



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 09:31 PM EDT

PDB ID : 6WO1
Title : Hybrid acetohydroxyacid synthase complex structure with *Cryptococcus neoformans* AHAS catalytic subunit and *Saccharomyces cerevisiae* AHAS regulatory subunit
Authors : Guddat, L.W.; Lonhienne, T.
Deposited on : 2020-04-23
Resolution : 3.30 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 : 2.35.1
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) : 2.35.1

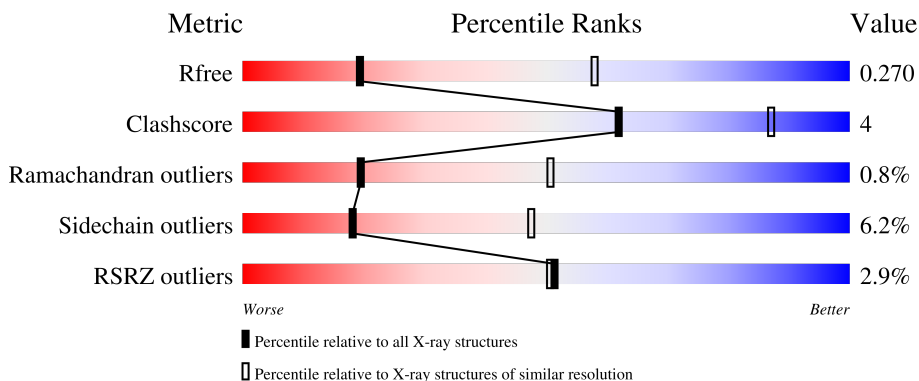
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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	718	
2	B	309	

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
4	8GF	A	1702	-	X	-	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5678 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.

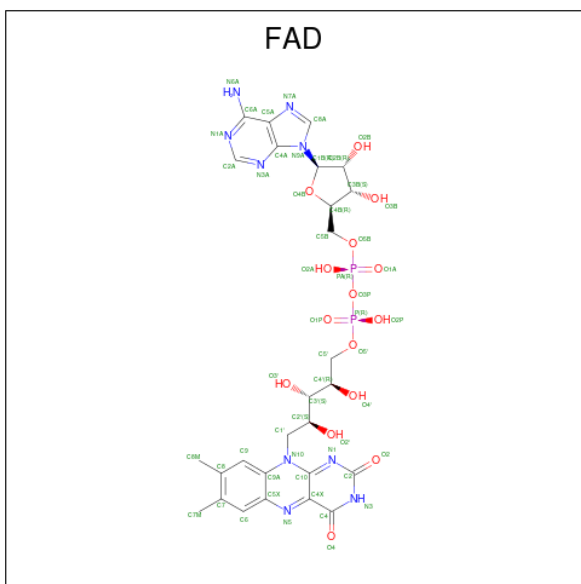
- Molecule 1 is a protein called Acetohydroxyacid synthase catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	551	4093	2601	696	773	23	0	0	0

- Molecule 2 is a protein called Acetohydroxyacid synthase regulatory subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	197	1506	945	270	283	8	0	0	0

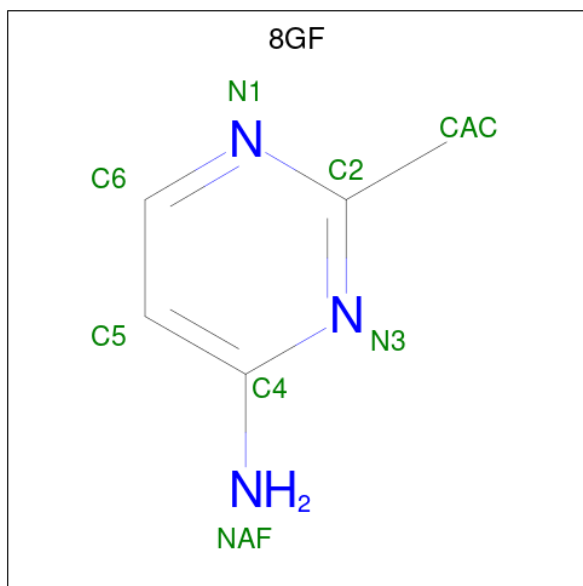
- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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

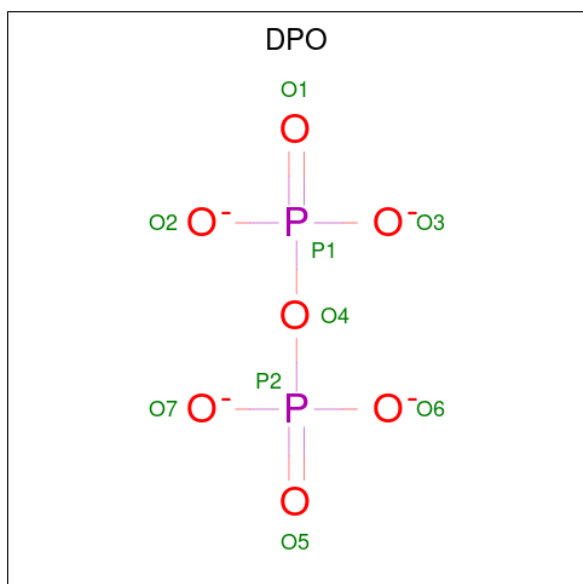
- Molecule 4 is 2-methylpyrimidin-4-amine (three-letter code: 8GF) (formula: $C_5H_7N_3$)

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



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	C N	0	0
			8	5 3		

- Molecule 5 is DIPHOSPHATE (three-letter code: DPO) (formula: O₇P₂) (labeled as "Ligand of Interest" by depositor).

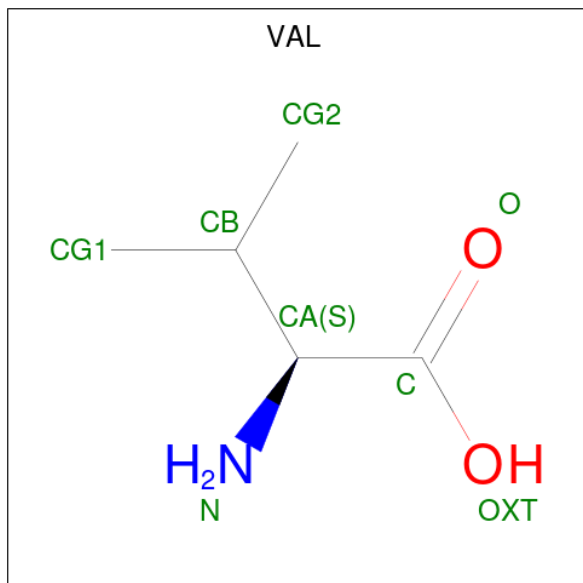


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O P	0	0
			9	7 2		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Mg 1 1	0	0

- Molecule 7 is VALINE (three-letter code: VAL) (formula: C₅H₁₁NO₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total C N O 8 5 1 2	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 4 3 2	Depositor
Cell constants a, b, c, α , β , γ	195.68Å 195.68Å 195.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.92 – 3.30 48.92 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.92-3.30) 99.5 (48.92-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 3.33Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.210 , 0.269 0.210 , 0.270	Depositor DCC
R_{free} test set	1978 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	92.7	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5678	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, DPO, MG, 8GF

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	0.27	0/4182	0.44	0/5688
2	B	0.27	0/1528	0.50	1/2075 (0.0%)
All	All	0.27	0/5710	0.46	1/7763 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	282	PRO	N-CA-CB	5.93	110.42	103.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4093	0	3987	36	0
2	B	1506	0	1488	12	0
3	A	53	0	30	4	0
4	A	8	0	0	1	0
5	A	9	0	0	0	0
6	A	1	0	0	0	0
7	B	8	0	8	1	0
All	All	5678	0	5513	49	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 4.

All (49) 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:449:GLU:OE2	3:A:1701:FAD:O3B	1.91	0.89
1:A:567:MET:H	4:A:1702:8GF:CAC	1.96	0.78
1:A:468:ASP:OD2	1:A:470:VAL:HG22	1.92	0.69
1:A:386:ARG:NH1	1:A:526:GLU:OE1	2.29	0.66
2:B:151:TYR:HB3	2:B:156:ILE:HG21	1.78	0.65
1:A:444:GLY:HA3	1:A:462:GLN:HG3	1.79	0.62
1:A:348:ALA:HB3	1:A:375:THR:HG22	1.80	0.62
2:B:129:ASP:N	2:B:129:ASP:OD1	2.33	0.61
1:A:185:ARG:NH2	1:A:350:ASN:OD1	2.37	0.58
2:B:85:GLN:HG2	2:B:87:GLU:HG3	1.87	0.57
1:A:457:LYS:HG2	1:A:458:ILE:HG13	1.88	0.55
1:A:187:SER:O	1:A:189:LYS:N	2.38	0.55
1:A:377:LEU:O	1:A:558:TRP:NE1	2.40	0.55
1:A:241:VAL:O	1:A:251:ASN:ND2	2.41	0.53
1:A:445:ILE:H	1:A:462:GLN:HG2	1.73	0.53
1:A:136:MET:HE3	1:A:141:VAL:HG21	1.92	0.52
1:A:387:ASP:OD2	1:A:389:LYS:N	2.36	0.50
2:B:166:ILE:HD13	2:B:262:VAL:HG21	1.94	0.50
2:B:156:ILE:HA	2:B:279:PRO:HA	1.94	0.48
1:A:195:VAL:O	1:A:222:CYS:HA	2.13	0.48
2:B:124:VAL:HG21	2:B:275:MET:HB3	1.97	0.47
1:A:384:ASP:OD1	1:A:386:ARG:HG2	2.15	0.47
1:A:387:ASP:OD2	1:A:389:LYS:HB2	2.15	0.47
1:A:477:VAL:HG22	1:A:478:PRO:HD3	1.97	0.47
2:B:198:GLU:HG2	2:B:202:LYS:HE3	1.97	0.46
3:A:1701:FAD:H9	3:A:1701:FAD:H1'1	1.77	0.46
1:A:395:GLY:HA3	3:A:1701:FAD:HN3	1.81	0.45
2:B:180:HIS:HA	2:B:183:THR:HB	1.99	0.45
1:A:537:THR:HG22	1:A:589:ILE:HB	1.98	0.45
1:A:206:THR:HB	1:A:207:PRO:HD3	1.98	0.45
1:A:218:MET:O	1:A:278:PRO:HA	2.17	0.45
1:A:346:ILE:HG21	1:A:362:LEU:HD11	1.98	0.45
2:B:101:ARG:HB3	2:B:101:ARG:HH21	1.82	0.45
1:A:710:HIS:H	1:A:711:PRO:CD	2.29	0.45
1:A:275:ARG:HE	3:A:1701:FAD:H3B	1.82	0.44
1:A:324:LEU:H	1:A:324:LEU:HD12	1.82	0.44
1:A:146:GLY:O	1:A:170:PRO:HD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:PRO:HG3	1:A:176:ALA:HB2	1.99	0.44
1:A:330:LEU:HD23	1:A:330:LEU:HA	1.88	0.43
2:B:110:VAL:HG21	7:B:401:VAL:HG21	2.01	0.43
1:A:493:LYS:HE3	1:A:493:LYS:HB2	1.83	0.43
1:A:132:PHE:O	1:A:136:MET:HG2	2.18	0.43
1:A:388:GLU:H	1:A:388:GLU:HG3	1.65	0.41
1:A:353:LEU:O	1:A:556:ARG:NH1	2.49	0.41
1:A:380:LEU:HG	1:A:552:TRP:CD1	2.55	0.41
2:B:226:THR:HA	2:B:229:THR:HG22	2.02	0.41
1:A:578:LYS:HA	1:A:586:VAL:HG21	2.02	0.41
1:A:697:MET:O	1:A:708:LEU:HB2	2.21	0.41
2:B:98:LEU:HD21	2:B:139:ILE:HD11	2.02	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	545/718 (76%)	518 (95%)	24 (4%)	3 (1%)	25	57
2	B	193/309 (62%)	183 (95%)	7 (4%)	3 (2%)	9	36
All	All	738/1027 (72%)	701 (95%)	31 (4%)	6 (1%)	19	51

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	710	HIS
2	B	75	ARG
2	B	182	HIS
1	A	188	GLY
1	A	506	SER
2	B	281	THR

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	420/593 (71%)	399 (95%)	21 (5%)	24	55
2	B	163/271 (60%)	148 (91%)	15 (9%)	9	31
All	All	583/864 (68%)	547 (94%)	36 (6%)	18	47

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	MET
1	A	144	VAL
1	A	167	PHE
1	A	208	MET
1	A	248	THR
1	A	324	LEU
1	A	350	ASN
1	A	388	GLU
1	A	401	TYR
1	A	407	GLN
1	A	417	VAL
1	A	447	HIS
1	A	449	GLU
1	A	460	GLU
1	A	476	LEU
1	A	477	VAL
1	A	590	ASP
1	A	692	GLU
1	A	705	HIS
1	A	708	LEU
1	A	710	HIS
2	B	84	VAL
2	B	85	GLN
2	B	91	LEU
2	B	101	ARG
2	B	105	ILE
2	B	107	SER

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Mol	Chain	Res	Type
2	B	125	LEU
2	B	126	GLN
2	B	128	GLN
2	B	129	ASP
2	B	136	ARG
2	B	165	ARG
2	B	171	THR
2	B	172	GLU
2	B	215	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	8GF	A	1702	-	8,8,8	6.05	3 (37%)	10,10,10	3.09	7 (70%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DPO	A	1703	6	6,8,8	0.75	0	13,13,13	1.13	1 (7%)
7	VAL	B	401	-	5,7,7	0.90	0	7,9,9	1.09	1 (14%)
3	FAD	A	1701	-	53,58,58	1.67	11 (20%)	68,89,89	1.35	13 (19%)

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
4	8GF	A	1702	-	-	-	0/1/1/1
5	DPO	A	1703	6	-	0/6/6/6	-
7	VAL	B	401	-	-	3/8/8/8	-
3	FAD	A	1701	-	-	8/30/50/50	0/6/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1702	8GF	CAC-C2	-14.22	1.09	1.49
4	A	1702	8GF	C4-NAF	-8.82	1.09	1.35
3	A	1701	FAD	C10-N1	5.07	1.43	1.33
3	A	1701	FAD	C2B-C1B	-4.66	1.46	1.53
3	A	1701	FAD	C4X-N5	4.55	1.39	1.30
3	A	1701	FAD	O3'-C3'	-2.80	1.36	1.43
3	A	1701	FAD	O4B-C1B	2.57	1.44	1.41
3	A	1701	FAD	C2-N1	2.56	1.42	1.36
3	A	1701	FAD	C1'-N10	-2.39	1.42	1.48
3	A	1701	FAD	C2'-C3'	2.37	1.57	1.53
4	A	1702	8GF	C5-C4	-2.31	1.35	1.41
3	A	1701	FAD	C2B-C3B	-2.12	1.47	1.53
3	A	1701	FAD	C1'-C2'	2.04	1.55	1.52
3	A	1701	FAD	C4X-C4	2.03	1.51	1.44

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1702	8GF	C6-N1-C2	5.25	120.11	115.74
4	A	1702	8GF	C6-C5-C4	4.85	119.72	116.91
3	A	1701	FAD	N3A-C2A-N1A	-3.97	122.47	128.68
4	A	1702	8GF	N1-C2-N3	-3.48	119.55	125.54
4	A	1702	8GF	C5-C6-N1	-2.89	120.37	123.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1701	FAD	C4-N3-C2	-2.88	120.32	125.64
5	A	1703	DPO	P2-O4-P1	-2.76	123.35	132.83
3	A	1701	FAD	C4X-C4-N3	2.75	120.17	113.19
4	A	1702	8GF	NAF-C4-N3	2.59	120.58	116.49
4	A	1702	8GF	CAC-C2-N1	2.50	119.89	117.14
3	A	1701	FAD	C4A-C5A-N7A	-2.47	106.83	109.40
3	A	1701	FAD	C9A-C5X-N5	-2.46	119.76	122.43
3	A	1701	FAD	O4-C4-C4X	-2.44	120.13	126.60
3	A	1701	FAD	C4X-C10-N1	-2.41	119.14	124.73
3	A	1701	FAD	C10-C4X-N5	-2.30	119.98	124.86
3	A	1701	FAD	C4-C4X-N5	2.27	121.46	118.23
3	A	1701	FAD	C3B-C2B-C1B	2.25	104.36	100.98
7	B	401	VAL	OXT-C-O	-2.21	119.07	124.09
4	A	1702	8GF	CAC-C2-N3	2.18	120.56	117.15
3	A	1701	FAD	C4'-C3'-C2'	-2.13	108.94	113.36
3	A	1701	FAD	C4X-C10-N10	2.12	119.58	116.48
3	A	1701	FAD	C10-N1-C2	2.06	121.03	116.90

There are no chirality outliers.

All (11) torsion outliers are listed below:

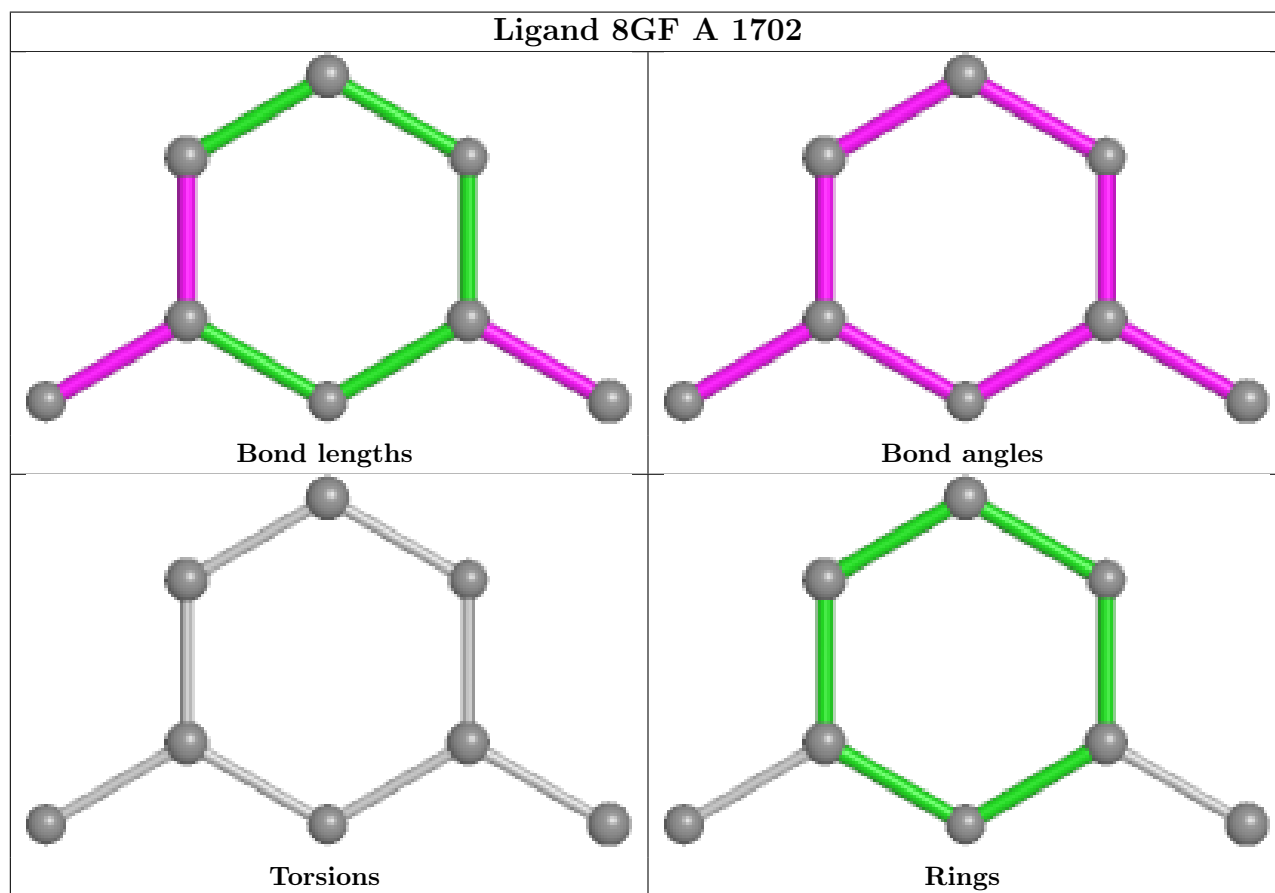
Mol	Chain	Res	Type	Atoms
3	A	1701	FAD	C3'-C4'-C5'-O5'
3	A	1701	FAD	O4'-C4'-C5'-O5'
3	A	1701	FAD	C5'-O5'-P-O1P
3	A	1701	FAD	C5'-O5'-P-O2P
3	A	1701	FAD	C5'-O5'-P-O3P
7	B	401	VAL	O-C-CA-CB
3	A	1701	FAD	P-O3P-PA-O5B
7	B	401	VAL	OXT-C-CA-CB
3	A	1701	FAD	O4B-C4B-C5B-O5B
3	A	1701	FAD	PA-O3P-P-O2P
7	B	401	VAL	C-CA-CB-CG2

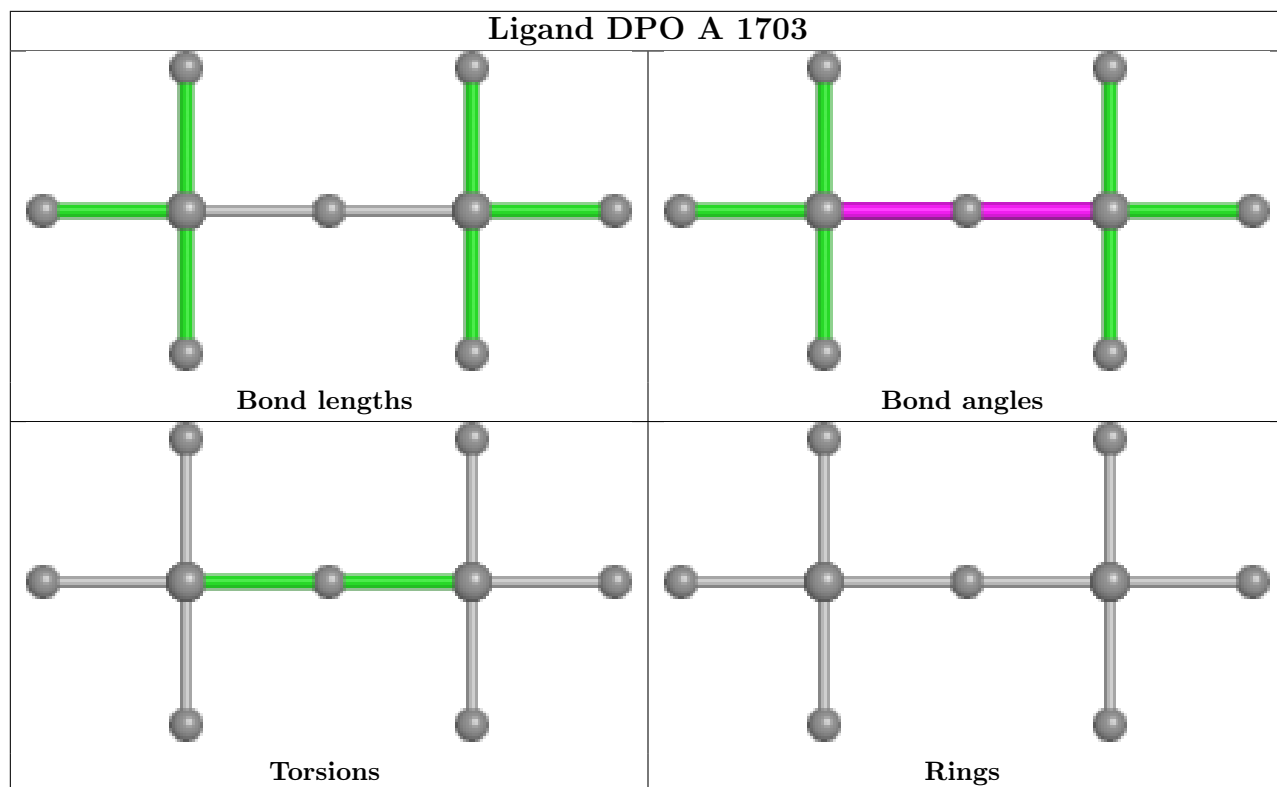
There are no ring outliers.

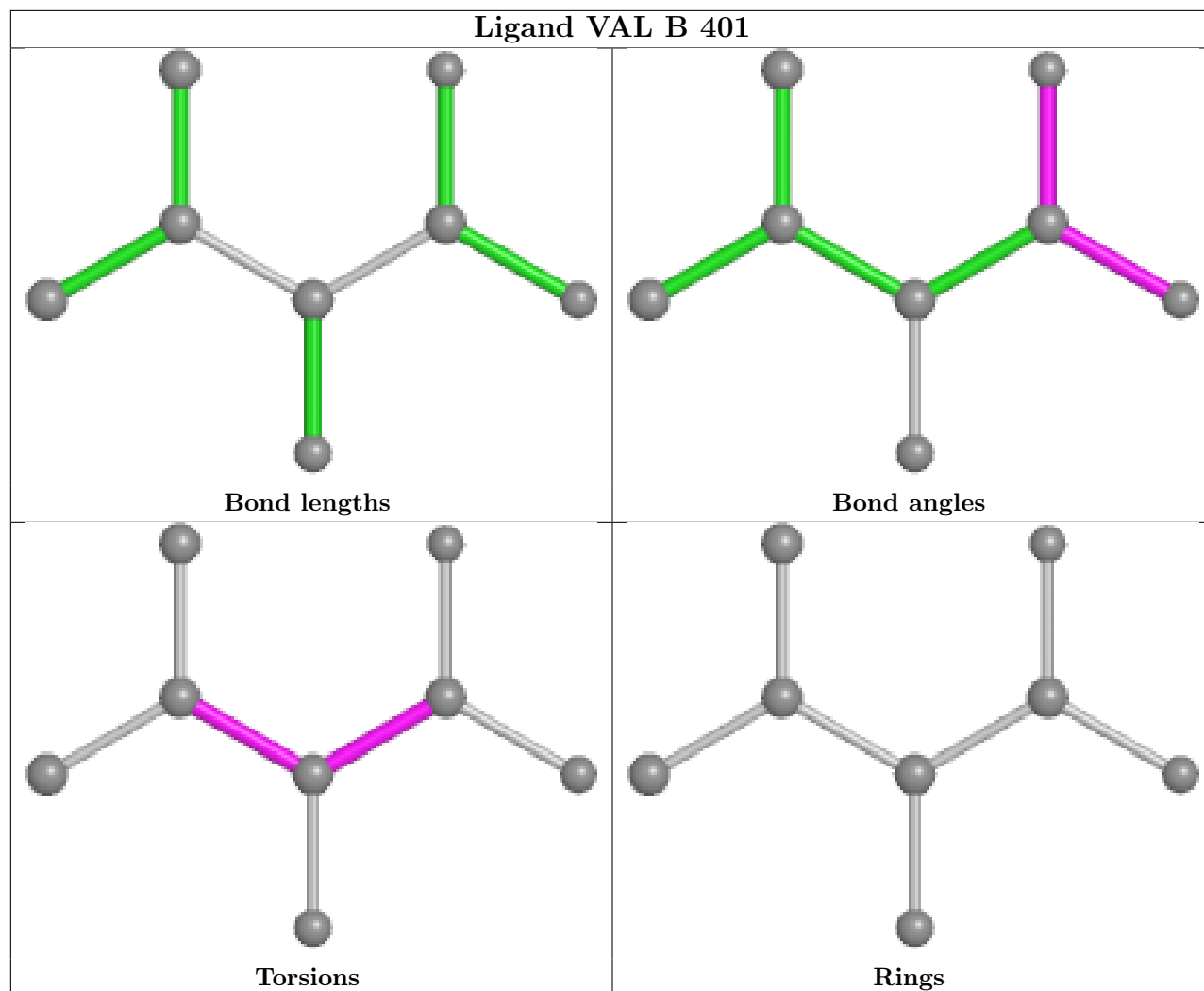
3 monomers are involved in 6 short contacts:

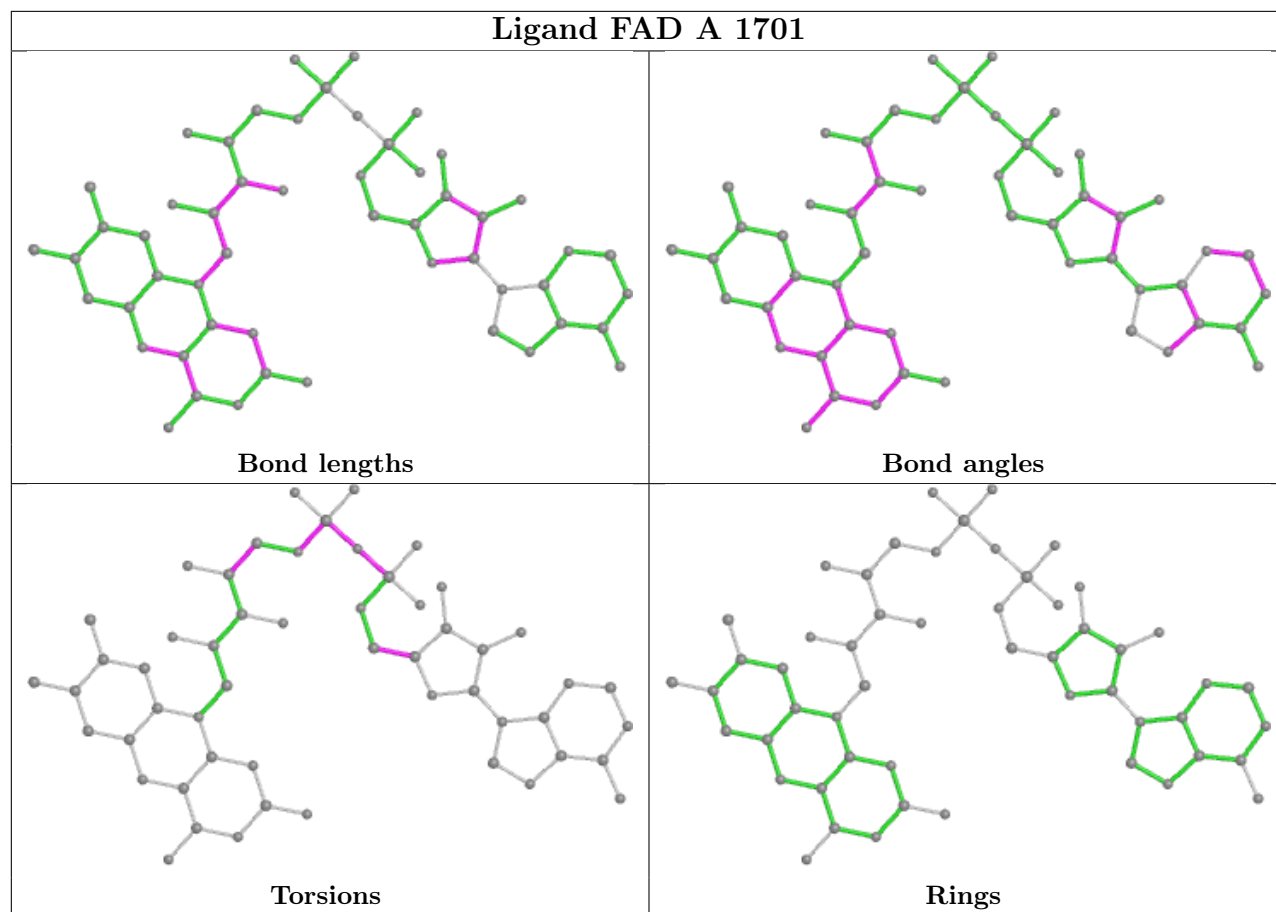
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1702	8GF	1	0
7	B	401	VAL	1	0
3	A	1701	FAD	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	551/718 (76%)	0.20	13 (2%) 59 56	53, 93, 129, 179	0
2	B	197/309 (63%)	0.37	9 (4%) 32 30	58, 85, 160, 180	0
All	All	748/1027 (72%)	0.24	22 (2%) 51 50	53, 92, 137, 180	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	199	ILE	4.4
2	B	203	GLN	3.3
2	B	198	GLU	3.1
1	A	121	TYR	3.1
2	B	201	GLU	2.9
1	A	705	HIS	2.9
1	A	203	ASN	2.7
1	A	292	LEU	2.5
2	B	183	THR	2.4
1	A	206	THR	2.4
1	A	120	ASP	2.4
1	A	296	ILE	2.4
2	B	200	ARG	2.3
2	B	204	PHE	2.3
1	A	297	PRO	2.3
1	A	512	LEU	2.2
1	A	709	LEU	2.2
1	A	568	GLY	2.2
2	B	153	ASN	2.1
1	A	210	ASP	2.1
1	A	499	TYR	2.1
2	B	227	ASN	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](#)

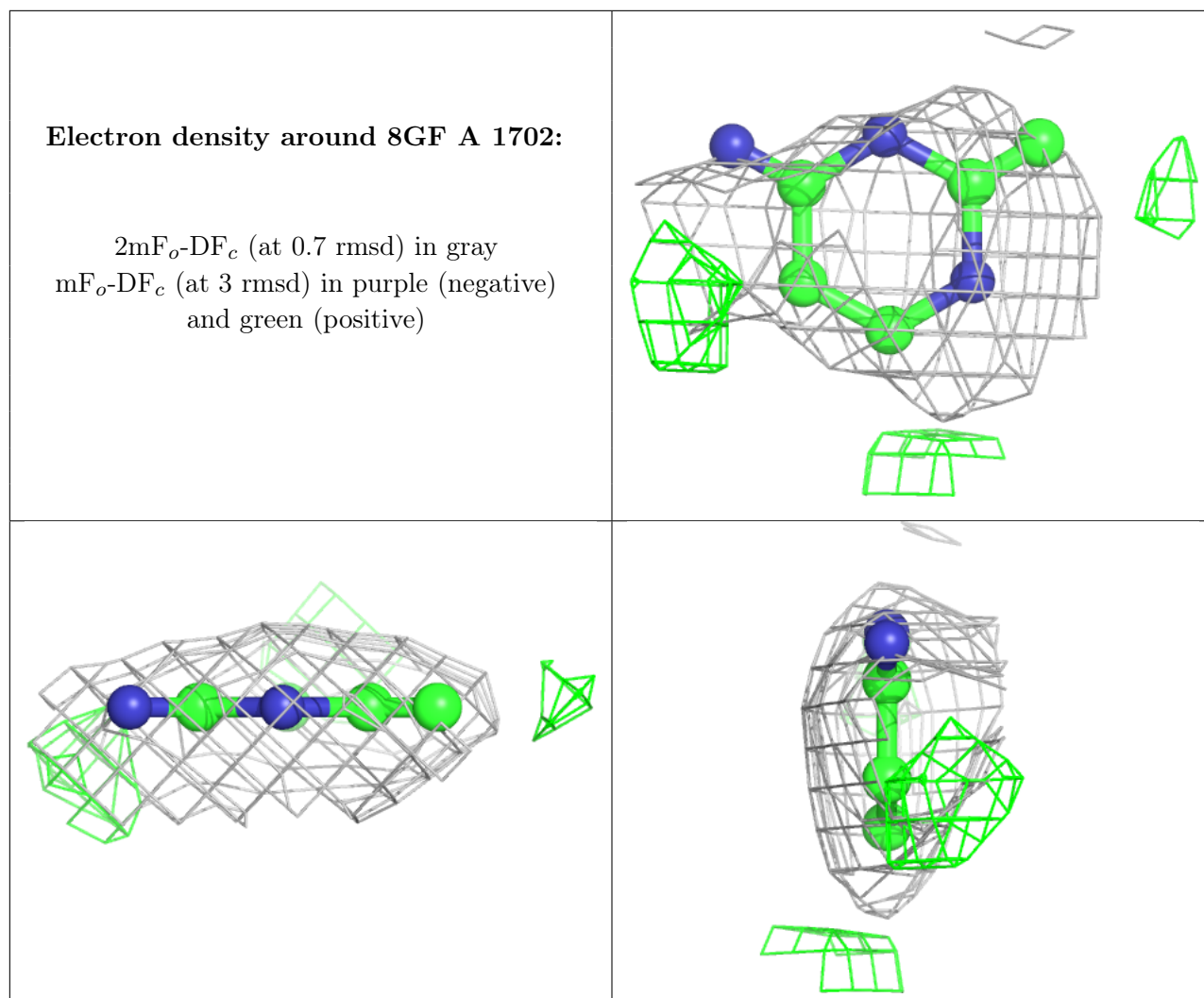
There are no monosaccharides in this entry.

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, 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.

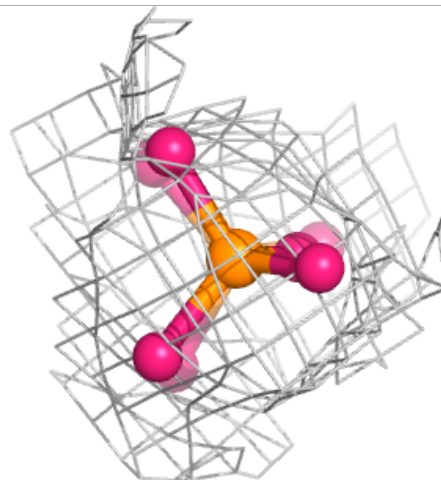
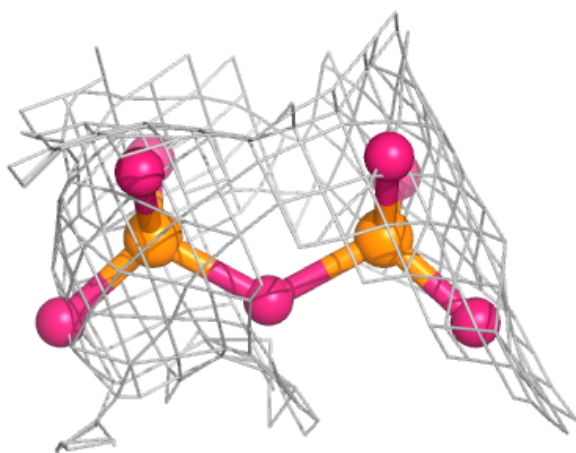
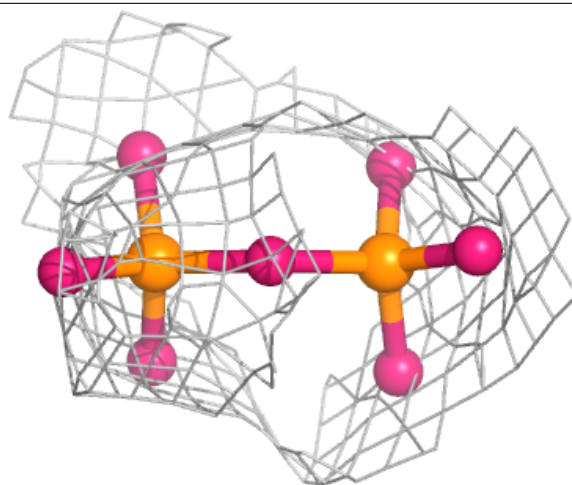
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	8GF	A	1702	8/8	0.80	0.46	103,106,109,112	0
5	DPO	A	1703	9/9	0.84	0.23	243,246,250,250	0
6	MG	A	1704	1/1	0.85	0.27	253,253,253,253	0
7	VAL	B	401	8/8	0.92	0.24	35,58,80,91	0
3	FAD	A	1701	53/53	0.94	0.25	63,79,98,106	0

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.



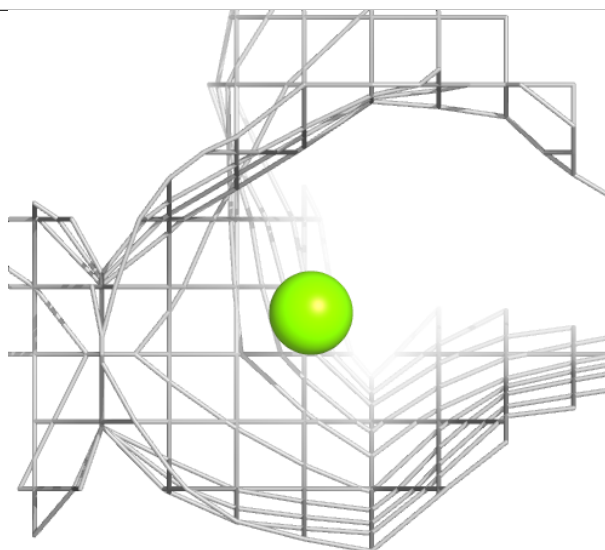
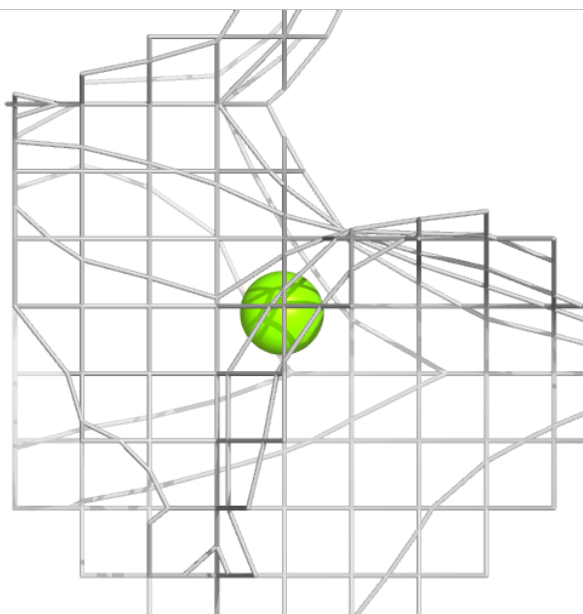
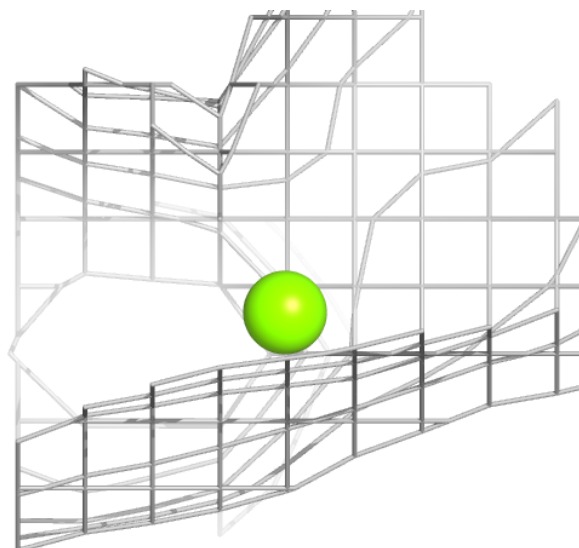
Electron density around DPO A 1703:

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



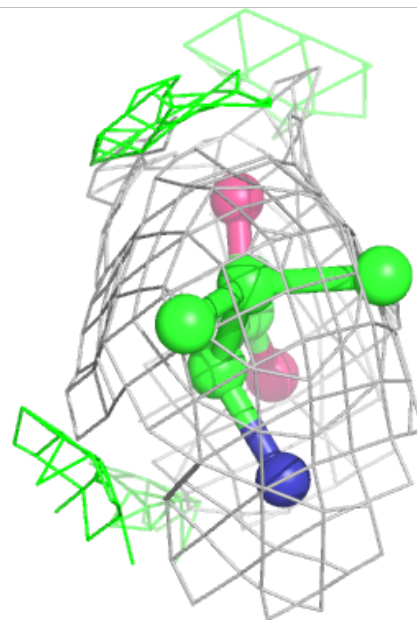
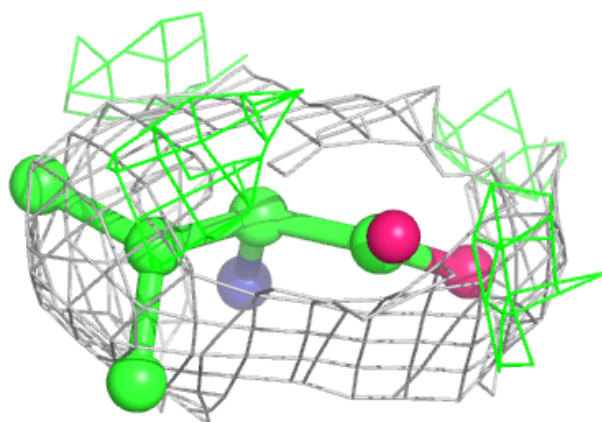
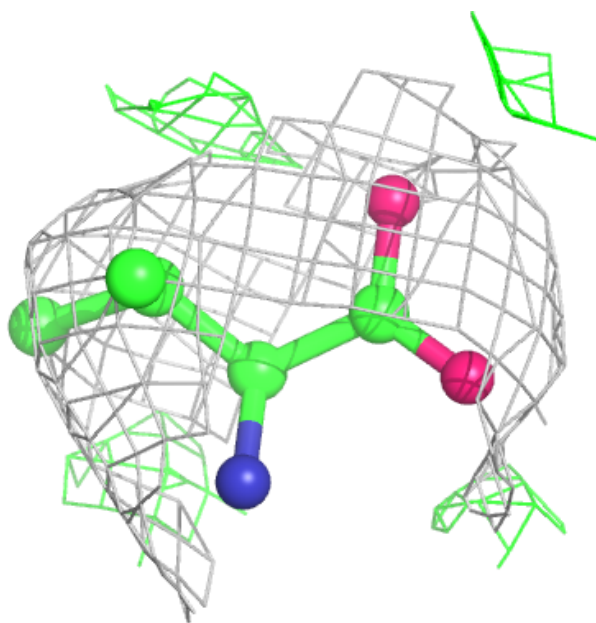
Electron density around MG A 1704:

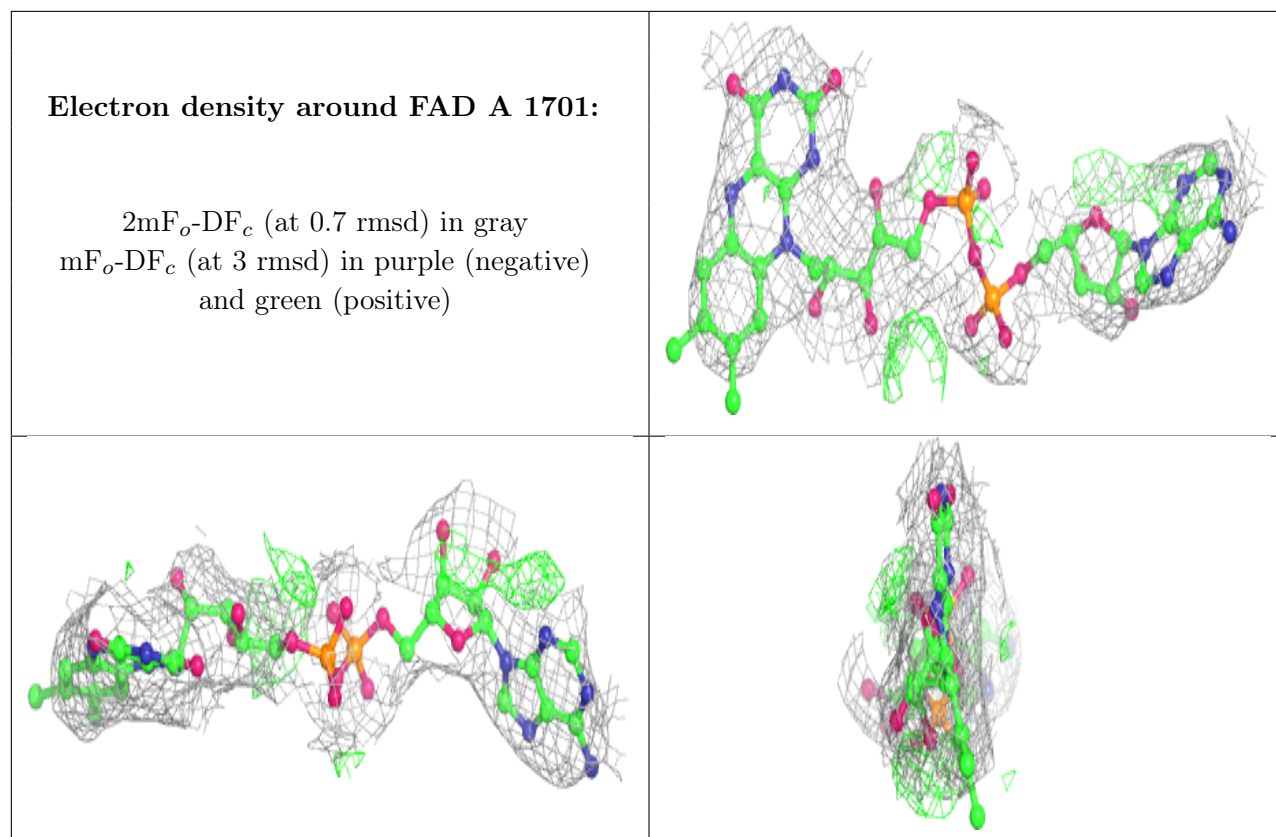
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around VAL 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)





6.5 Other polymers [i](#)

There are no such residues in this entry.