Full wwPDB X-ray Structure Validation Report ⓘ

Nov 23, 2023 – 02:34 AM JST

PDB ID	:	7YTV
Title	:	Crystal structure of Aspergillus fumigatus Thioredoxin reductase in complex with auranofin
Authors	:	Lin, W.
Deposited on	:	2022-08-16
Resolution	:	3.87 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

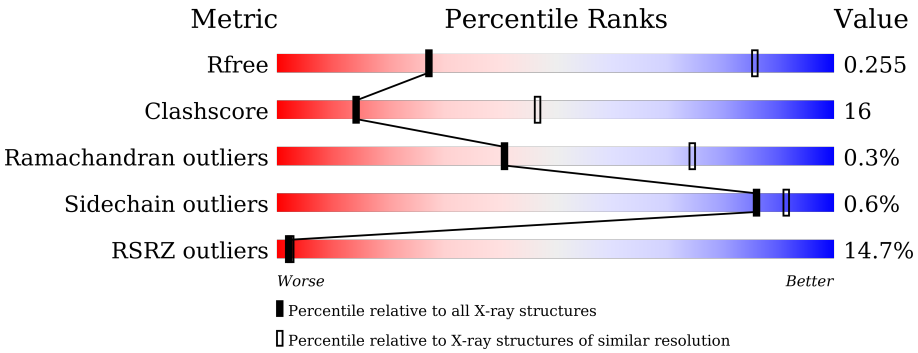
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.87 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1048 (4.10-3.62)
Clashscore	141614	1015 (4.08-3.64)
Ramachandran outliers	138981	1069 (4.10-3.62)
Sidechain outliers	138945	1062 (4.10-3.62)
RSRZ outliers	127900	1206 (4.12-3.60)

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	373	12% 58% 26% 15%
1	B	373	13% 62% 23% . 14%

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
3	FAD	A	402	-	-	-	X

2 Entry composition

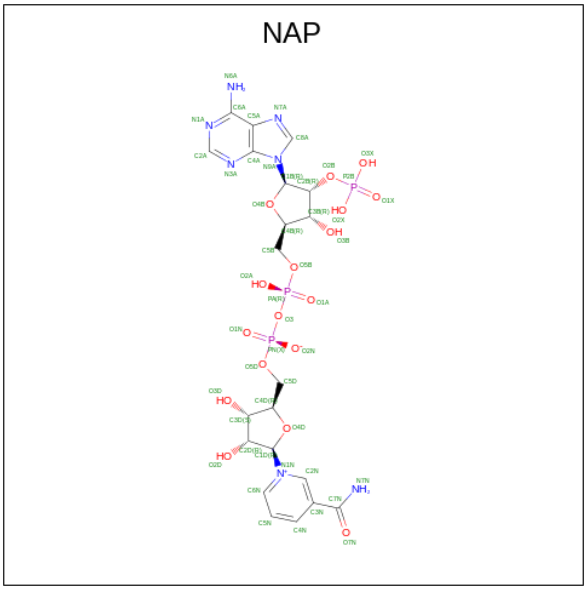
There are 4 unique types of molecules in this entry. The entry contains 4858 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thioredoxin reductase.

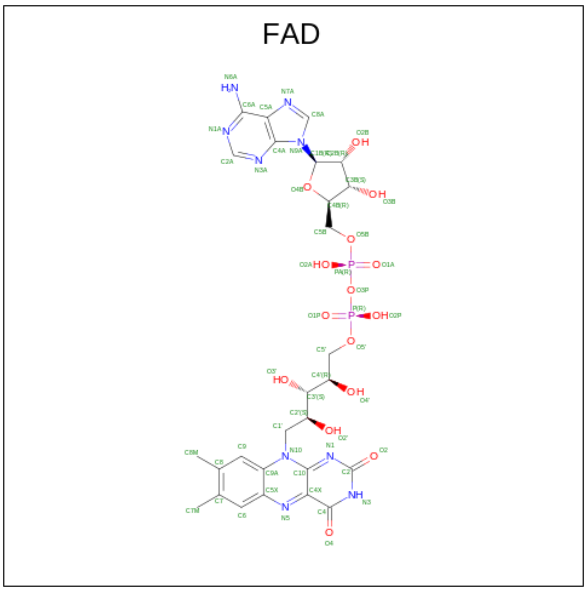
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2337	1467	402	458	10			
1	B	321	Total	C	N	O	S	0	0	0
			2352	1481	404	456	11			

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is water.

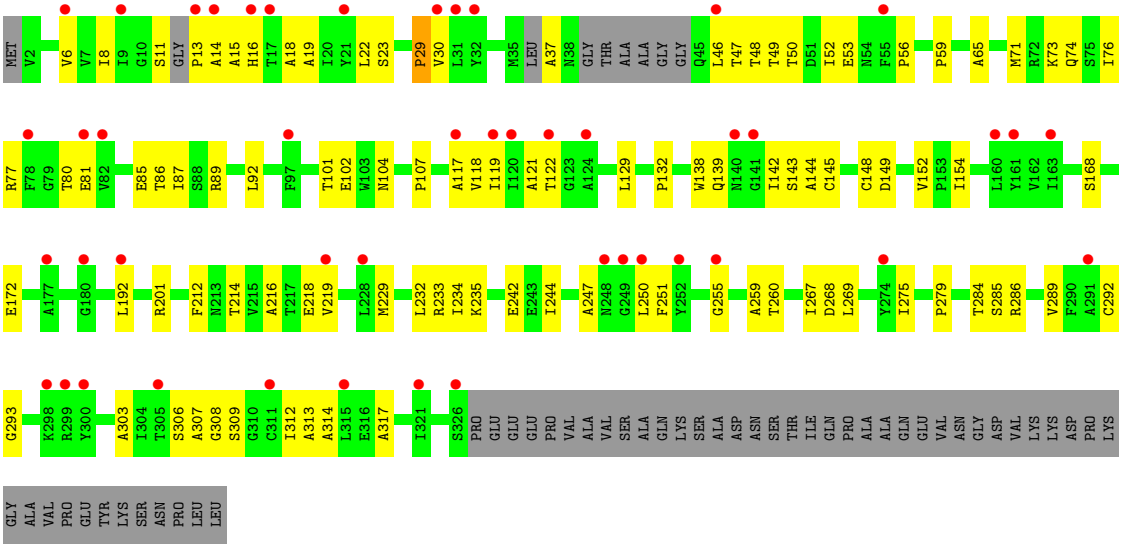
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		

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.

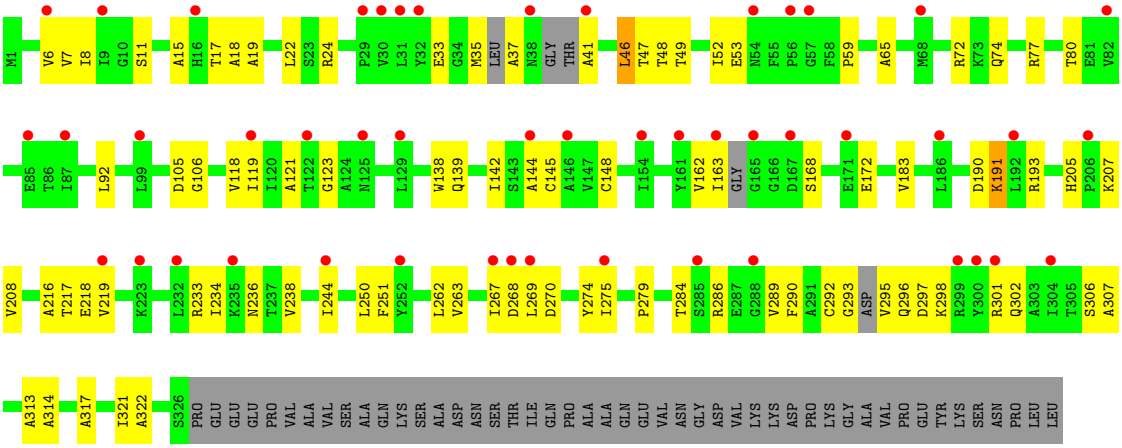
• Molecule 1: Thioredoxin reductase

Chain A:



• Molecule 1: Thioredoxin reductase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	152.81Å 152.81Å 135.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	59.43 – 3.87 132.34 – 3.87	Depositor EDS
% Data completeness (in resolution range)	99.6 (59.43-3.87) 99.7 (132.34-3.87)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.34 (at 3.89Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.230 , 0.245 0.241 , 0.255	Depositor DCC
R_{free} test set	840 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å^2)	139.5	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 122.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.049 for -h,-k,l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	4858	wwPDB-VP
Average B, all atoms (Å^2)	143.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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: NAP, FAD

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	0.24	0/2375	0.43	0/3225
1	B	0.24	0/2391	0.43	0/3249
All	All	0.24	0/4766	0.43	0/6474

There are no bond length outliers.

There are no bond angle outliers.

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	2337	0	2268	78	0
1	B	2352	0	2286	76	0
2	A	31	0	10	2	0
2	B	31	0	10	2	0
3	A	53	0	30	9	0
3	B	53	0	30	7	0
4	A	1	0	0	0	0
All	All	4858	0	4634	146	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 16.

All (146) 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:46:LEU:HD13	1:B:52:ILE:HD11	1.17	1.16
1:B:46:LEU:CD1	1:B:52:ILE:CD1	2.25	1.14
1:B:46:LEU:CD1	1:B:52:ILE:HD11	1.78	1.12
1:A:13:PRO:HD2	3:A:402:FAD:O1P	1.48	1.11
1:A:46:LEU:HD22	1:A:52:ILE:HD11	1.36	1.07
1:A:14:ALA:HB2	1:A:306:SER:O	1.54	1.06
1:B:46:LEU:HD13	1:B:52:ILE:CD1	1.85	1.04
1:B:145:CYS:HG	1:B:148:CYS:HG	1.08	1.00
1:B:46:LEU:HD11	1:B:52:ILE:CD1	1.98	0.93
1:A:122:THR:HG21	1:A:259:ALA:O	1.68	0.93
1:A:15:ALA:HB2	1:A:121:ALA:HB3	1.53	0.90
1:B:46:LEU:HD11	1:B:52:ILE:HD13	1.55	0.88
1:A:14:ALA:CB	1:A:306:SER:O	2.22	0.87
1:A:46:LEU:HD22	1:A:52:ILE:CD1	2.05	0.85
1:A:14:ALA:HB2	1:A:306:SER:C	2.00	0.79
1:A:19:ALA:HB1	1:A:80:THR:HG21	1.66	0.77
1:A:14:ALA:HA	1:A:307:ALA:HA	1.70	0.73
1:B:49:THR:HA	1:B:139:GLN:HG3	1.72	0.71
1:A:89:ARG:HH22	1:A:107:PRO:HA	1.56	0.70
1:A:46:LEU:HG	3:A:402:FAD:O3'	1.90	0.69
1:B:190:ASP:OD1	1:B:191:LYS:HG3	1.94	0.68
1:A:216:ALA:HB1	1:A:232:LEU:HD11	1.75	0.68
1:A:139:GLN:NE2	1:A:149:ASP:OD2	2.26	0.67
1:A:48:THR:HB	1:A:138:TRP:HE1	1.61	0.66
1:A:279:PRO:HB2	1:B:279:PRO:HB2	1.78	0.65
1:B:47:THR:HG22	1:B:65:ALA:HA	1.79	0.65
1:B:317:ALA:O	1:B:321:ILE:HG12	1.97	0.64
1:A:49:THR:HA	1:A:139:GLN:HG3	1.80	0.64
1:B:275:ILE:HB	1:B:295:VAL:HG12	1.80	0.64
1:A:101:THR:HB	1:A:104:ASN:HB2	1.80	0.63
1:A:122:THR:CG2	1:A:259:ALA:O	2.46	0.63
1:B:193:ARG:NH1	2:B:401:NAP:O2A	2.32	0.63
1:B:92:LEU:O	1:B:286:ARG:NH1	2.32	0.62
1:A:59:PRO:O	1:B:77:ARG:NH1	2.34	0.61
1:A:50:THR:HG23	1:A:139:GLN:HB3	1.84	0.60
1:A:92:LEU:O	1:A:286:ARG:NH1	2.33	0.60
1:A:13:PRO:CD	3:A:402:FAD:O1P	2.38	0.60
1:A:143:SER:OG	1:A:144:ALA:N	2.35	0.60
1:B:148:CYS:SG	1:B:302:GLN:NE2	2.75	0.60
1:A:6:VAL:HG22	1:A:117:ALA:HB3	1.83	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ALA:HB3	3:A:402:FAD:HM83	1.84	0.59
1:B:163:ILE:HG13	1:B:250:LEU:HD11	1.84	0.59
1:B:6:VAL:HG11	1:B:22:LEU:HD13	1.83	0.59
1:B:193:ARG:NH2	2:B:401:NAP:O1X	2.35	0.59
1:A:119:ILE:HD11	1:A:317:ALA:HB2	1.85	0.58
1:A:18:ALA:HA	1:A:314:ALA:HB2	1.86	0.58
1:A:47:THR:HG22	1:A:65:ALA:HA	1.85	0.58
1:B:270:ASP:OD2	1:B:298:LYS:NZ	2.37	0.57
1:A:77:ARG:NH1	1:B:59:PRO:O	2.36	0.57
1:B:8:ILE:HD11	1:B:22:LEU:HD12	1.85	0.57
1:A:148:CYS:O	1:B:24:ARG:NH1	2.38	0.56
1:A:15:ALA:HB2	1:A:121:ALA:CB	2.31	0.56
1:A:86:THR:HB	1:A:102:GLU:HB2	1.86	0.56
1:B:275:ILE:HB	1:B:295:VAL:CG1	2.35	0.56
1:B:297:ASP:OD2	1:B:301:ARG:N	2.32	0.56
1:A:30:VAL:HG12	1:A:81:GLU:HB3	1.87	0.56
1:B:48:THR:HB	1:B:138:TRP:HE1	1.71	0.56
1:B:267:ILE:HD11	1:B:289:VAL:HG21	1.88	0.55
1:B:15:ALA:HB2	1:B:121:ALA:HB3	1.87	0.55
1:B:144:ALA:HB3	3:B:402:FAD:HM83	1.89	0.55
1:B:46:LEU:CD1	1:B:52:ILE:HD13	2.16	0.54
1:B:119:ILE:HD11	1:B:317:ALA:HB2	1.89	0.54
1:A:201:ARG:HH21	1:B:322:ALA:HB1	1.73	0.54
1:B:11:SER:H	3:B:402:FAD:H4B	1.72	0.54
1:B:105:ASP:OD1	1:B:106:GLY:N	2.41	0.54
1:B:145:CYS:HB3	1:B:148:CYS:SG	2.48	0.54
1:A:267:ILE:HD11	1:A:289:VAL:HG11	1.90	0.53
1:B:11:SER:OG	1:B:33:GLU:O	2.26	0.52
1:B:142:ILE:HD11	1:B:219:VAL:HG11	1.92	0.52
1:A:308:GLY:O	1:A:312:ILE:HG12	2.10	0.52
1:B:33:GLU:OE1	1:B:72:ARG:NH1	2.42	0.52
1:A:8:ILE:HD11	1:A:22:LEU:HD22	1.92	0.52
1:B:7:VAL:O	1:B:118:VAL:HA	2.09	0.52
1:B:205:HIS:HE1	1:B:207:LYS:HB2	1.76	0.51
1:B:263:VAL:HG21	1:B:269:LEU:HD21	1.92	0.51
1:B:18:ALA:HA	1:B:314:ALA:HB2	1.93	0.51
1:A:23:SER:HB3	1:A:80:THR:HG23	1.93	0.50
1:A:37:ALA:HB1	1:A:86:THR:HG23	1.93	0.50
1:A:168:SER:O	1:A:172:GLU:HG3	2.12	0.50
1:A:73:LYS:HA	1:A:76:ILE:HG12	1.95	0.49
1:B:217:THR:N	1:B:233:ARG:O	2.44	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:THR:HG21	1:B:307:ALA:HB1	1.95	0.49
1:A:85:GLU:HB3	1:A:101:THR:HG23	1.95	0.48
1:A:142:ILE:HD11	1:A:219:VAL:HG11	1.96	0.48
1:B:216:ALA:HA	1:B:234:ILE:HA	1.96	0.47
1:A:56:PRO:HB3	1:B:17:THR:HG22	1.97	0.47
1:B:145:CYS:CB	1:B:148:CYS:SG	3.03	0.47
1:A:13:PRO:HD2	3:A:402:FAD:P	2.50	0.46
1:A:22:LEU:HD23	1:A:29:PRO:HB3	1.96	0.46
1:B:284:THR:OG1	1:B:289:VAL:O	2.24	0.46
1:A:303:ALA:HB2	3:A:402:FAD:H3'	1.97	0.46
1:B:293:GLY:HA2	1:B:306:SER:HA	1.98	0.46
1:B:296:GLN:HA	1:B:296:GLN:OE1	2.16	0.46
1:B:123:GLY:HA3	3:B:402:FAD:O2A	2.14	0.46
1:A:255:GLY:N	2:A:401:NAP:H52A	2.30	0.46
1:A:145:CYS:CB	1:A:148:CYS:SG	3.02	0.46
1:B:46:LEU:HD21	3:B:402:FAD:C10	2.46	0.46
1:B:190:ASP:C	1:B:191:LYS:HG3	2.36	0.45
1:B:290:PHE:HE1	1:B:317:ALA:HA	1.80	0.45
1:A:11:SER:H	3:A:402:FAD:H4B	1.80	0.45
1:B:19:ALA:HB1	1:B:80:THR:HG21	1.99	0.45
1:B:168:SER:O	1:B:172:GLU:HG3	2.16	0.45
1:A:233:ARG:HE	1:A:244:ILE:HD11	1.82	0.44
1:A:53:GLU:O	1:B:74:GLN:NE2	2.50	0.44
1:A:284:THR:OG1	1:A:289:VAL:HG13	2.16	0.44
1:B:205:HIS:CE1	1:B:207:LYS:HB2	2.52	0.44
1:B:292:CYS:HB3	1:B:313:ALA:HB2	1.99	0.44
1:A:118:VAL:HG13	1:A:289:VAL:HG23	1.99	0.44
1:A:214:THR:HG21	1:A:234:ILE:HD12	2.00	0.44
1:A:269:LEU:HD11	1:A:275:ILE:HG12	1.98	0.44
1:B:148:CYS:HB3	3:B:402:FAD:C4	2.47	0.44
1:B:8:ILE:HG12	1:B:119:ILE:HB	1.99	0.44
1:A:232:LEU:HD22	1:A:250:LEU:HD12	2.00	0.44
1:A:119:ILE:HD12	1:A:313:ALA:HB1	2.00	0.44
1:A:292:CYS:HB3	1:A:313:ALA:HB2	2.00	0.44
1:B:37:ALA:HB2	1:B:41:ALA:N	2.33	0.44
1:A:13:PRO:CD	3:A:402:FAD:H5'2	2.47	0.44
1:B:270:ASP:OD1	1:B:274:TYR:N	2.51	0.44
1:B:218:GLU:OE2	1:B:233:ARG:NE	2.48	0.43
1:B:162:VAL:HG22	1:B:251:PHE:HB2	2.00	0.43
1:B:236:ASN:HD21	1:B:238:VAL:HG22	1.82	0.43
1:B:262:LEU:HD11	3:B:402:FAD:N6A	2.33	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:HIS:HB3	1:A:71:MET:HG2	1.99	0.43
1:B:268:ASP:N	1:B:268:ASP:OD1	2.51	0.43
1:A:129:LEU:HD23	1:A:129:LEU:HA	1.86	0.43
1:A:268:ASP:HB2	1:A:285:SER:HB2	2.00	0.43
1:B:262:LEU:HD11	3:B:402:FAD:H61A	1.83	0.43
1:A:29:PRO:HG2	1:A:80:THR:HG22	2.01	0.43
1:A:8:ILE:HG12	1:A:119:ILE:HB	2.00	0.43
1:B:233:ARG:HG3	1:B:244:ILE:HG12	2.00	0.42
1:A:192:LEU:HG	1:A:212:PHE:HE1	1.83	0.42
1:A:87:ILE:O	3:A:402:FAD:N6A	2.53	0.42
1:A:152:VAL:HG12	1:A:154:ILE:H	1.85	0.42
1:B:145:CYS:CB	1:B:148:CYS:HG	2.31	0.42
1:B:183:VAL:O	1:B:208:VAL:HA	2.19	0.42
1:A:74:GLN:NE2	1:B:53:GLU:O	2.53	0.41
1:A:121:ALA:HB2	1:A:292:CYS:SG	2.60	0.41
1:A:293:GLY:H	1:A:309:SER:HB3	1.85	0.41
1:A:132:PRO:HG2	1:A:218:GLU:HG3	2.02	0.41
1:A:14:ALA:HB1	1:A:306:SER:O	2.17	0.41
1:A:143:SER:HB3	1:A:251:PHE:CE1	2.55	0.41
1:A:235:LYS:HB2	1:A:242:GLU:HG2	2.01	0.41
1:B:190:ASP:OD1	1:B:191:LYS:N	2.54	0.41
1:A:229:MET:HE3	1:A:247:ALA:O	2.21	0.41
1:B:119:ILE:HD12	1:B:313:ALA:HB1	2.03	0.40
1:A:255:GLY:H	2:A:401:NAP:H52A	1.85	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	309/373 (83%)	285 (92%)	23 (7%)	1 (0%)	41 74

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	311/373 (83%)	293 (94%)	17 (6%)	1 (0%)	41 74
All	All	620/746 (83%)	578 (93%)	40 (6%)	2 (0%)	41 74

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	191	LYS
1	A	29	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	236/294 (80%)	235 (100%)	1 (0%)	91 94
1	B	235/294 (80%)	233 (99%)	2 (1%)	78 88
All	All	471/588 (80%)	468 (99%)	3 (1%)	86 91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	260	THR
1	B	35	MET
1	B	46	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	FAD	A	402	-	53,58,58	2.61	17 (32%)	68,89,89	1.95	17 (25%)
2	NAP	A	401	-	27,33,52	2.14	7 (25%)	35,52,80	2.35	5 (14%)
2	NAP	B	401	-	27,33,52	2.13	7 (25%)	35,52,80	2.39	6 (17%)
3	FAD	B	402	-	53,58,58	2.59	17 (32%)	68,89,89	1.95	16 (23%)

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	FAD	A	402	-	-	18/30/50/50	0/6/6/6
2	NAP	A	401	-	-	10/17/37/67	0/3/3/5
2	NAP	B	401	-	-	10/17/37/67	0/3/3/5
3	FAD	B	402	-	-	13/30/50/50	0/6/6/6

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	FAD	C4X-N5	7.58	1.45	1.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	FAD	C4X-N5	7.44	1.45	1.30
3	B	402	FAD	C10-N1	6.69	1.46	1.33
3	A	402	FAD	C10-N1	6.68	1.46	1.33
2	A	401	NAP	O4B-C4B	-6.15	1.31	1.45
2	B	401	NAP	O4B-C4B	-6.00	1.31	1.45
3	A	402	FAD	O4B-C4B	-5.98	1.31	1.45
3	B	402	FAD	O4B-C4B	-5.86	1.31	1.45
3	A	402	FAD	C5X-N5	5.29	1.49	1.39
3	A	402	FAD	C9A-N10	5.26	1.50	1.41
3	B	402	FAD	C9A-N10	5.24	1.50	1.41
3	B	402	FAD	C5X-N5	5.24	1.49	1.39
3	B	402	FAD	C2-N1	5.06	1.48	1.36
3	A	402	FAD	C2-N1	5.06	1.48	1.36
2	B	401	NAP	P2B-O2B	4.84	1.68	1.59
2	A	401	NAP	P2B-O2B	4.80	1.68	1.59
3	A	402	FAD	C2-N3	4.51	1.49	1.39
3	B	402	FAD	C2-N3	4.30	1.49	1.39
3	A	402	FAD	C4-N3	4.12	1.46	1.38
3	B	402	FAD	C10-N10	3.97	1.46	1.37
3	B	402	FAD	C4-N3	3.95	1.46	1.38
3	A	402	FAD	C10-N10	3.92	1.45	1.37
2	A	401	NAP	C2A-N3A	3.45	1.37	1.32
2	B	401	NAP	C2A-N3A	3.44	1.37	1.32
2	A	401	NAP	O3B-C3B	-3.37	1.35	1.43
2	B	401	NAP	O3B-C3B	-3.36	1.35	1.43
3	A	402	FAD	C6A-N6A	3.05	1.45	1.34
3	B	402	FAD	C6A-N6A	3.04	1.45	1.34
3	A	402	FAD	O3B-C3B	-3.00	1.35	1.43
3	B	402	FAD	O3B-C3B	-2.98	1.36	1.43
3	B	402	FAD	O2-C2	-2.96	1.18	1.24
3	B	402	FAD	O2B-C2B	2.93	1.49	1.43
2	A	401	NAP	PN-O5D	2.93	1.66	1.54
2	B	401	NAP	PN-O5D	2.92	1.66	1.54
3	A	402	FAD	O2B-C2B	2.91	1.49	1.43
3	A	402	FAD	O2-C2	-2.89	1.18	1.24
3	A	402	FAD	C2A-N3A	2.81	1.36	1.32
3	B	402	FAD	C5A-C4A	-2.79	1.33	1.40
3	B	402	FAD	C2A-N3A	2.79	1.36	1.32
3	A	402	FAD	C5A-C4A	-2.76	1.33	1.40
3	B	402	FAD	O4-C4	-2.56	1.18	1.23
3	A	402	FAD	O4-C4	-2.54	1.18	1.23
2	B	401	NAP	C5A-C4A	-2.43	1.34	1.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAP	C5A-C4A	-2.38	1.34	1.40
2	B	401	NAP	C6A-N6A	2.18	1.42	1.34
2	A	401	NAP	C6A-N6A	2.14	1.41	1.34
3	A	402	FAD	C4X-C4	2.06	1.52	1.44
3	B	402	FAD	C4X-C4	2.03	1.52	1.44

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	401	NAP	C5A-C6A-N6A	9.13	134.22	120.35
2	B	401	NAP	C5A-C6A-N6A	9.10	134.18	120.35
3	B	402	FAD	C7M-C7-C6	-7.85	104.97	119.49
3	A	402	FAD	C7M-C7-C6	-7.67	105.31	119.49
3	B	402	FAD	C7M-C7-C8	6.93	134.95	120.74
3	A	402	FAD	C7M-C7-C8	6.81	134.70	120.74
2	B	401	NAP	N6A-C6A-N1A	-6.11	105.90	118.57
2	A	401	NAP	N6A-C6A-N1A	-6.06	105.99	118.57
2	B	401	NAP	N3A-C2A-N1A	-5.54	120.02	128.68
3	B	402	FAD	N3A-C2A-N1A	-5.45	120.16	128.68
3	A	402	FAD	N3A-C2A-N1A	-5.35	120.32	128.68
2	A	401	NAP	N3A-C2A-N1A	-5.31	120.38	128.68
2	B	401	NAP	C1B-N9A-C4A	-4.90	118.03	126.64
2	A	401	NAP	C1B-N9A-C4A	-4.74	118.31	126.64
3	A	402	FAD	C4-N3-C2	-3.57	119.05	125.64
3	B	402	FAD	N6A-C6A-N1A	-3.54	111.23	118.57
3	A	402	FAD	N6A-C6A-N1A	-3.43	111.45	118.57
3	B	402	FAD	C4-N3-C2	-3.32	119.50	125.64
2	A	401	NAP	PA-O3-PN	-2.82	123.15	132.83
3	A	402	FAD	C4X-C4-N3	2.67	119.98	113.19
3	B	402	FAD	C4X-C4-N3	2.58	119.75	113.19
3	A	402	FAD	P-O3P-PA	-2.55	124.08	132.83
3	B	402	FAD	P-O3P-PA	-2.50	124.25	132.83
2	B	401	NAP	PA-O3-PN	-2.49	124.28	132.83
3	A	402	FAD	O4-C4-C4X	-2.48	120.03	126.60
3	A	402	FAD	C4X-C10-N10	2.47	120.09	116.48
3	A	402	FAD	C8M-C8-C9	2.46	124.04	119.49
3	B	402	FAD	C4X-C10-N10	2.46	120.07	116.48
3	B	402	FAD	O4-C4-C4X	-2.42	120.17	126.60
3	A	402	FAD	C10-C4X-N5	-2.39	119.80	124.86
2	B	401	NAP	O4B-C1B-C2B	2.35	110.66	106.59
3	B	402	FAD	C8M-C8-C9	2.31	123.76	119.49
3	B	402	FAD	C10-C4X-N5	-2.27	120.04	124.86

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	402	FAD	C5A-C6A-N1A	2.26	125.47	120.35
3	B	402	FAD	C5'-C4'-C3'	-2.23	107.90	112.20
3	A	402	FAD	C5A-C6A-N1A	2.23	125.41	120.35
3	A	402	FAD	C4'-C3'-C2'	-2.20	108.78	113.36
3	A	402	FAD	C4X-C10-N1	-2.20	119.62	124.73
3	A	402	FAD	C8M-C8-C7	-2.18	116.28	120.74
3	A	402	FAD	C4-C4X-C10	2.17	120.44	116.79
3	B	402	FAD	C4X-C10-N1	-2.16	119.72	124.73
3	B	402	FAD	C4-C4X-C10	2.14	120.38	116.79
3	B	402	FAD	C8M-C8-C7	-2.03	116.58	120.74
3	A	402	FAD	C9A-C5X-N5	-2.00	120.25	122.43

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAP	C5B-O5B-PA-O1A
2	A	401	NAP	C5B-O5B-PA-O2A
2	A	401	NAP	C5B-O5B-PA-O3
2	A	401	NAP	PN-O3-PA-O5B
2	B	401	NAP	C5B-O5B-PA-O1A
2	B	401	NAP	C5B-O5B-PA-O2A
2	B	401	NAP	C5B-O5B-PA-O3
2	B	401	NAP	O4B-C4B-C5B-O5B
3	A	402	FAD	C5B-O5B-PA-O1A
3	A	402	FAD	N10-C1'-C2'-O2'
3	A	402	FAD	N10-C1'-C2'-C3'
3	A	402	FAD	C1'-C2'-C3'-O3'
3	A	402	FAD	C1'-C2'-C3'-C4'
3	A	402	FAD	O4'-C4'-C5'-O5'
3	A	402	FAD	C5'-O5'-P-O1P
3	A	402	FAD	C5'-O5'-P-O2P
3	A	402	FAD	PA-O3P-P-O5'
3	B	402	FAD	C5B-O5B-PA-O1A
3	B	402	FAD	C5B-O5B-PA-O2A
3	B	402	FAD	C3B-C4B-C5B-O5B
3	B	402	FAD	N10-C1'-C2'-O2'
3	B	402	FAD	N10-C1'-C2'-C3'
3	B	402	FAD	C1'-C2'-C3'-O3'
3	B	402	FAD	C1'-C2'-C3'-C4'
3	B	402	FAD	O4B-C4B-C5B-O5B
3	A	402	FAD	C2'-C3'-C4'-O4'

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	B	401	NAP	C3B-C2B-O2B-P2B
3	A	402	FAD	O3'-C3'-C4'-O4'
2	B	401	NAP	C3B-C4B-C5B-O5B
3	A	402	FAD	O2'-C2'-C3'-C4'
3	A	402	FAD	O3'-C3'-C4'-C5'
3	A	402	FAD	C2'-C3'-C4'-C5'
3	A	402	FAD	O2'-C2'-C3'-O3'
2	B	401	NAP	C1B-C2B-O2B-P2B
3	A	402	FAD	C3'-C4'-C5'-O5'
2	A	401	NAP	C3B-C2B-O2B-P2B
3	B	402	FAD	PA-O3P-P-O1P
3	B	402	FAD	PA-O3P-P-O5'
2	A	401	NAP	C2B-O2B-P2B-O1X
2	A	401	NAP	PA-O3-PN-O2N
2	A	401	NAP	C2B-O2B-P2B-O2X
3	B	402	FAD	O2'-C2'-C3'-C4'
2	B	401	NAP	PN-O3-PA-O1A
3	A	402	FAD	O4B-C4B-C5B-O5B
3	B	402	FAD	O2'-C2'-C3'-O3'
2	A	401	NAP	O4B-C4B-C5B-O5B
2	B	401	NAP	PN-O3-PA-O5B
2	B	401	NAP	C2B-O2B-P2B-O1X
2	A	401	NAP	C1B-C2B-O2B-P2B
3	A	402	FAD	C5'-O5'-P-O3P
3	B	402	FAD	C5B-O5B-PA-O3P

There are no ring outliers.

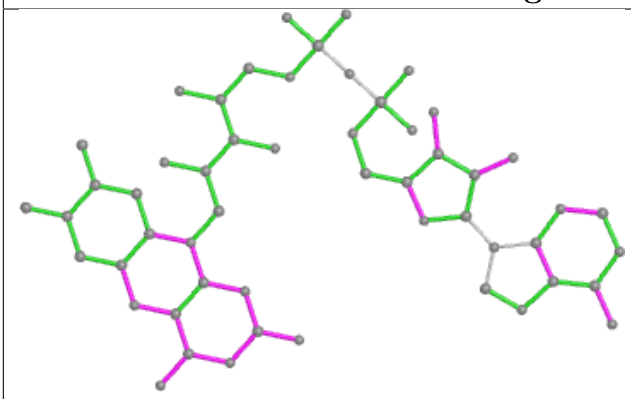
4 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	FAD	9	0
2	A	401	NAP	2	0
2	B	401	NAP	2	0
3	B	402	FAD	7	0

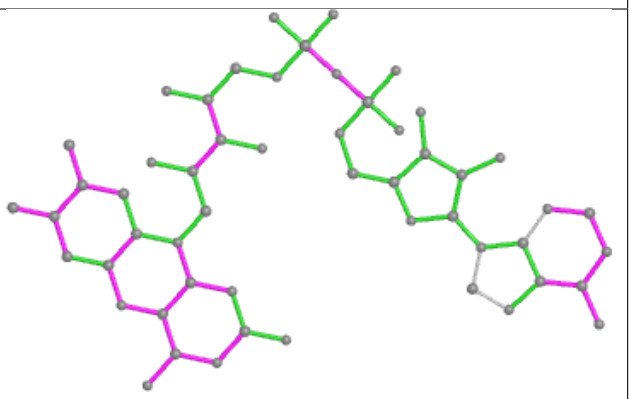
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

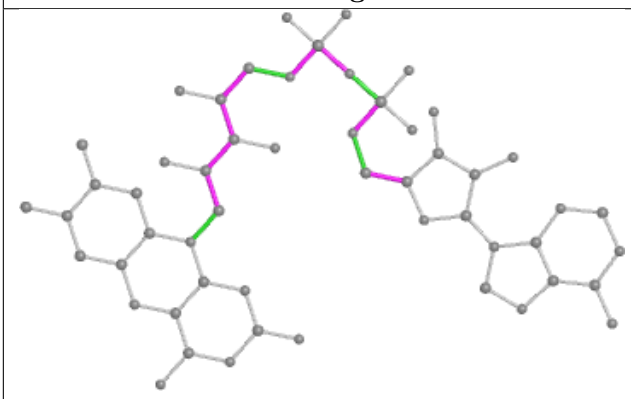
Ligand FAD A 402



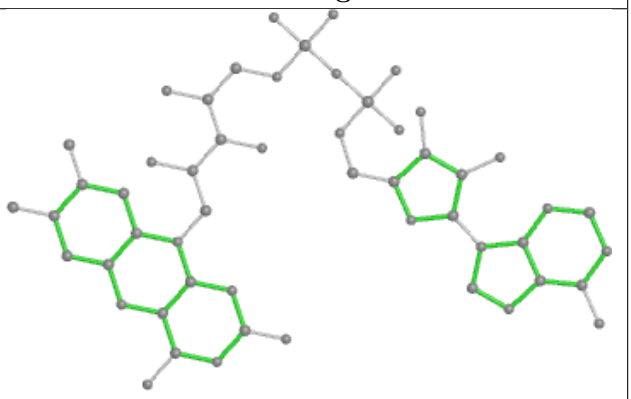
Bond lengths



Bond angles

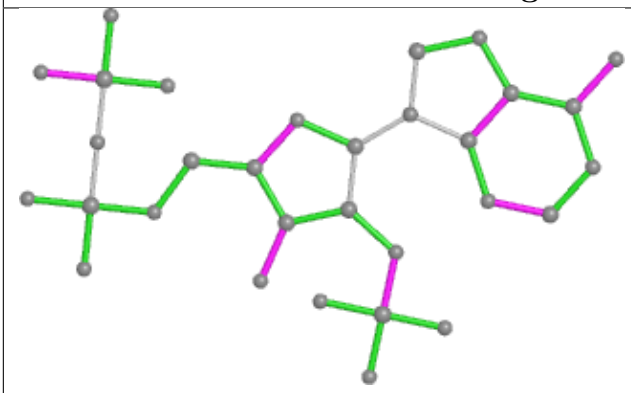


Torsions

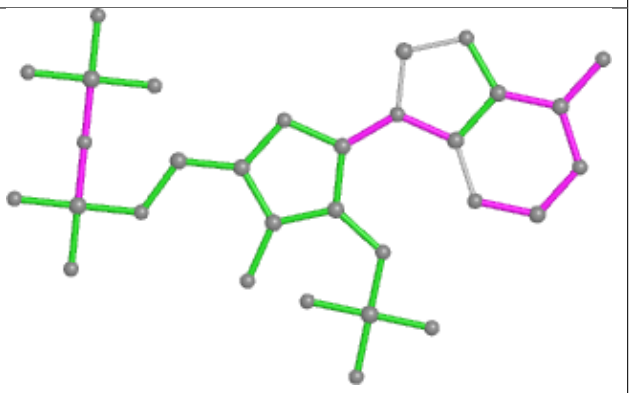


Rings

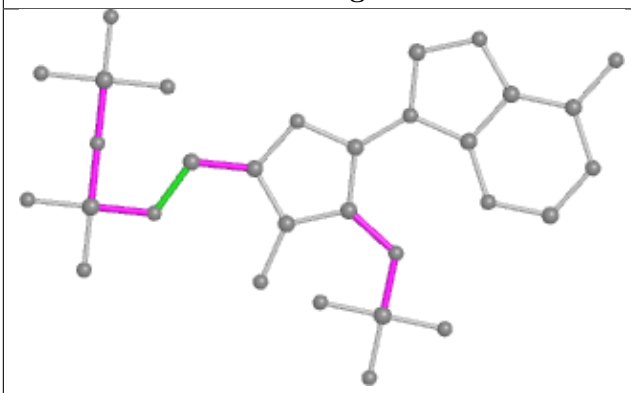
Ligand NAP A 401



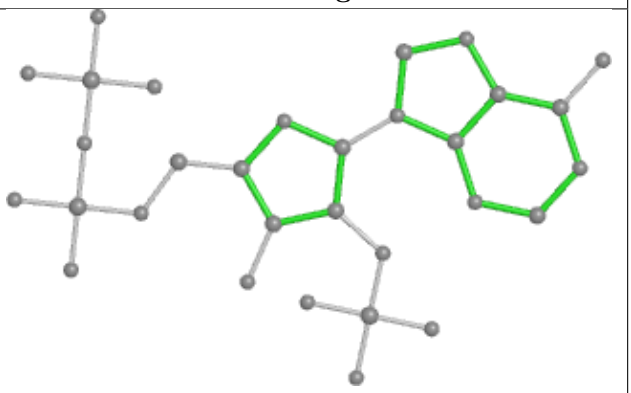
Bond lengths



Bond angles

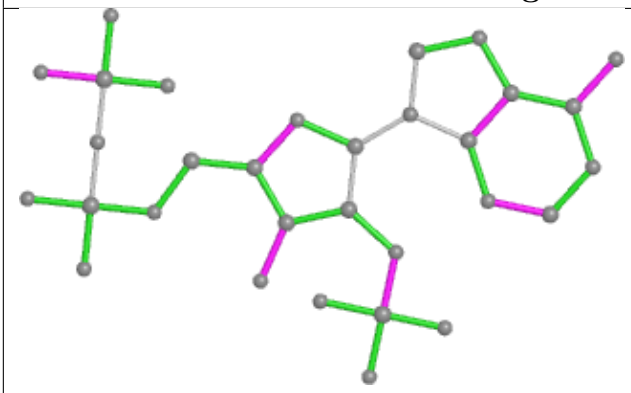


Torsions

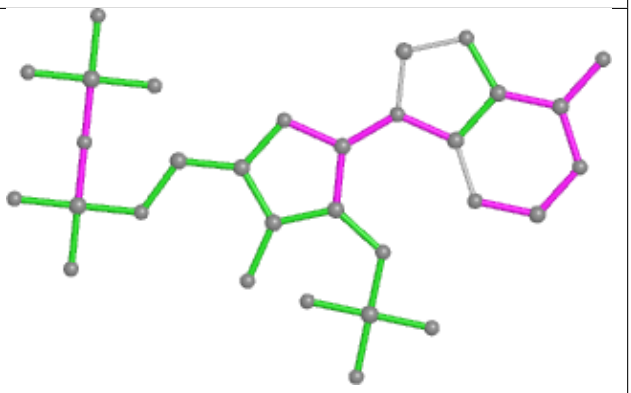


Rings

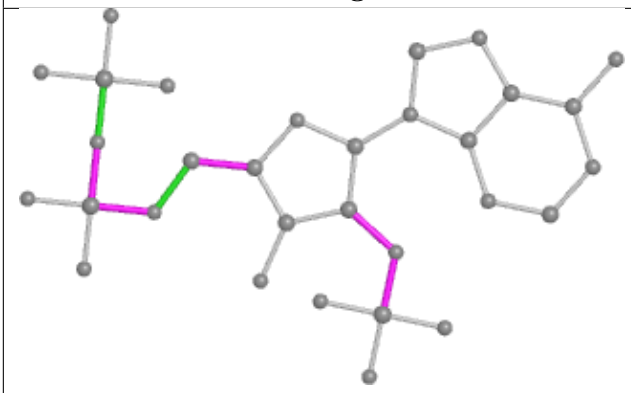
Ligand NAP B 401



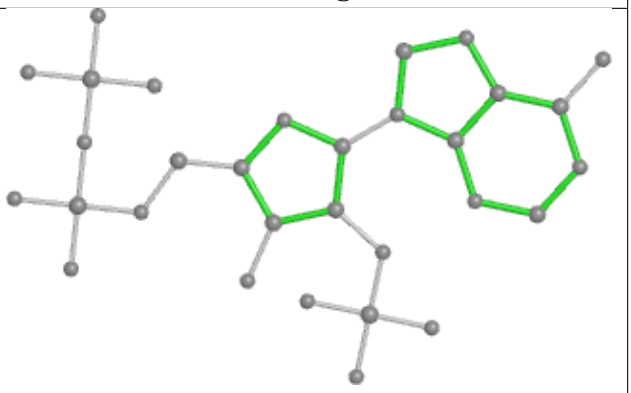
Bond lengths



Bond angles

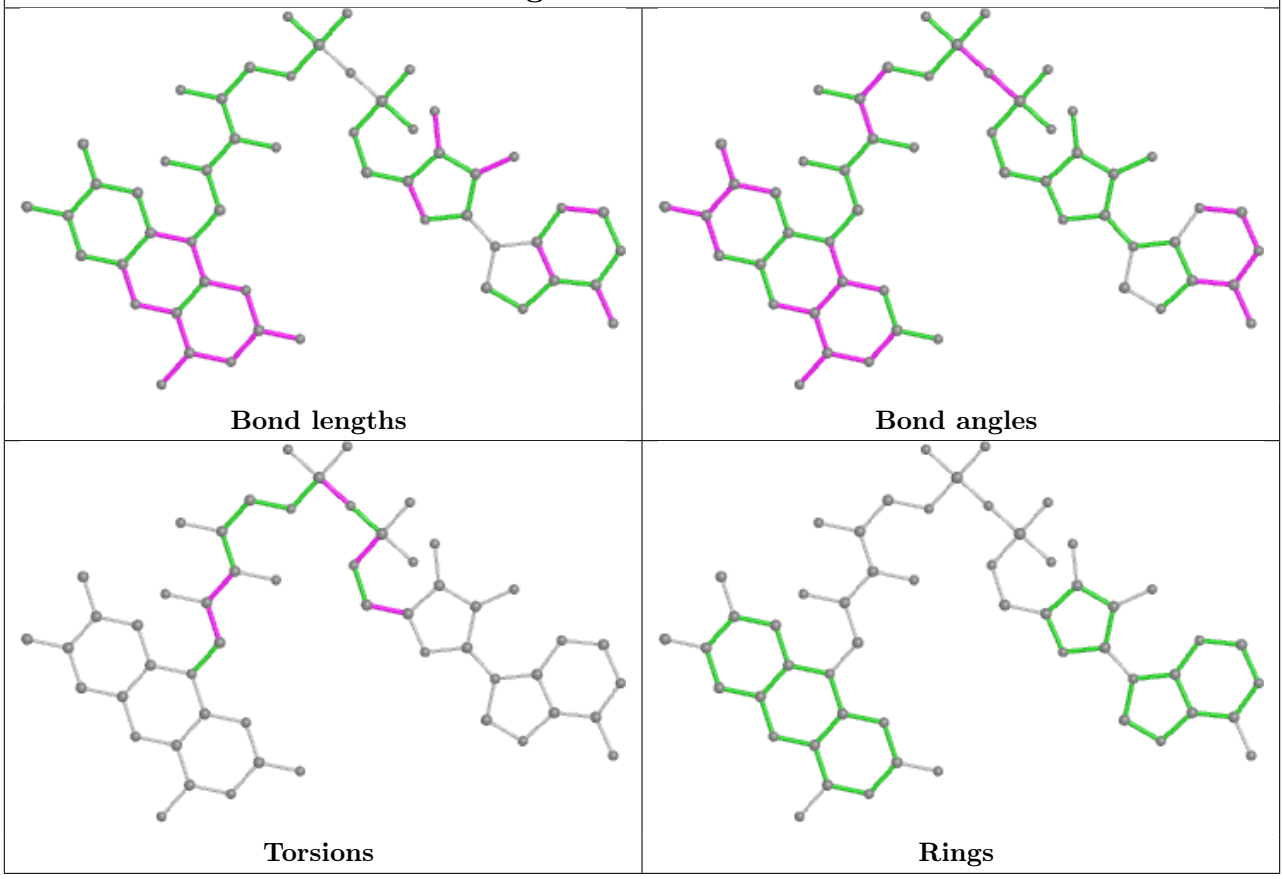


Torsions



Rings

Ligand FAD B 402



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	317/373 (84%)	1.01	46 (14%)	2	2	90, 140, 175, 218	0
1	B	321/373 (86%)	1.12	48 (14%)	2	2	84, 140, 180, 222	0
All	All	638/746 (85%)	1.07	94 (14%)	2	2	84, 140, 178, 222	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	301	ARG	4.9
1	B	192	LEU	4.6
1	B	54	ASN	4.5
1	B	244	ILE	4.1
1	B	32	TYR	3.9
1	A	31	LEU	3.7
1	B	223	LYS	3.4
1	A	192	LEU	3.4
1	A	117	ALA	3.4
1	B	9	ILE	3.4
1	B	163	ILE	3.4
1	B	56	PRO	3.3
1	B	186	LEU	3.3
1	A	291	ALA	3.3
1	A	17	THR	3.3
1	B	30	VAL	3.2
1	B	299	ARG	3.2
1	B	87	ILE	3.2
1	A	177	ALA	3.1
1	A	141	GLY	3.1
1	B	6	VAL	3.1
1	A	248	ASN	3.1
1	B	300	TYR	3.0
1	B	269	LEU	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	250	LEU	2.9
1	A	9	ILE	2.9
1	B	275	ILE	2.9
1	B	99	LEU	2.8
1	B	165	GLY	2.8
1	A	300	TYR	2.8
1	B	41	ALA	2.8
1	A	161	TYR	2.8
1	A	299	ARG	2.8
1	A	55	PHE	2.7
1	A	13	PRO	2.7
1	A	255	GLY	2.7
1	B	144	ALA	2.7
1	A	249	GLY	2.7
1	B	232	LEU	2.7
1	B	304	ILE	2.7
1	A	321	ILE	2.7
1	B	171	GLU	2.6
1	B	119	ILE	2.6
1	A	30	VAL	2.6
1	B	235	LYS	2.6
1	B	31	LEU	2.6
1	B	122	THR	2.6
1	A	119	ILE	2.5
1	A	46	LEU	2.5
1	B	16	HIS	2.5
1	A	16	HIS	2.5
1	A	305	THR	2.5
1	B	57	GLY	2.5
1	A	81	GLU	2.5
1	A	298	LYS	2.4
1	A	219	VAL	2.4
1	B	146	ALA	2.4
1	A	140	ASN	2.4
1	B	167	ASP	2.4
1	A	252	TYR	2.4
1	B	206	PRO	2.4
1	A	180	GLY	2.3
1	B	285	SER	2.3
1	A	326	SER	2.3
1	B	288	GLY	2.3
1	B	38	ASN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	129	LEU	2.2
1	A	122	THR	2.2
1	B	268	ASP	2.2
1	A	6	VAL	2.2
1	A	274	TYR	2.2
1	A	163	ILE	2.1
1	A	82	VAL	2.1
1	B	85	GLU	2.1
1	B	82	VAL	2.1
1	A	97	PHE	2.1
1	A	124	ALA	2.1
1	B	219	VAL	2.1
1	A	78	PHE	2.1
1	A	311	CYS	2.1
1	B	161	TYR	2.1
1	A	14	ALA	2.1
1	A	228	LEU	2.1
1	A	120	ILE	2.0
1	B	68	MET	2.0
1	B	125	ASN	2.0
1	B	29	PRO	2.0
1	A	315	LEU	2.0
1	B	267	ILE	2.0
1	B	252	TYR	2.0
1	B	154	ILE	2.0
1	A	21	TYR	2.0
1	A	32	TYR	2.0
1	A	160	LEU	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 monosaccharides in this entry.

6.4 Ligands (i)

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

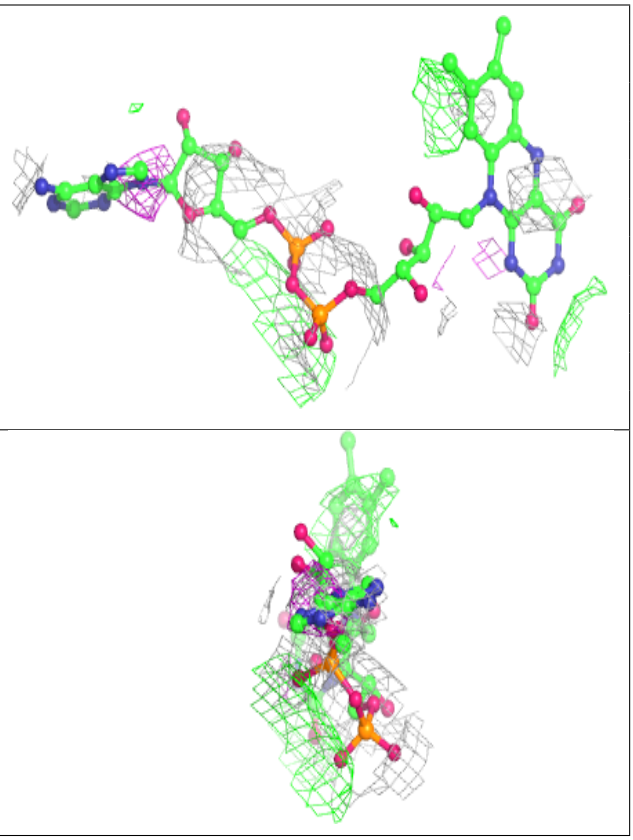
median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

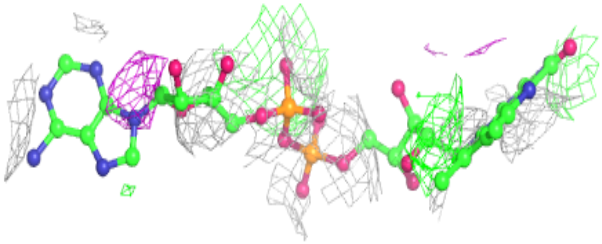
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	FAD	A	402	53/53	0.75	0.96	103,166,203,207	53
2	NAP	B	401	31/48	0.77	0.34	137,174,233,284	0
2	NAP	A	401	31/48	0.82	0.26	142,185,216,230	0
3	FAD	B	402	53/53	0.87	0.78	98,161,191,197	53

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 402:

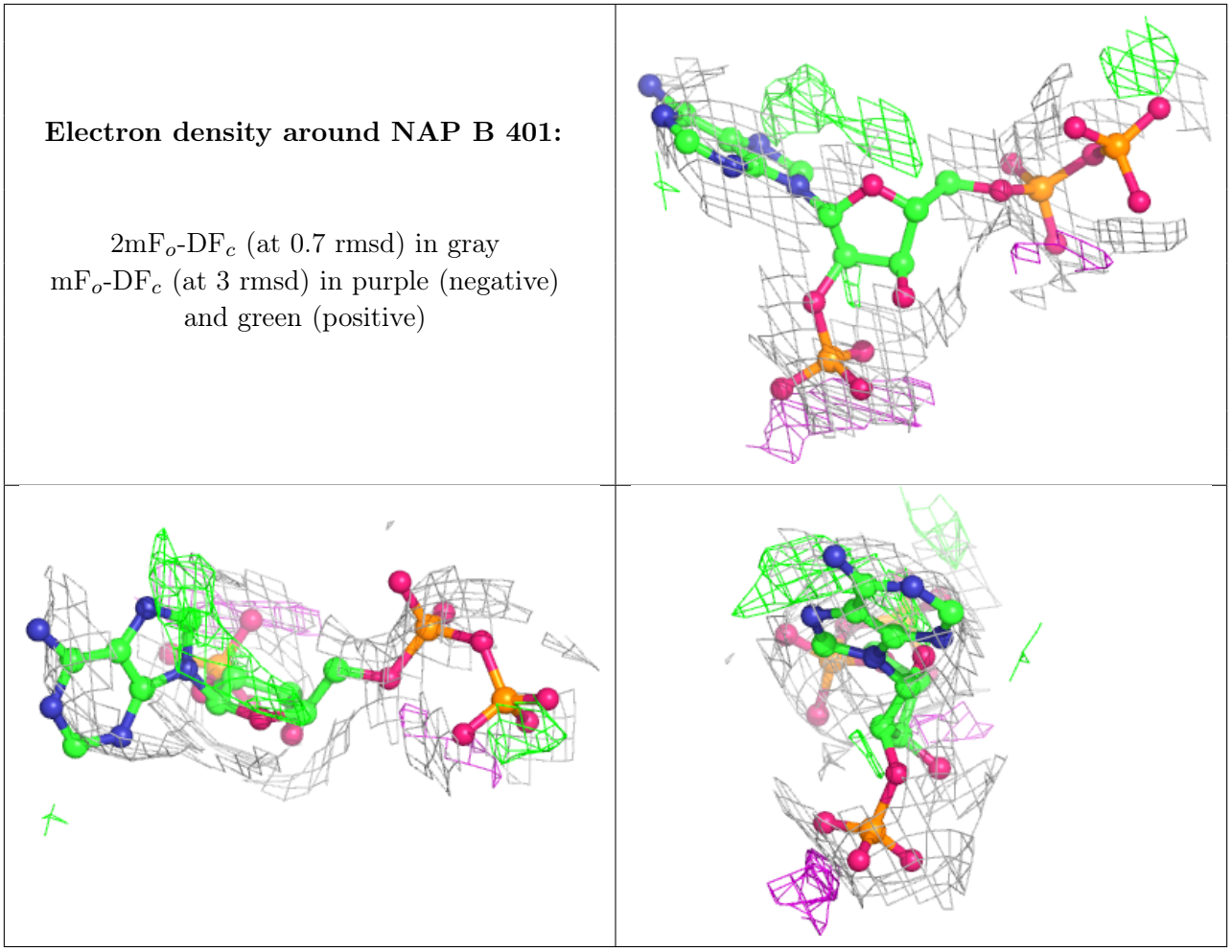
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





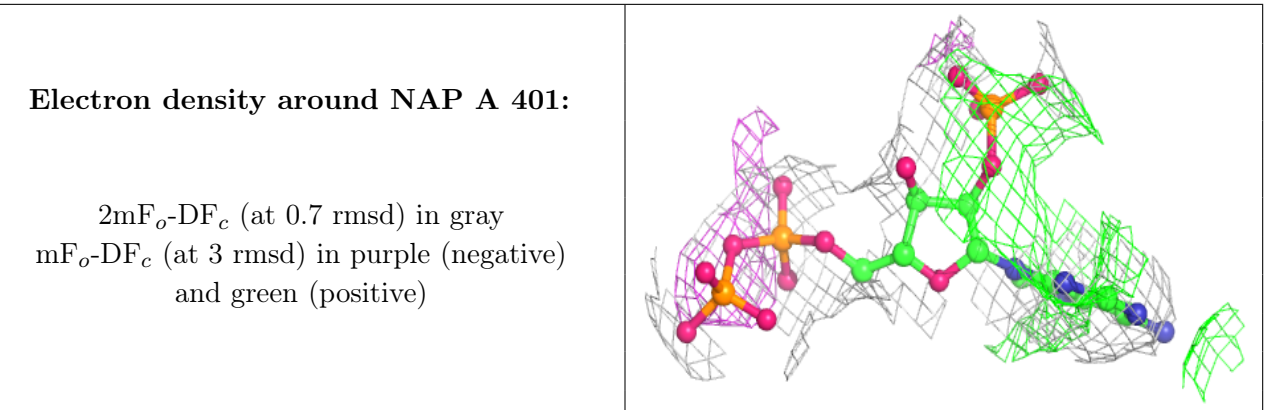
Electron density around NAP B 401:

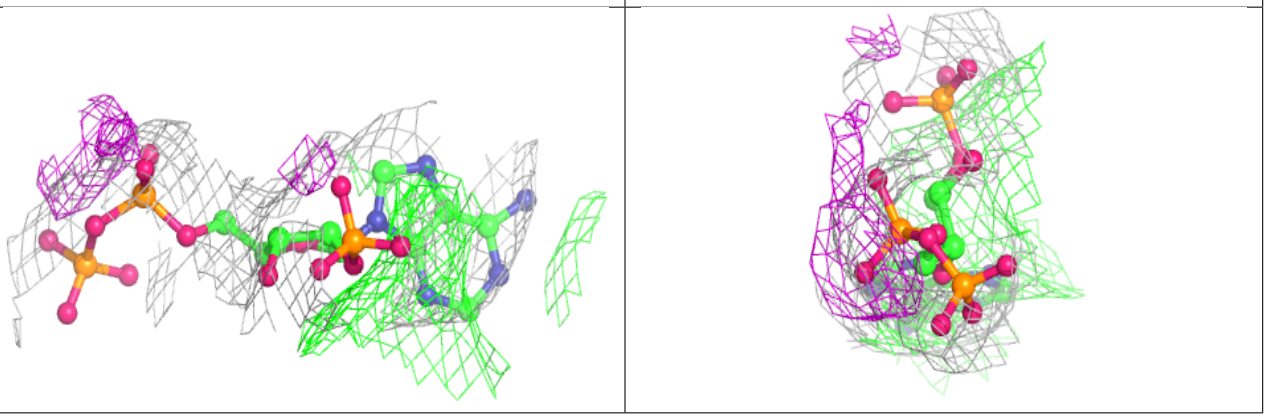
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around NAP A 401:

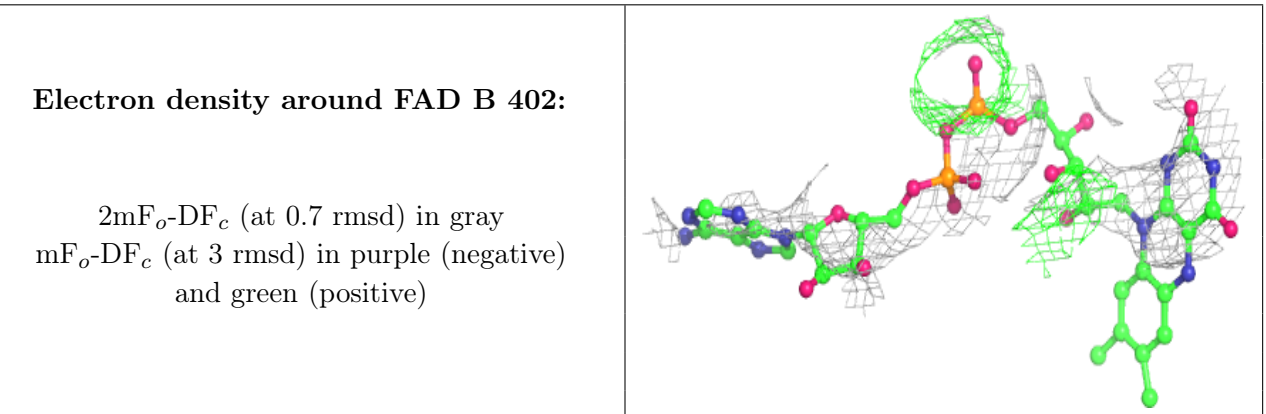
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

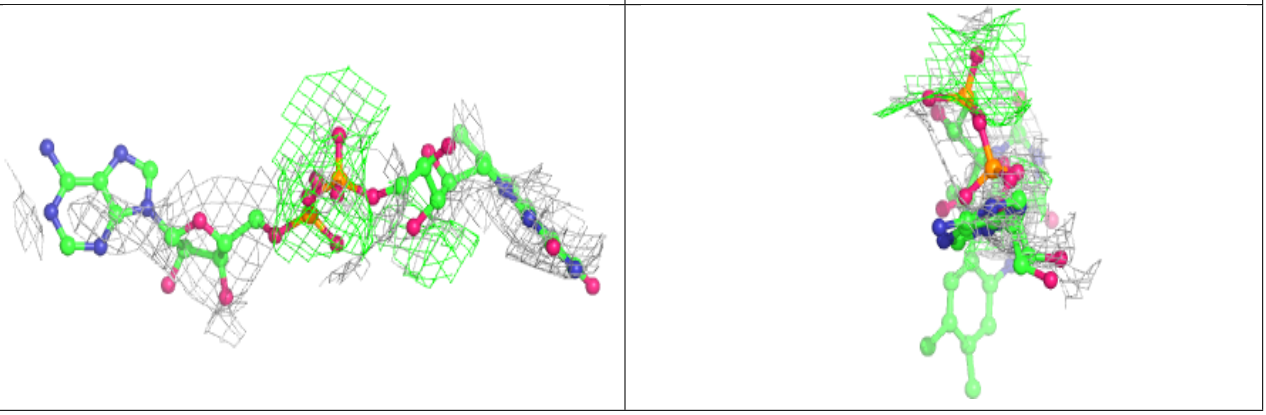




Electron density around FAD B 402:

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





6.5 Other polymers (i)

There are no such residues in this entry.