



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:43 PM UTC

PDB ID : 7V8O / pdb_00007v8o
Title : Crystal structure of cyclohexanone monooxygenase from T. municipale mutant L437T complexed with NADP⁺ and FAD in space group of P21221
Authors : Li, T.; Li, G.Y.; Yin, H.
Deposited on : 2021-08-23
Resolution : 2.72 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

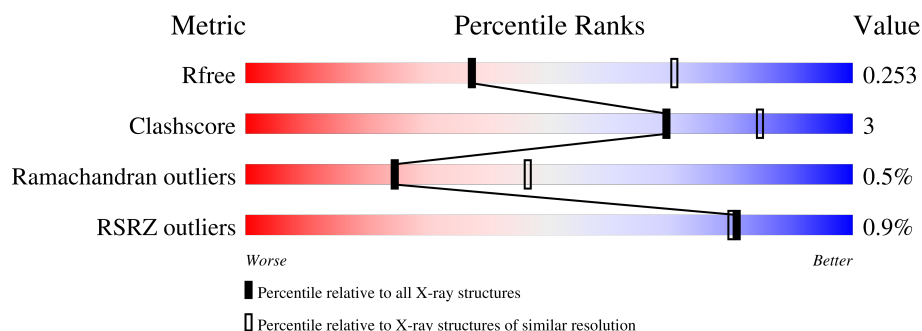
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	541	86% 11% .
1	B	541	90% 7% .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 8576 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.

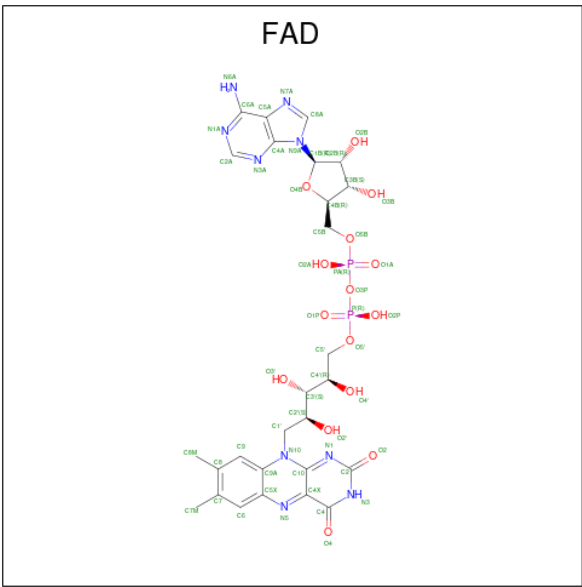
- Molecule 1 is a protein called Cyclohexanone Monooxygenase from Thermocrispum municipale.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	529	Total	C	N	O	S	0	4	0
			4184	2642	721	807	14			
1	B	529	Total	C	N	O	S	0	4	0
			4179	2639	718	807	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	437	THR	LEU	engineered mutation	UNP A0A1L1QK40
B	437	THR	LEU	engineered mutation	UNP A0A1L1QK40

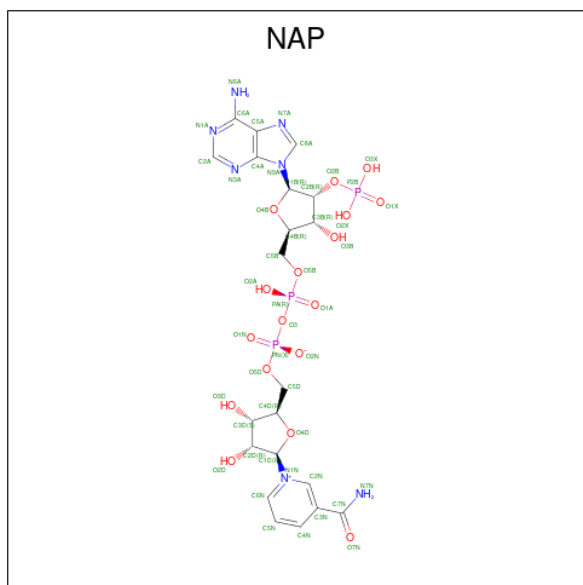
- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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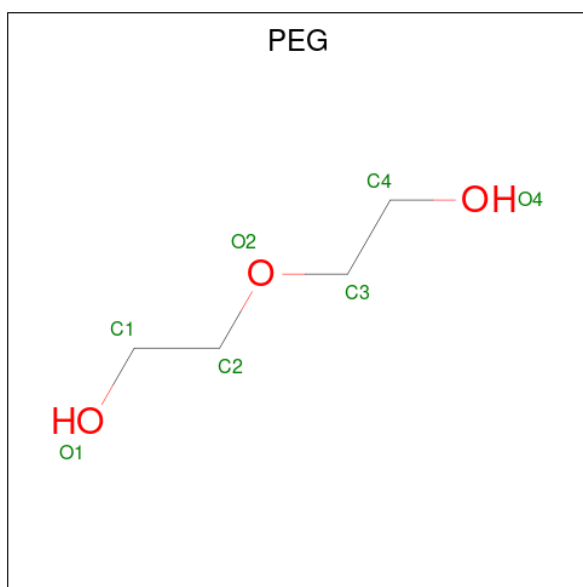
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $\text{C}_{21}\text{H}_{28}\text{N}_7\text{O}_{17}\text{P}_3$).



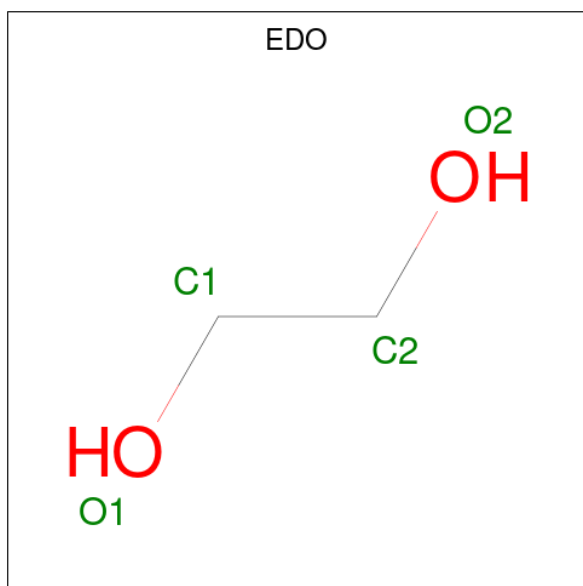
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 40	C 15	N 6	O 16	P 3	0	0
3	B	1	Total 40	C 15	N 6	O 16	P 3	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Na 1	0	0

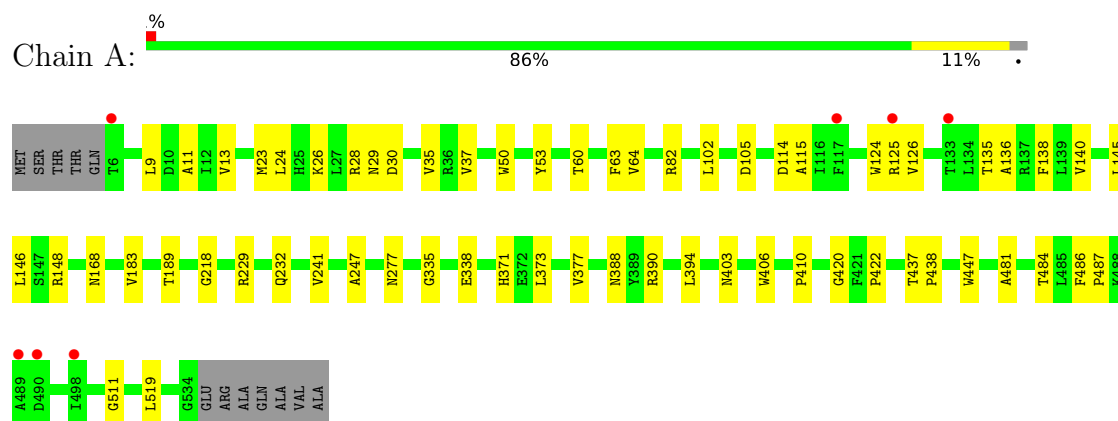
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	6	Total 6	O 6	0	0
7	B	9	Total 9	O 9	0	0

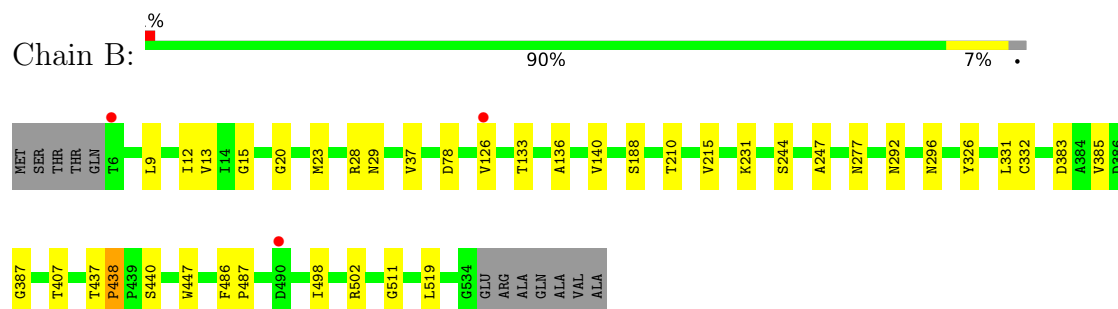
3 Residue-property plots

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

- Molecule 1: Cyclohexanone Monooxygenase from *Thermocrispum municipale*



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants a, b, c, α , β , γ	68.30Å 112.52Å 156.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.15 – 2.72 28.15 – 2.72	Depositor EDS
% Data completeness (in resolution range)	98.3 (28.15-2.72) 98.3 (28.15-2.72)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.31 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.220 , 0.281 (Not available) , 0.253	Depositor DCC
R_{free} test set	1656 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.463	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 12.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8576	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 75.01 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3714e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: FAD, PEG, NAP, EDO, NA

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	1.01	0/4283	1.43	1/5820 (0.0%)
1	B	1.01	0/4278	1.44	4/5814 (0.1%)
All	All	1.01	0/8561	1.44	5/11634 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	511	GLY	CA-C-O	-5.74	118.27	122.23
1	B	511	GLY	CA-C-O	-5.54	118.41	122.23
1	B	438	PRO	N-CA-C	5.22	117.07	110.70
1	B	440	SER	CA-C-N	5.02	127.07	120.60
1	B	440	SER	C-N-CA	5.02	127.07	120.60

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	4184	0	4025	38	0
1	B	4179	0	4017	19	0
2	A	53	0	31	8	0
2	B	53	0	31	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	40	0	19	0	0
3	B	40	0	19	1	0
4	B	7	0	10	0	0
5	B	4	0	6	0	0
6	B	1	0	0	0	0
7	A	6	0	0	0	0
7	B	9	0	0	0	0
All	All	8576	0	8158	58	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 3.

All (58) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:THR:CB	2:A:601:FAD:C7M	2.73	0.67
1:A:60:THR:HB	2:A:601:FAD:C7M	2.27	0.65
1:A:60:THR:HB	2:A:601:FAD:HM73	1.80	0.64
1:B:28:ARG:HH21	1:B:29:ASN:HD21	1.46	0.63
1:A:390:ARG:NH1	1:A:403:ASN:OD1	2.35	0.59
1:A:60:THR:OG1	2:A:601:FAD:C7M	2.51	0.58
1:A:28:ARG:HH21	1:A:29:ASN:HD21	1.51	0.57
1:B:292:ASN:OD1	1:B:292:ASN:C	2.49	0.56
1:A:11:ALA:HA	1:A:138:PHE:O	2.07	0.54
1:B:12:ILE:HG21	1:B:126:VAL:HG21	1.89	0.54
1:B:296:ASN:HD21	1:B:326:TYR:H	1.56	0.54
1:A:9:LEU:O	1:A:136:ALA:HA	2.08	0.54
1:A:115:ALA:HB3	1:A:394:LEU:HD13	1.91	0.53
1:A:124:TRP:O	1:A:135:THR:HA	2.10	0.52
1:B:9:LEU:O	1:B:136:ALA:HA	2.09	0.52
1:B:78[B]:ASP:OD1	1:B:231:LYS:NZ	2.41	0.51
1:A:125[A]:ARG:CZ	1:A:135:THR:HG23	2.40	0.51
1:A:437:THR:H	2:A:601:FAD:HM82	1.75	0.51
1:A:447:TRP:CD1	1:A:519:LEU:HB3	2.46	0.51
1:B:407:THR:O	1:B:502:ARG:NE	2.43	0.50
2:A:601:FAD:C9	2:A:601:FAD:O3'	2.60	0.50
1:B:447:TRP:CD1	1:B:519:LEU:HB3	2.47	0.49
1:A:63:PHE:CD2	1:A:64:VAL:HG13	2.48	0.49
1:B:188:SER:OG	1:B:331:LEU:HD22	2.12	0.49
1:A:241:VAL:HG13	1:A:247:ALA:HB1	1.95	0.48
1:A:53:TYR:CZ	1:A:189:THR:HG23	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ARG:O	1:A:232:GLN:HB2	2.15	0.47
1:A:406:TRP:CG	1:A:410:PRO:HA	2.50	0.47
1:A:11:ALA:O	1:A:35:VAL:HA	2.15	0.47
1:A:60:THR:OG1	2:A:601:FAD:HM72	2.15	0.46
1:A:13:VAL:O	1:A:37:VAL:HA	2.16	0.46
1:B:15:GLY:O	1:B:20:GLY:HA3	2.16	0.46
1:B:210:THR:OG1	3:B:602:NAP:O2X	2.32	0.45
1:A:183:VAL:HA	1:A:377:VAL:O	2.17	0.45
1:A:148:ARG:NH1	1:A:388:ASN:OD1	2.42	0.44
1:A:24:LEU:HG	1:A:102:LEU:HD22	1.99	0.44
1:A:23:MET:HG2	1:A:140:VAL:HG11	1.99	0.44
1:B:126:VAL:O	1:B:133:THR:HA	2.18	0.44
1:A:486:PHE:N	1:A:487:PRO:CD	2.82	0.43
1:B:486:PHE:N	1:B:487:PRO:CD	2.81	0.43
1:B:23:MET:HG2	1:B:140:VAL:HG11	2.01	0.43
1:A:60:THR:H	2:A:601:FAD:HM73	1.84	0.43
1:B:215:VAL:O	1:B:332:CYS:HA	2.19	0.42
1:A:145:LEU:C	1:A:146:LEU:HG	2.45	0.42
1:B:244:SER:OG	1:B:247:ALA:HA	2.20	0.42
1:A:28:ARG:NH1	1:A:105:ASP:OD2	2.51	0.42
1:A:481:ALA:O	1:A:484:THR:N	2.46	0.42
1:B:13:VAL:O	1:B:37:VAL:HA	2.20	0.42
1:A:82:ARG:NH1	1:A:218:GLY:O	2.53	0.41
1:A:114:ASP:O	1:A:126:VAL:HA	2.19	0.41
1:B:437:THR:N	1:B:438:PRO:CD	2.83	0.41
1:A:371:HIS:O	1:A:373:LEU:HG	2.20	0.41
1:A:420:GLY:C	1:A:422:PRO:HD3	2.45	0.41
1:A:335:GLY:HA2	1:A:338:GLU:OE1	2.19	0.41
1:A:26:LYS:O	1:A:30:ASP:HB2	2.22	0.40
1:A:437:THR:N	1:A:438:PRO:CD	2.85	0.40
1:B:383:ASP:HA	1:B:498:ILE:CD1	2.51	0.40
1:A:50:TRP:HB3	1:A:168:ASN:O	2.22	0.40

There are no symmetry-related clashes.

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	427/541 (79%)	397 (93%)	29 (7%)	1 (0%)	43	66
1	B	419/541 (77%)	391 (93%)	25 (6%)	3 (1%)	18	39
All	All	846/1082 (78%)	788 (93%)	54 (6%)	4 (0%)	24	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	ASN
1	B	277	ASN
1	B	387	GLY
1	B	385	VAL

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	A	601	-	58,58,58	0.69	1 (1%)	85,89,89	0.70	2 (2%)
5	EDO	B	604	-	3,3,3	0.19	0	2,2,2	0.23	0
4	PEG	B	603	-	6,6,6	0.23	0	5,5,5	0.10	0
3	NAP	B	602	-	42,43,52	0.44	0	62,67,80	0.84	2 (3%)
2	FAD	B	601	-	58,58,58	0.59	0	85,89,89	0.65	1 (1%)
3	NAP	A	602	-	42,43,52	0.63	1 (2%)	62,67,80	0.79	2 (3%)

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	FAD	A	601	-	-	4/34/50/50	0/6/6/6
5	EDO	B	604	-	-	1/1/1/1	-
4	PEG	B	603	-	-	3/4/4/4	-
3	NAP	B	602	-	-	1/27/59/67	0/4/4/5
2	FAD	B	601	-	-	6/34/50/50	0/6/6/6
3	NAP	A	602	-	-	1/27/59/67	0/4/4/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	C9-C8	-2.45	1.36	1.39
3	A	602	NAP	PA-O3	2.29	1.62	1.59

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	NAP	P2B-O2B-C2B	-4.65	111.01	123.43
3	A	602	NAP	P2B-O2B-C2B	-4.34	111.85	123.43
2	A	601	FAD	C1'-C2'-C3'	2.38	116.11	109.66
3	B	602	NAP	C3D-C2D-C1D	2.36	105.92	101.46
2	A	601	FAD	C4-N3-C2	-2.19	121.76	125.64
3	A	602	NAP	C3D-C2D-C1D	2.16	105.55	101.46
2	B	601	FAD	C4-N3-C2	-2.07	121.96	125.64

There are no chirality outliers.

All (16) torsion outliers are listed below:

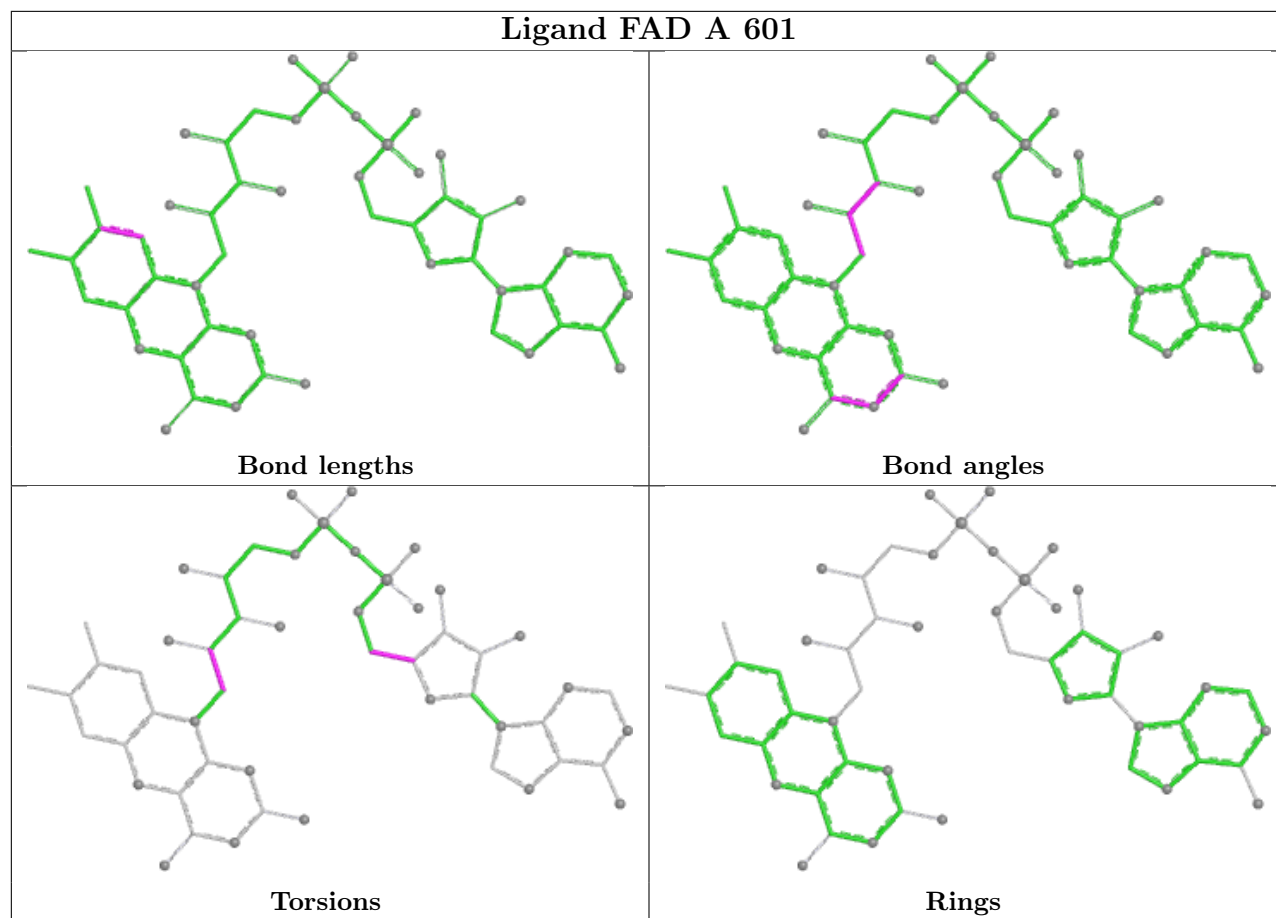
Mol	Chain	Res	Type	Atoms
2	A	601	FAD	N10-C1'-C2'-O2'
2	A	601	FAD	N10-C1'-C2'-C3'
2	B	601	FAD	N10-C1'-C2'-O2'
2	B	601	FAD	N10-C1'-C2'-C3'
3	B	602	NAP	PA-O3-PN-O5D
4	B	603	PEG	O1-C1-C2-O2
4	B	603	PEG	O2-C3-C4-O4
2	B	601	FAD	O4B-C4B-C5B-O5B
2	A	601	FAD	O4B-C4B-C5B-O5B
2	B	601	FAD	C3B-C4B-C5B-O5B
4	B	603	PEG	C1-C2-O2-C3
2	A	601	FAD	C3B-C4B-C5B-O5B
2	B	601	FAD	C2'-C3'-C4'-C5'
3	A	602	NAP	C2B-O2B-P2B-O1X
5	B	604	EDO	O1-C1-C2-O2
2	B	601	FAD	O3'-C3'-C4'-O4'

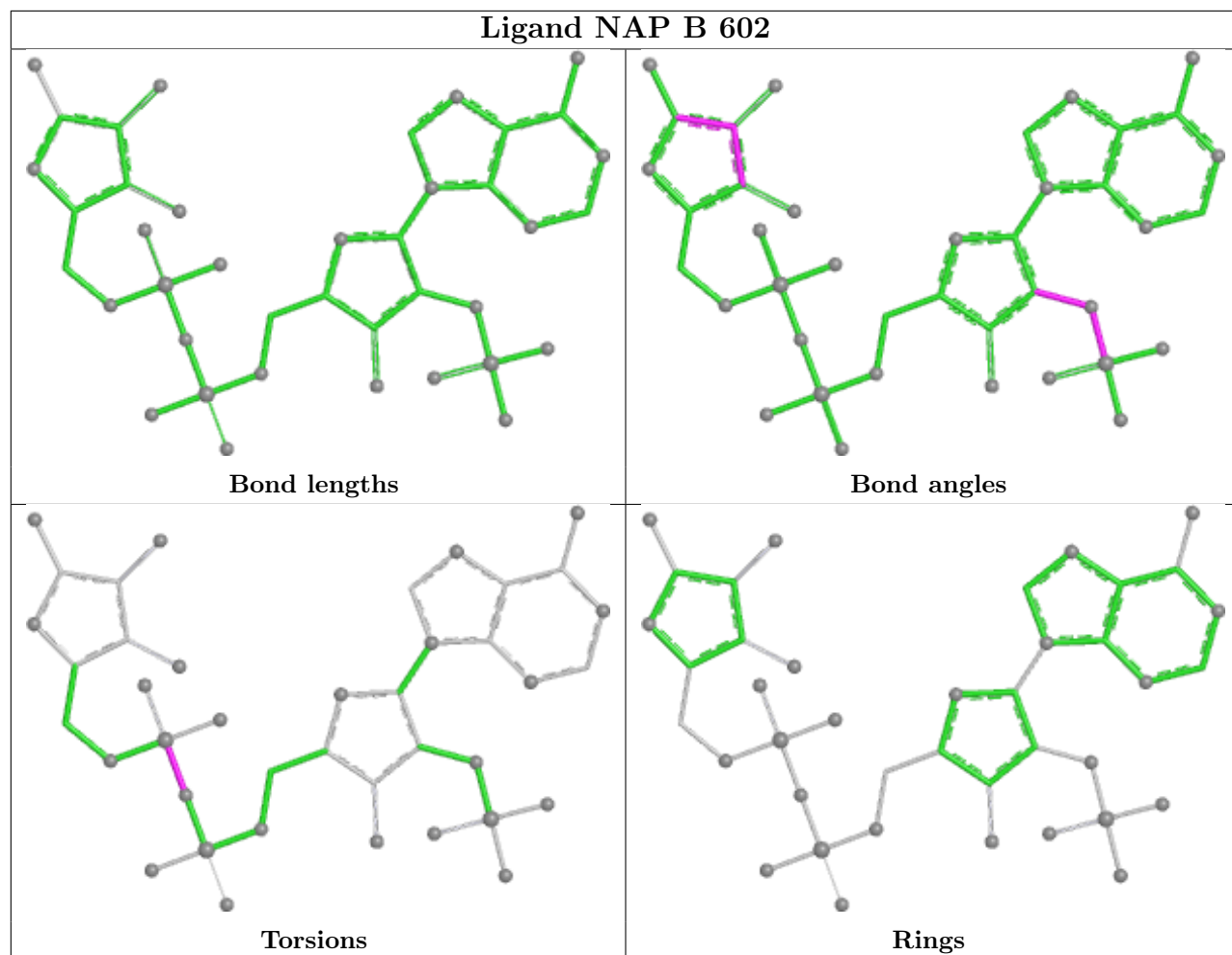
There are no ring outliers.

2 monomers are involved in 9 short contacts:

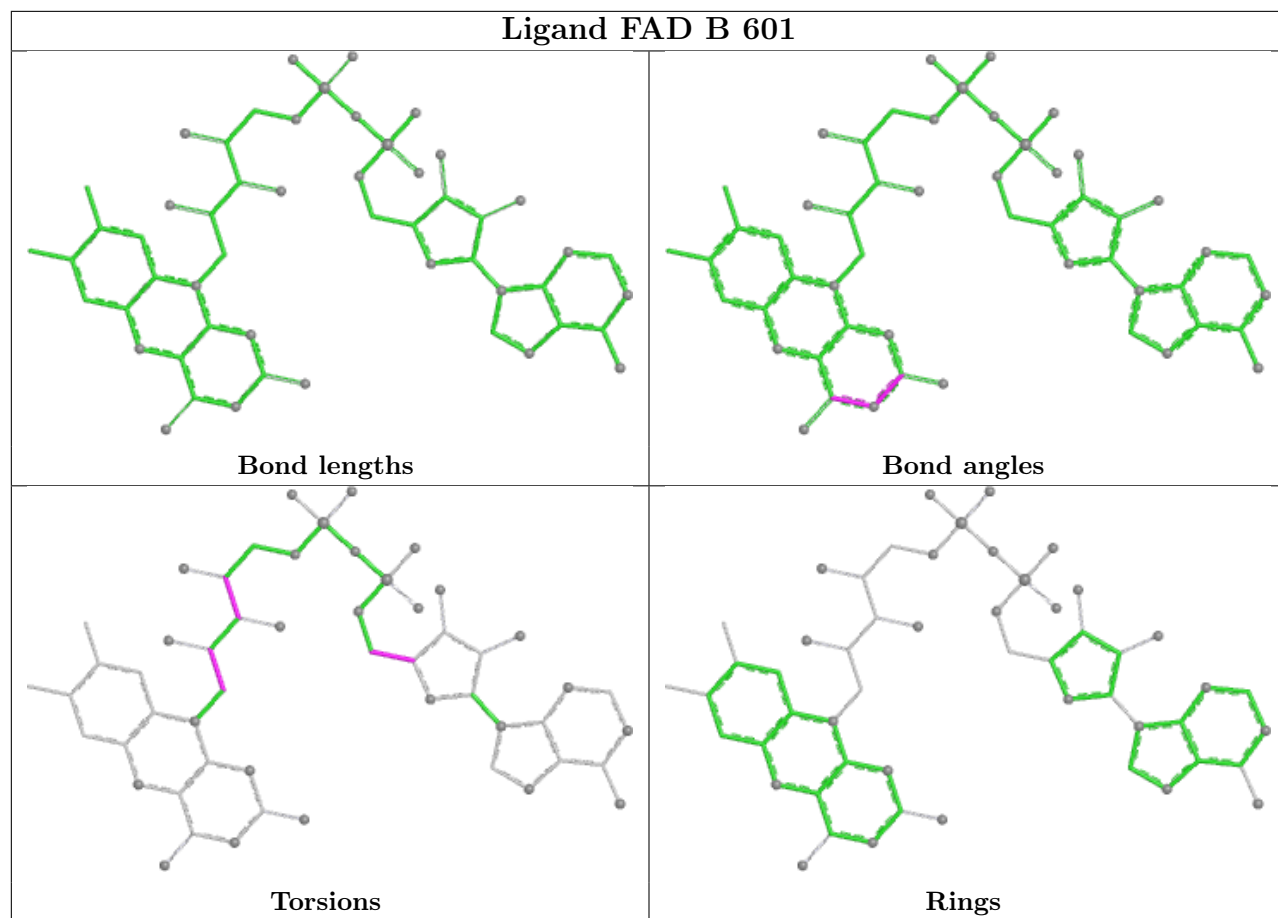
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	8	0
3	B	602	NAP	1	0

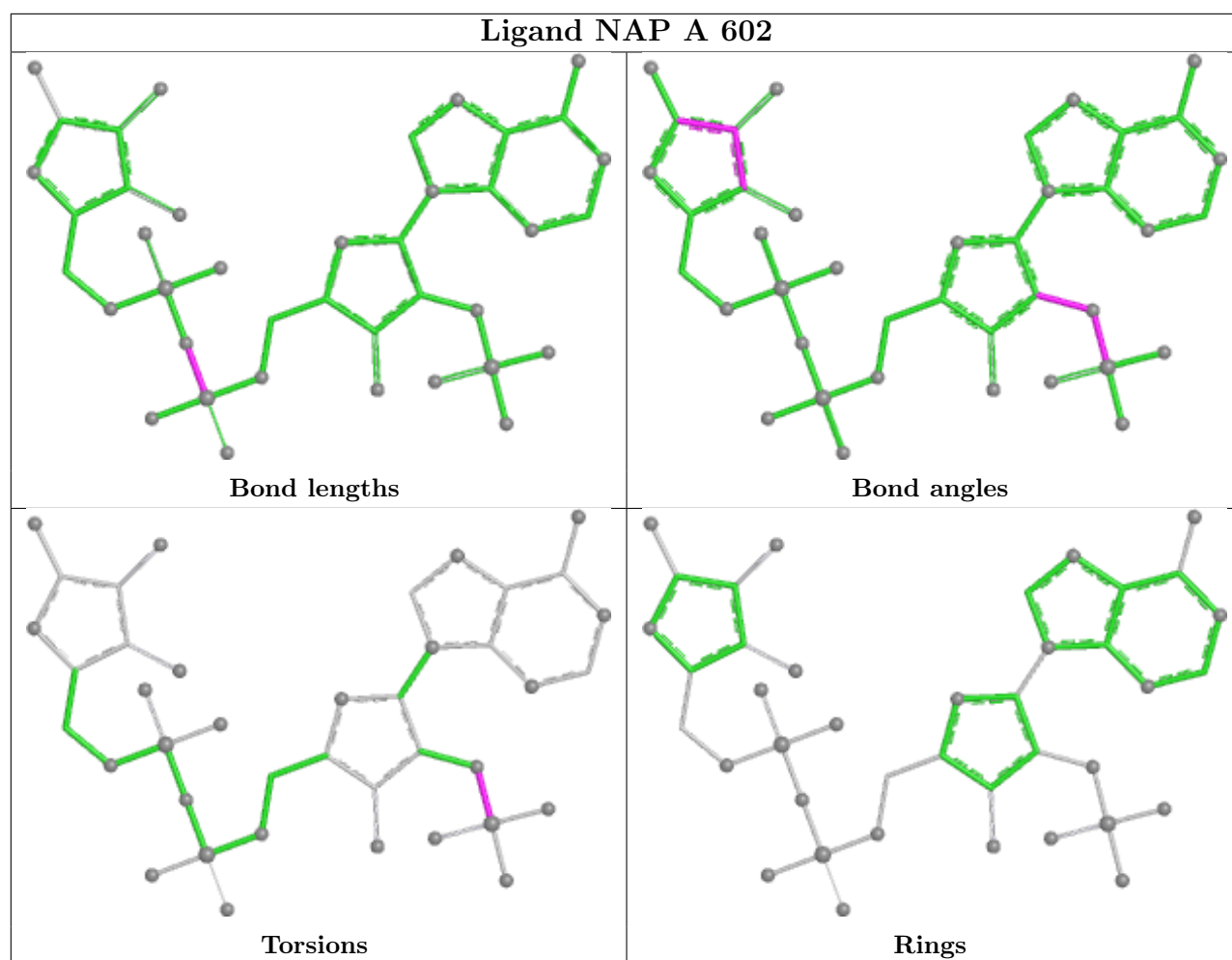
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 B 601





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 [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	529/541 (97%)	0.13	7 (1%) 75 73	17, 36, 54, 65	4 (0%)
1	B	529/541 (97%)	-0.09	3 (0%) 85 85	13, 31, 48, 62	4 (0%)
All	All	1058/1082 (97%)	0.02	10 (0%) 81 80	13, 33, 51, 65	8 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	489	ALA	2.9
1	A	498	ILE	2.7
1	A	6	THR	2.6
1	B	126	VAL	2.5
1	A	125[A]	ARG	2.4
1	B	490	ASP	2.4
1	A	133	THR	2.3
1	A	117	PHE	2.3
1	B	6	THR	2.2
1	A	490	ASP	2.2

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 ⓘ

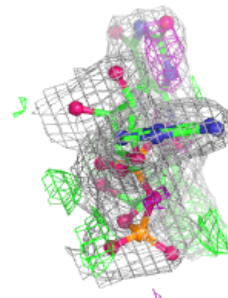
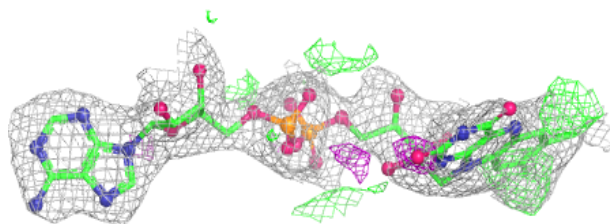
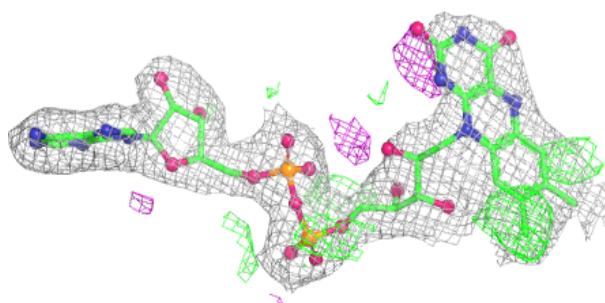
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	EDO	B	604	4/4	0.79	0.14	31,34,34,35	0
4	PEG	B	603	7/7	0.88	0.10	37,40,42,44	0
2	FAD	A	601	53/53	0.89	0.12	26,31,41,47	0
6	NA	B	605	1/1	0.90	0.09	41,41,41,41	0
3	NAP	A	602	40/48	0.92	0.10	35,43,57,58	0
3	NAP	B	602	40/48	0.93	0.10	28,33,44,44	0
2	FAD	B	601	53/53	0.97	0.07	20,22,29,29	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.

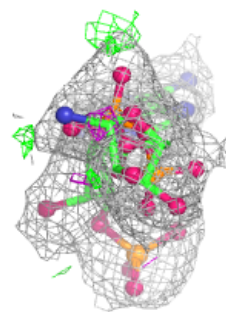
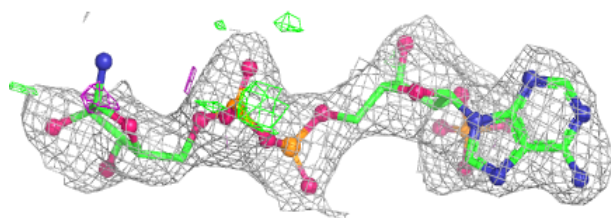
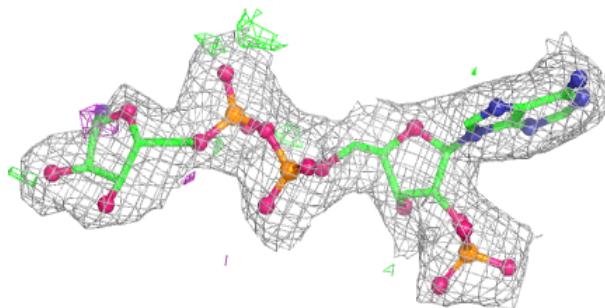
Electron density around FAD A 601:

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

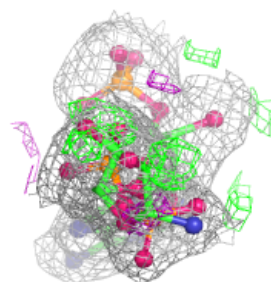
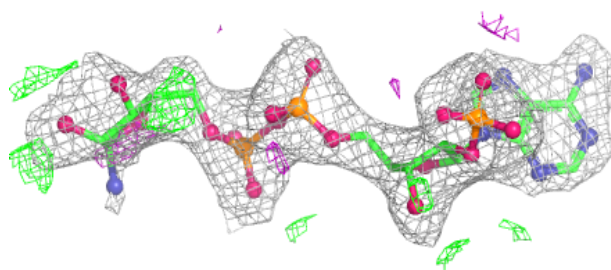
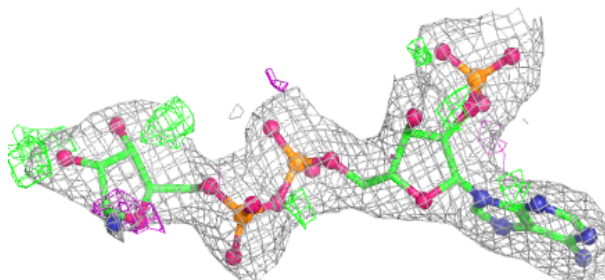


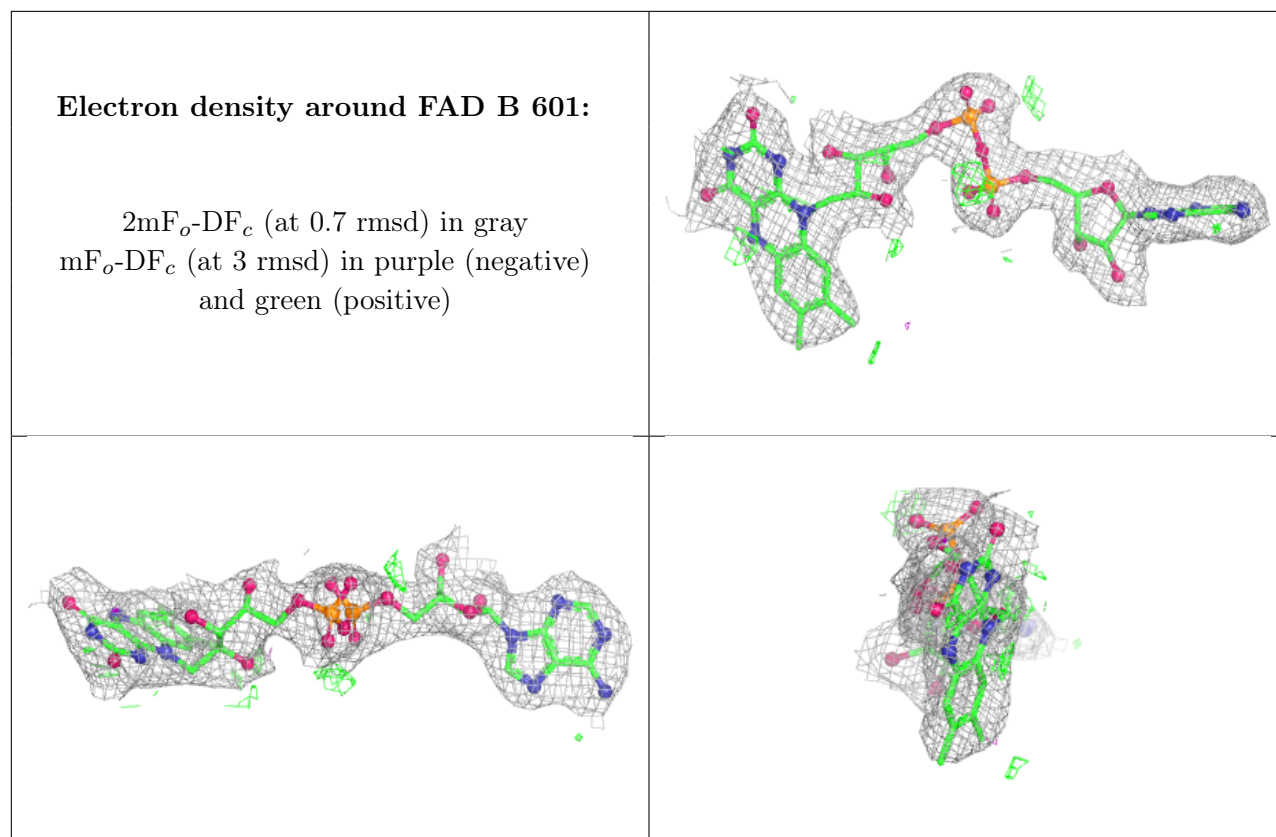
Electron density around NAP A 602:

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

**Electron density around NAP B 602:**

$2mF_o-DF_c$ (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.