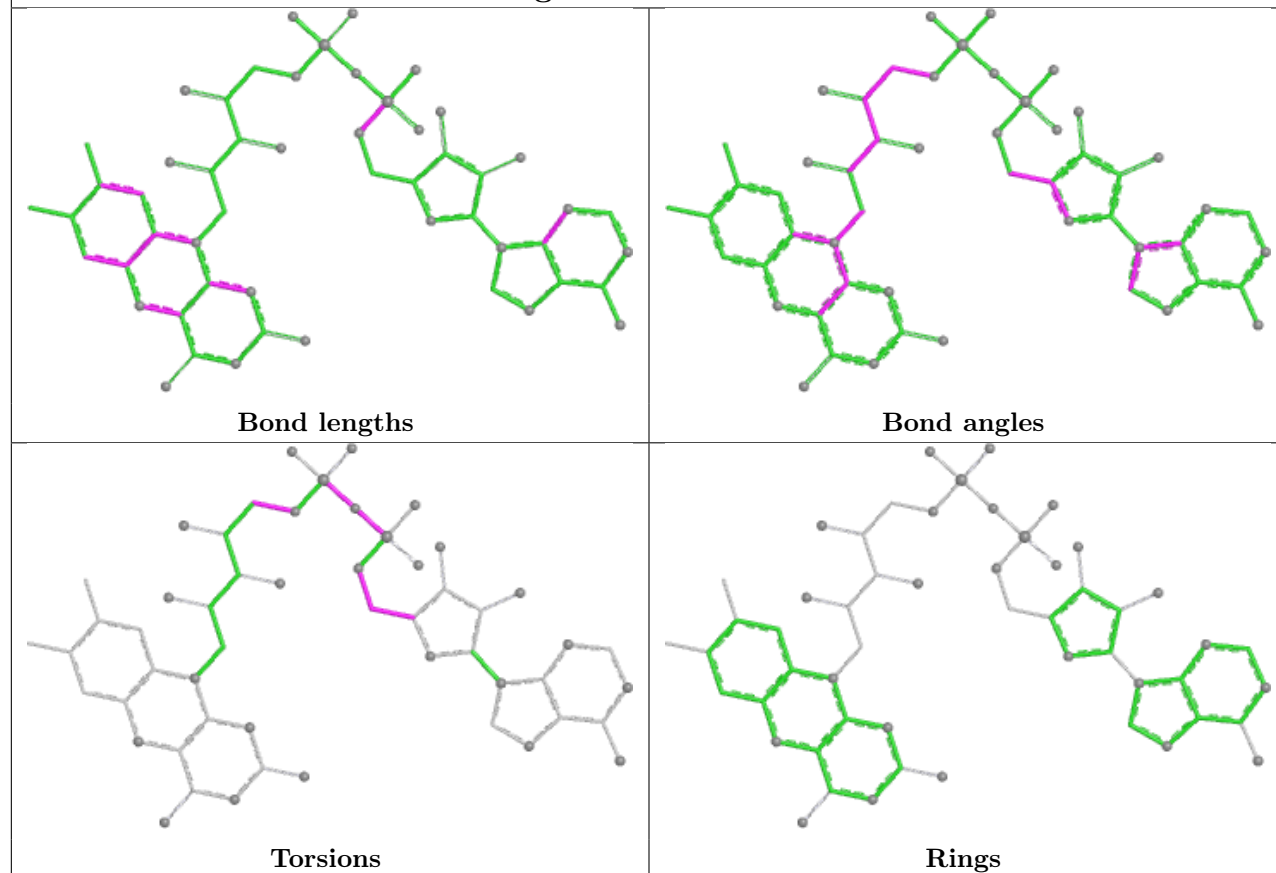


Ligand UMP A 241

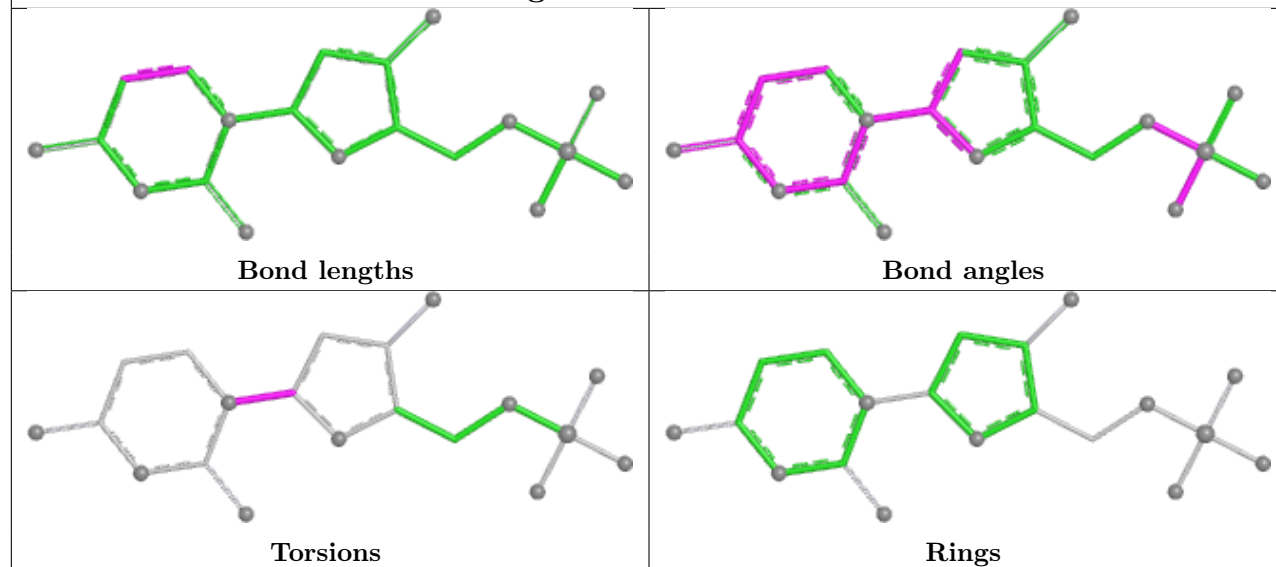




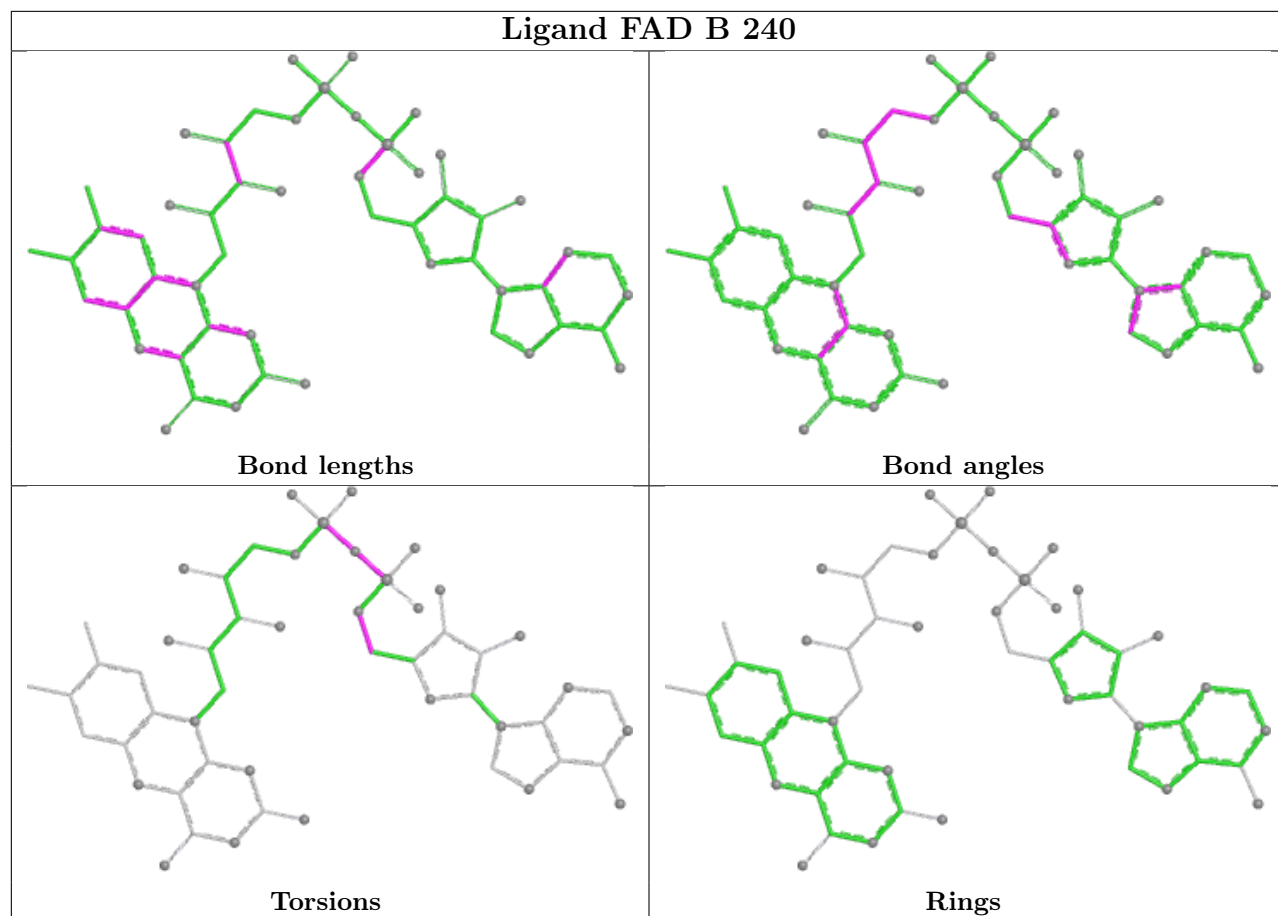
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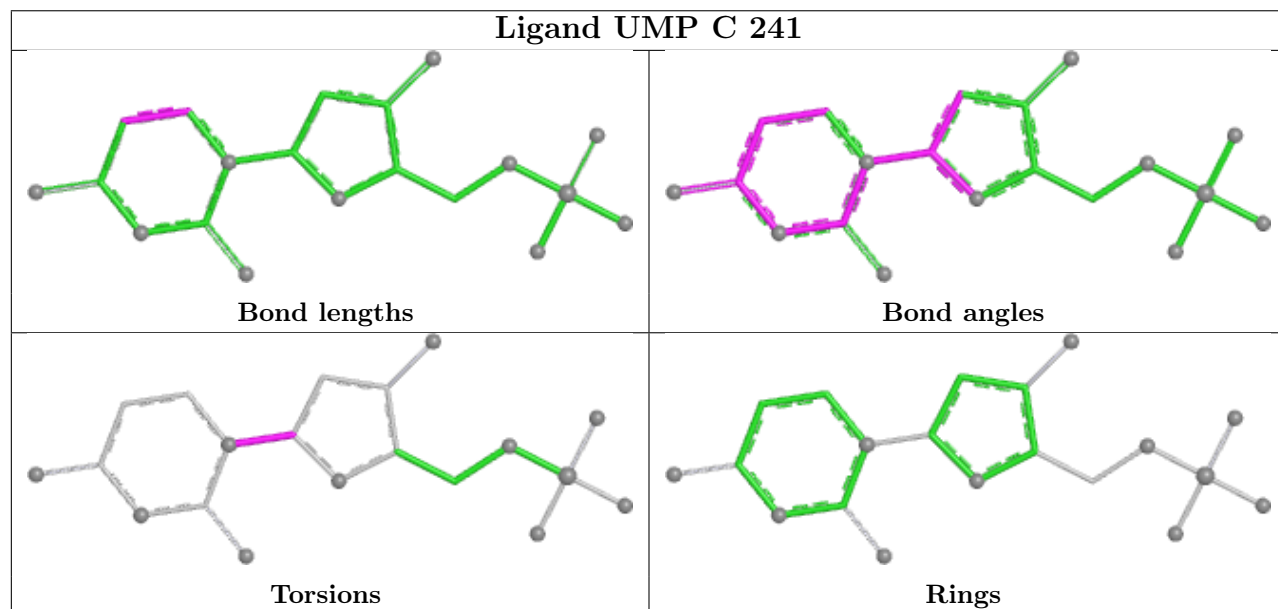
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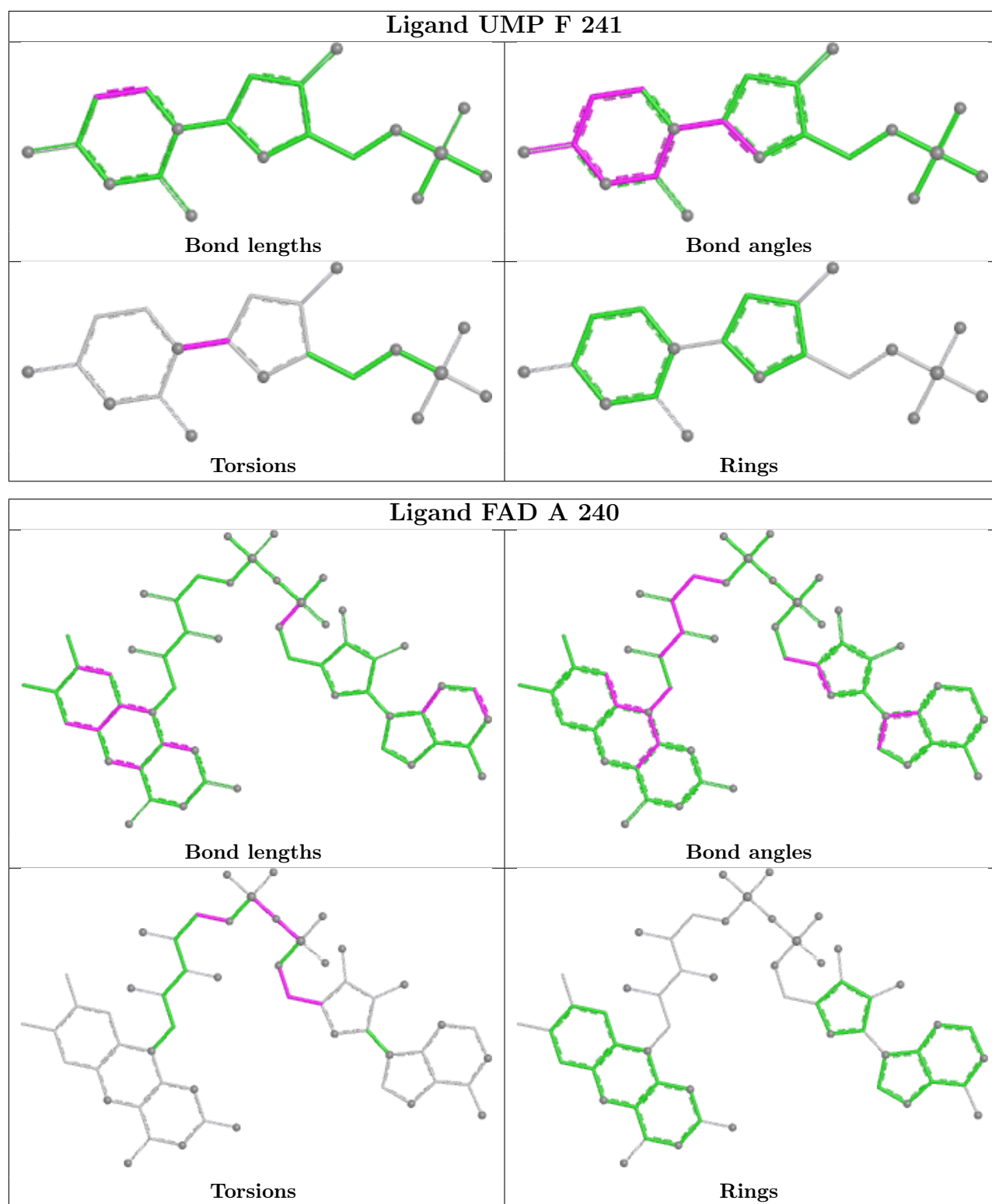


Ligand FAD B 240



Ligand UMP C 241





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	215/216 (99%)	-0.34	1 (0%) 87 85	19, 30, 53, 79	0
1	B	207/216 (95%)	-0.40	0 100 100	15, 30, 57, 72	0
1	C	207/216 (95%)	-0.52	0 100 100	14, 25, 43, 54	0
1	D	207/216 (95%)	-0.36	2 (0%) 79 76	18, 32, 51, 67	0
1	E	207/216 (95%)	-0.23	0 100 100	20, 34, 53, 65	0
1	F	207/216 (95%)	-0.19	0 100 100	20, 37, 57, 71	0
All	All	1250/1296 (96%)	-0.34	3 (0%) 91 89	14, 31, 53, 79	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	230	LYS	2.4
1	A	52	ASP	2.2
1	D	120	SER	2.1

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

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

6.3 Carbohydrates [i](#)

There are no 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,

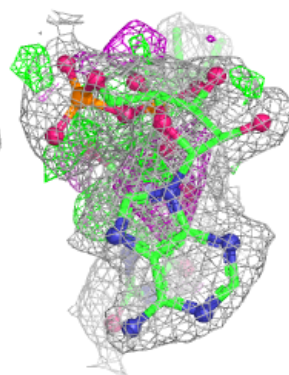
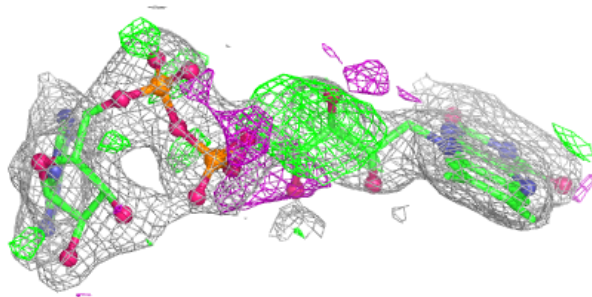
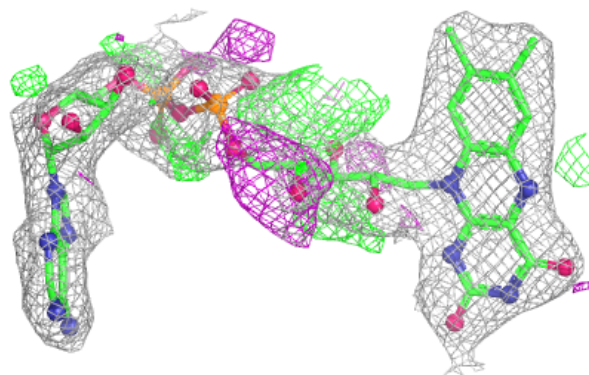
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	D	3	5/5	0.73	0.28	117,117,118,118	0
4	SO4	C	9	5/5	0.78	0.15	83,84,84,84	0
4	SO4	E	1	5/5	0.79	0.17	89,89,90,90	0
4	SO4	D	6	5/5	0.82	0.14	90,90,90,91	0
4	SO4	A	11	5/5	0.82	0.18	82,82,82,83	0
4	SO4	F	7	5/5	0.83	0.14	96,96,96,96	0
4	SO4	A	5	5/5	0.85	0.15	79,80,81,81	0
4	SO4	C	8	5/5	0.86	0.14	74,75,76,76	0
4	SO4	A	10	5/5	0.86	0.13	69,70,70,71	0
3	FAD	A	240	53/53	0.88	0.12	18,27,47,50	0
3	FAD	B	240	53/53	0.88	0.12	27,34,51,57	0
4	SO4	E	2	5/5	0.88	0.28	119,119,120,120	0
3	FAD	F	240	53/53	0.88	0.13	29,36,50,54	0
3	FAD	C	240	53/53	0.89	0.12	17,25,40,44	0
3	FAD	E	240	53/53	0.90	0.12	22,27,45,50	0
3	FAD	D	240	53/53	0.91	0.11	18,26,41,48	0
2	UMP	D	241	20/20	0.92	0.11	39,40,44,44	0
2	UMP	E	241	20/20	0.92	0.09	28,32,38,41	0
4	SO4	C	4	5/5	0.93	0.08	71,71,72,73	0
2	UMP	F	241	20/20	0.93	0.10	34,38,42,43	0
2	UMP	C	241	20/20	0.94	0.08	26,28,32,33	0
2	UMP	A	241	20/20	0.94	0.09	24,28,32,35	0
2	UMP	B	241	20/20	0.94	0.09	22,26,32,33	0

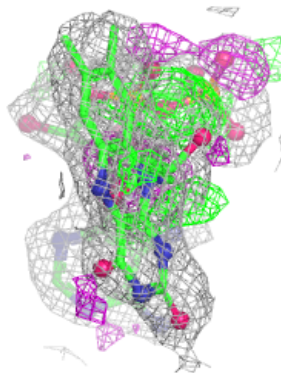
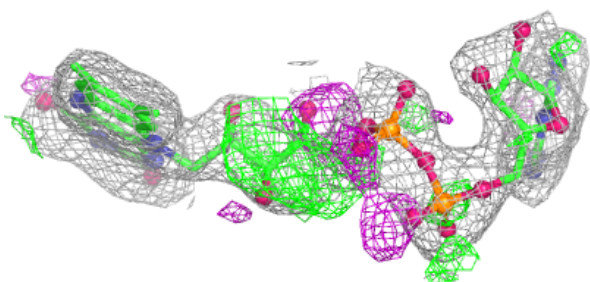
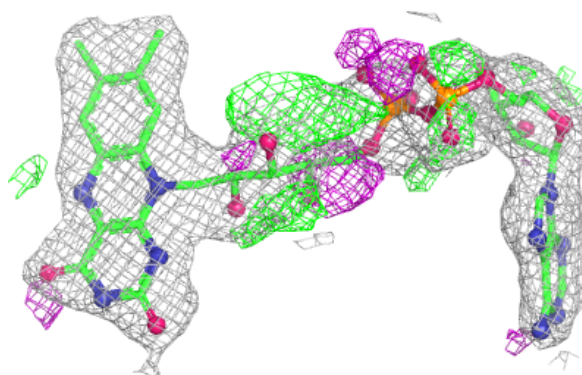
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD A 240:

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

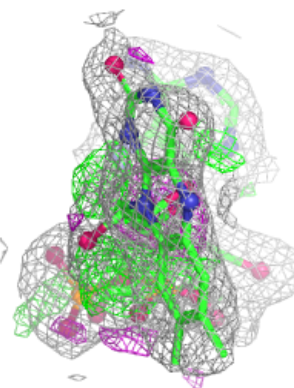
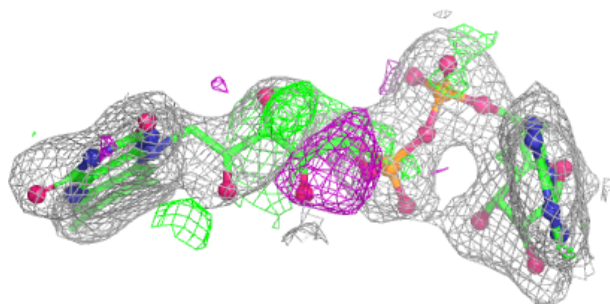
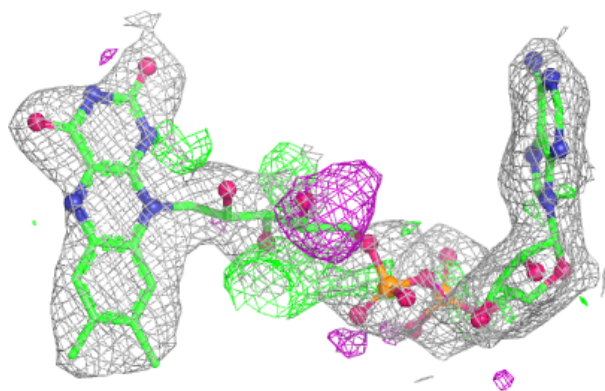
**Electron density around FAD B 240:**

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

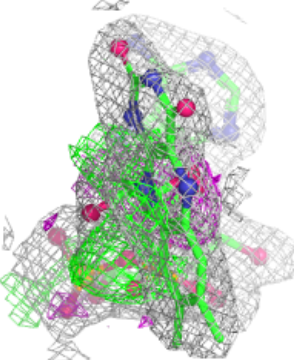
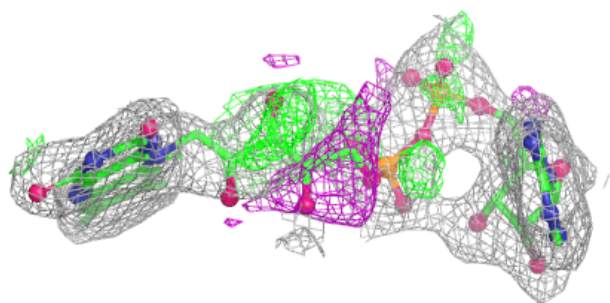
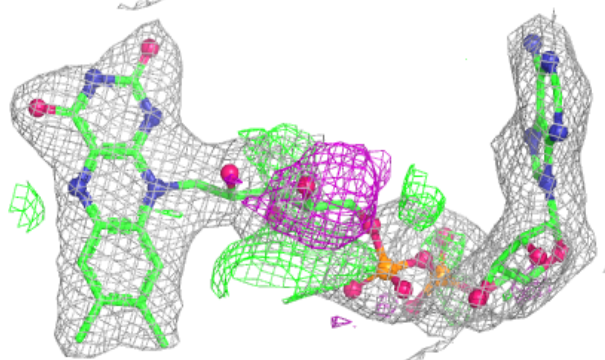


Electron density around FAD F 240:

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

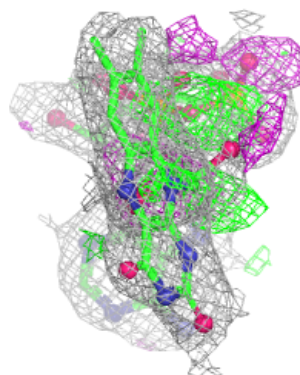
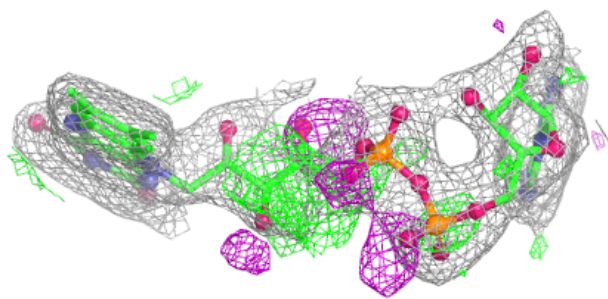
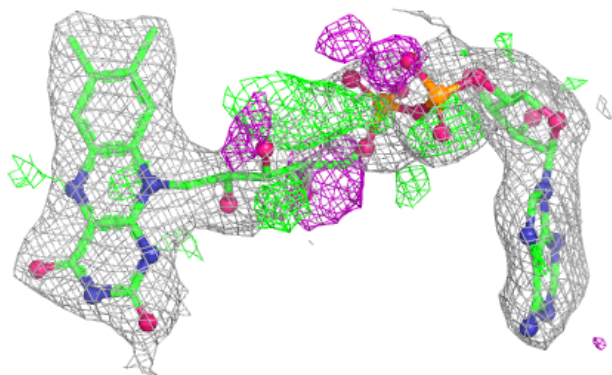
**Electron density around FAD C 240:**

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

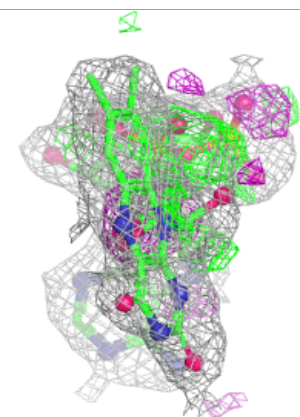
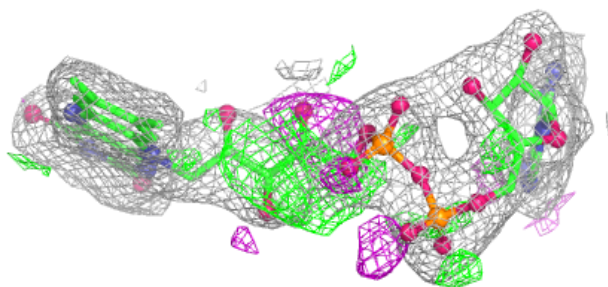
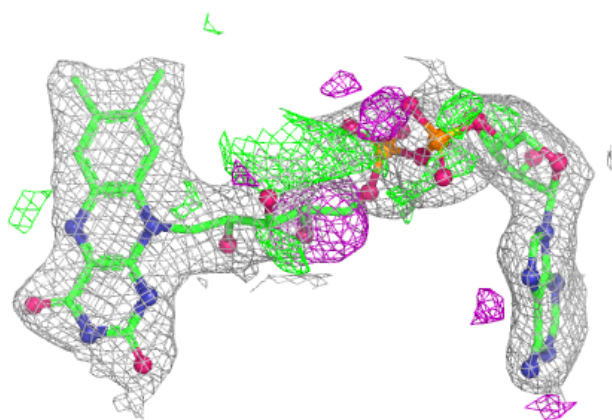


Electron density around FAD E 240:

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

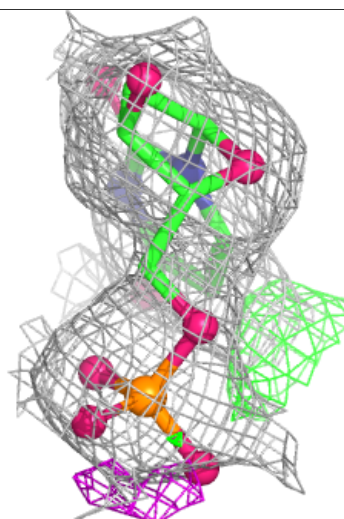
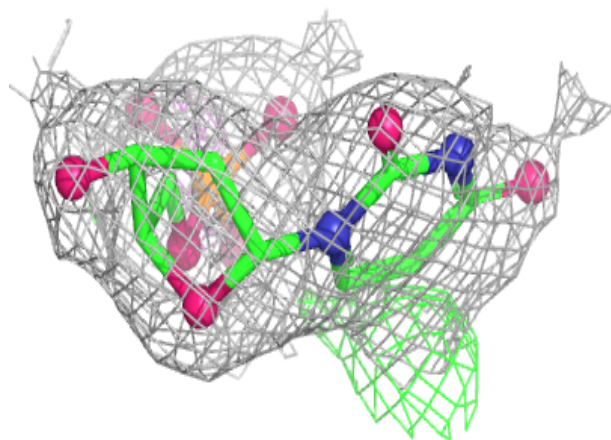
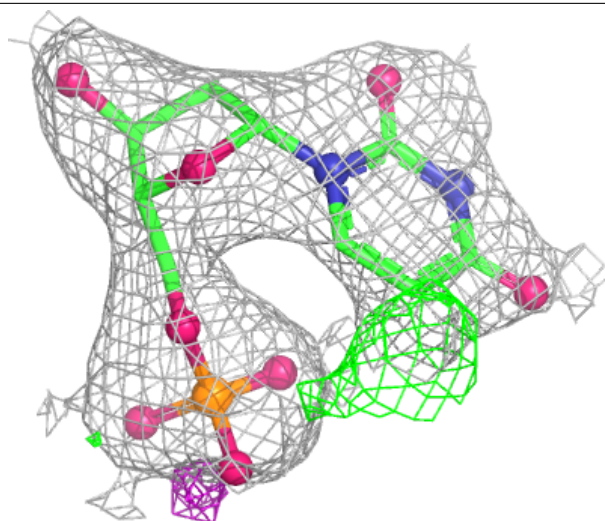
**Electron density around FAD D 240:**

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



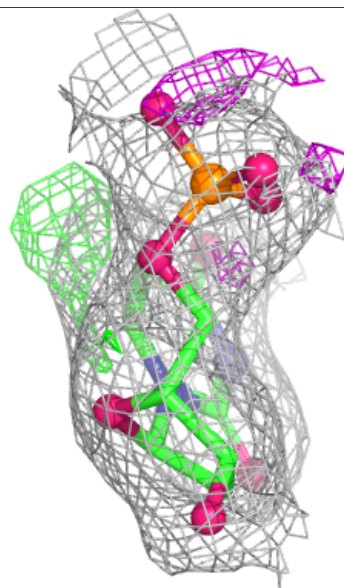
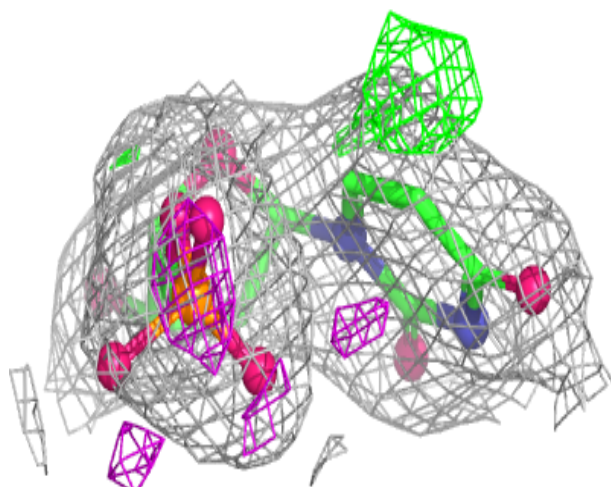
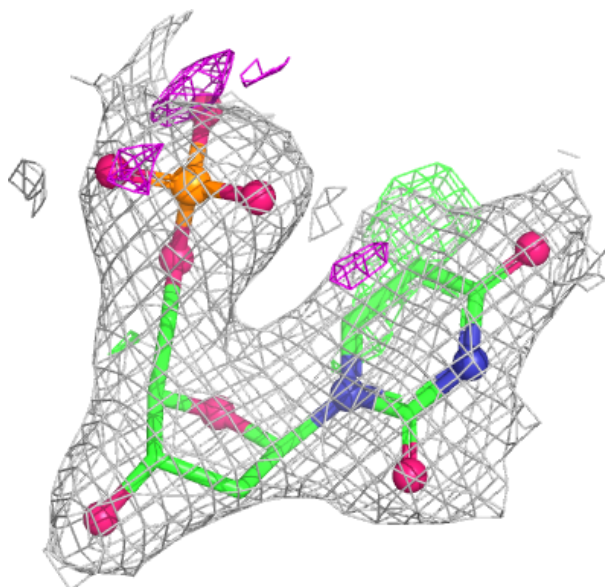
Electron density around UMP D 241:

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



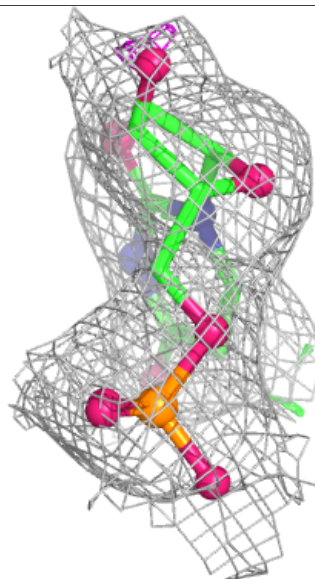
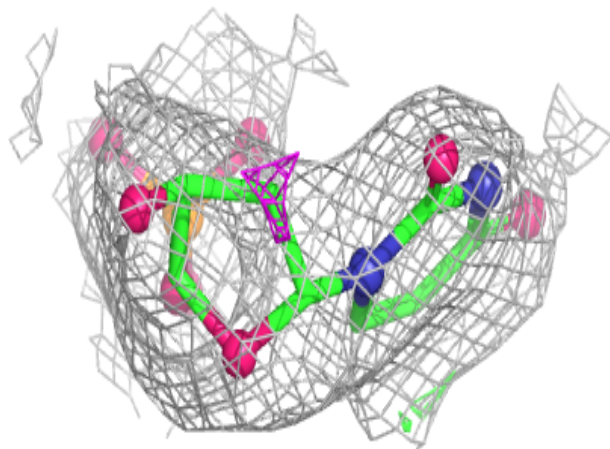
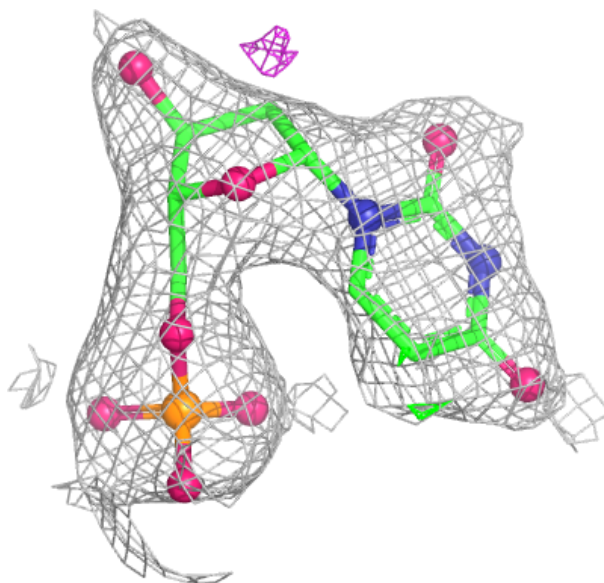
Electron density around UMP E 241:

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



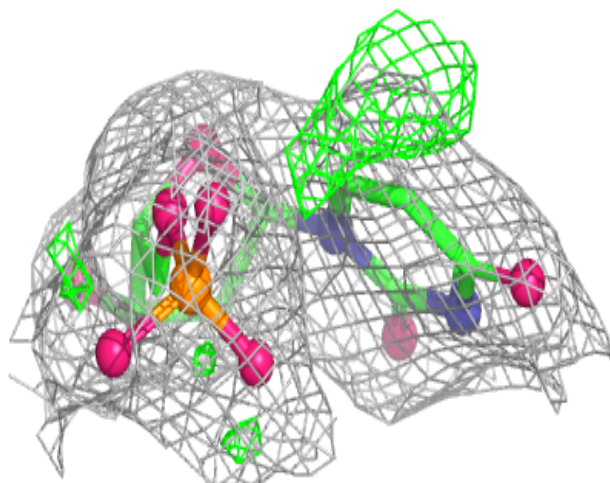
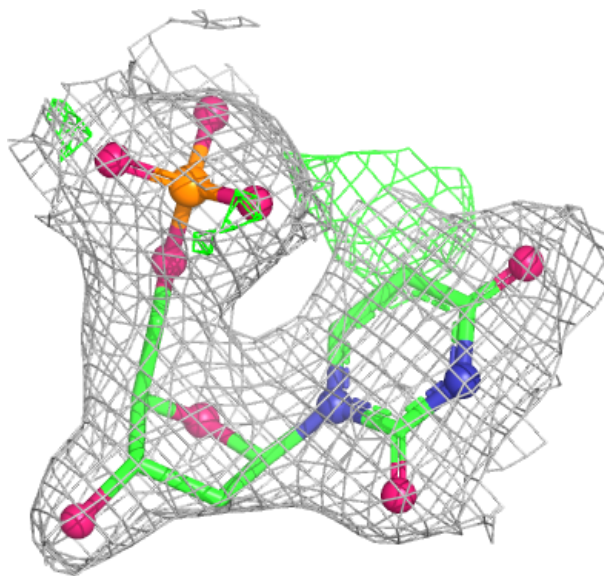
Electron density around UMP F 241:

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



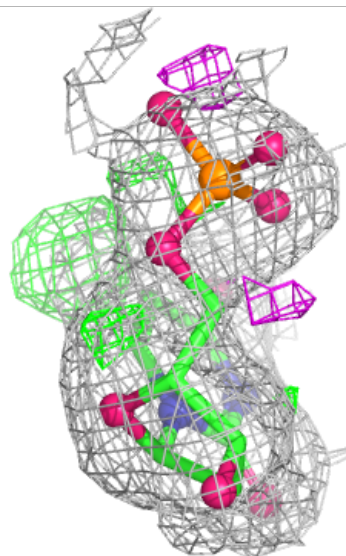
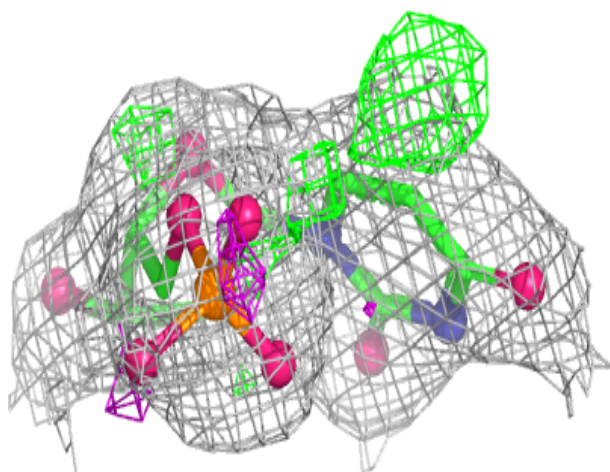
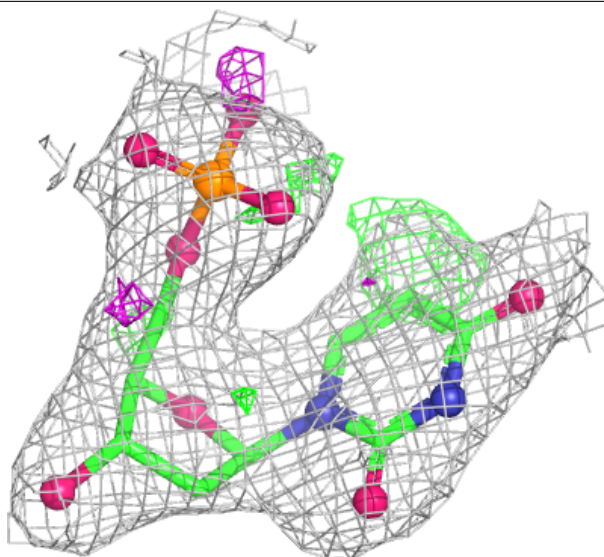
Electron density around UMP C 241:

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



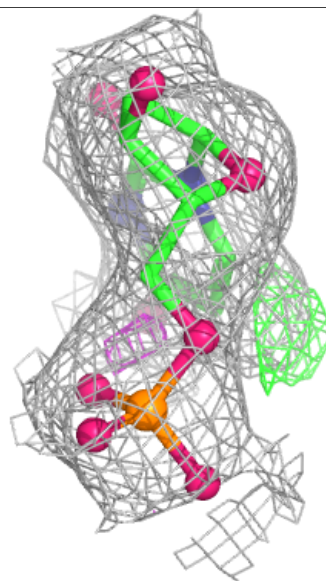
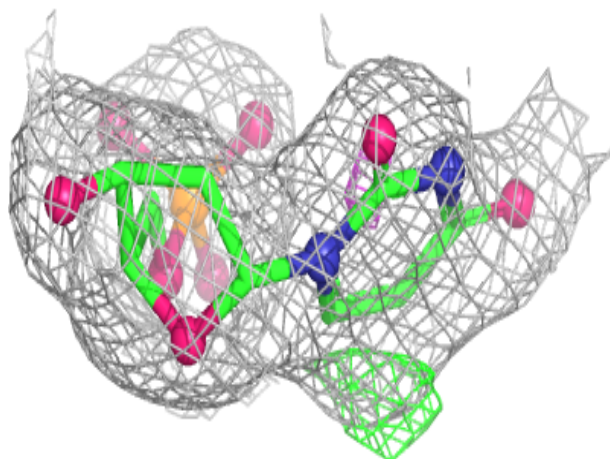
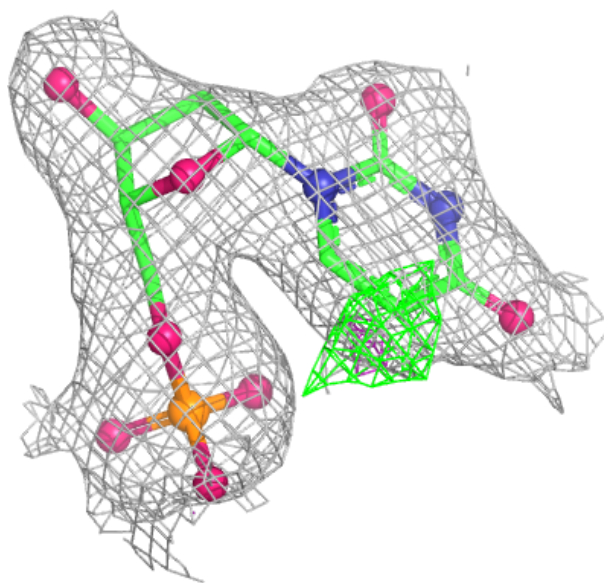
Electron density around UMP A 241:

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



Electron density around UMP B 241:

$2mF_o-DF_c$ (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.