



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2025 – 11:33 am BST

PDB ID : 9R2J / pdb\_00009r2j  
Title : Crystal structure of human MAO B in complex with (E)-3-(4-nitrophenyl)-1-(3-(trifluoromethyl)phenyl)prop-2-en-1-one (chalcone inhibitor, 4a)  
Authors : Marchese, S.; Binda, C.  
Deposited on : 2025-04-30  
Resolution : 1.80 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](mailto: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>.

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

MolProbity : 4-5-2 with Phenix2.0rc1  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.45.1

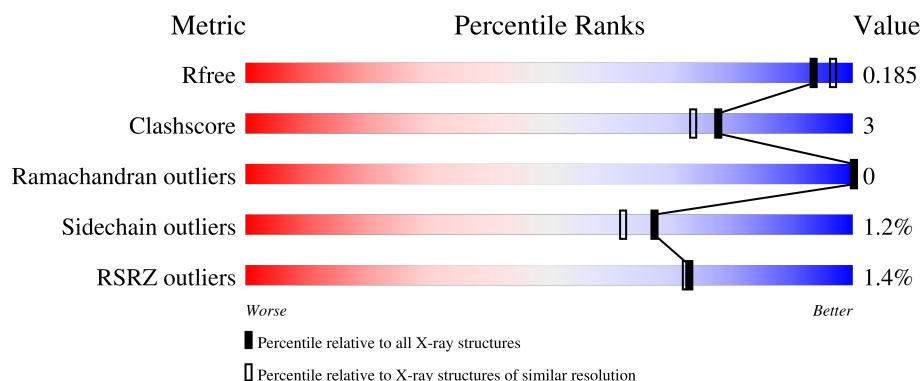
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	164625	7108 (1.80-1.80)
Clashscore	180529	8162 (1.80-1.80)
Ramachandran outliers	177936	8077 (1.80-1.80)
Sidechain outliers	177891	8076 (1.80-1.80)
RSRZ outliers	164620	7108 (1.80-1.80)

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  $\geq 3$ , 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	AAA	519	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div></div> </div> </div>
1	BBB	519	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>6%</div> <div>5%</div> </div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8968 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 Amine oxidase [flavin-containing] B.

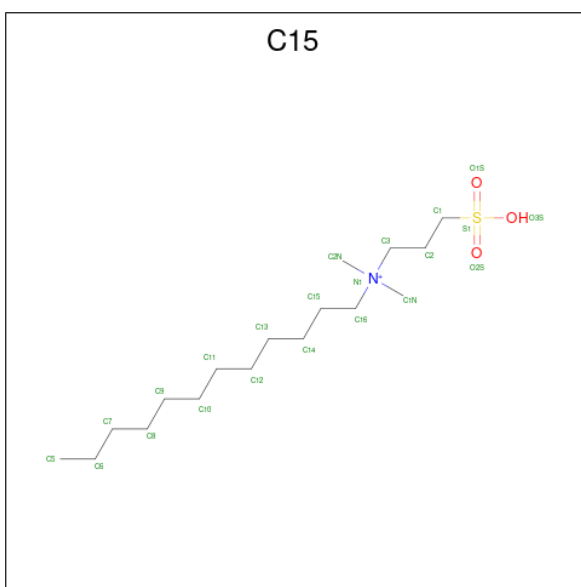
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	500	Total	C	N	O	S	0	2	0
			3988	2549	683	732	24			
1	BBB	495	Total	C	N	O	S	0	1	0
			3949	2524	677	724	24			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ) (labeled as "Ligand of Interest" by depositor).



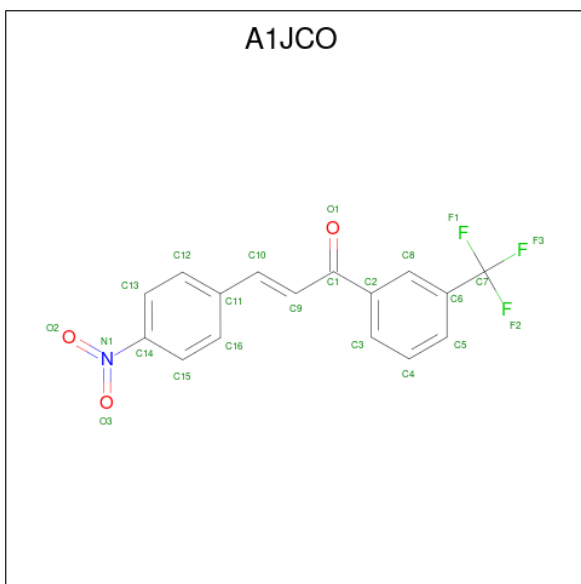
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	BBB	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is N-DODECYL-N,N-DIMETHYL-3-AMMONIO-1-PROPANESULFONATE (CCD ID: C15) (formula:  $C_{17}H_{38}NO_3S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	AAA	1	Total	C	N	O	S	0	0
			15	10	1	3	1		
3	BBB	1	Total	C	N	O	S	0	0
			11	6	1	3	1		

- Molecule 4 is ( {E})-3-(4-nitrophenyl)-1-[3-(trifluoromethyl)phenyl]prop-2-en-1-one (CCD ID: A1JCO) (formula: C<sub>16</sub>H<sub>10</sub>F<sub>3</sub>NO<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	AAA	1	Total	C	F	N	O	0	0
			23	16	3	1	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	BBB	1	Total	C	F	N	O	0	0
			23	16	3	1	3		

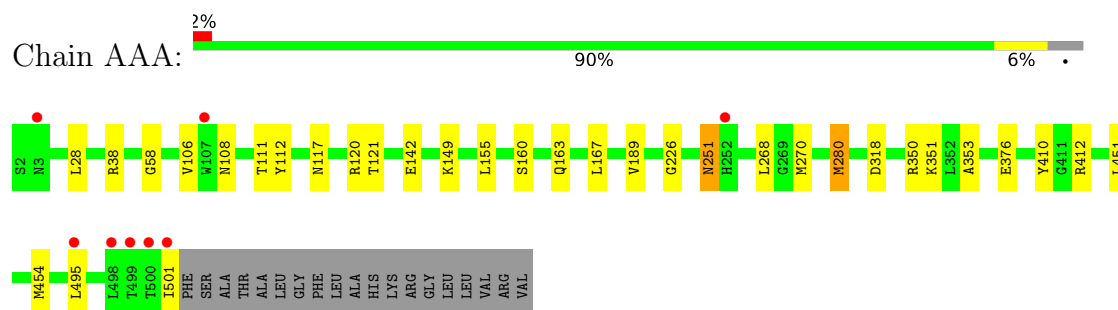
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	408	Total	O	0	0
			408	408		
5	BBB	445	Total	O	0	0
			445	445		

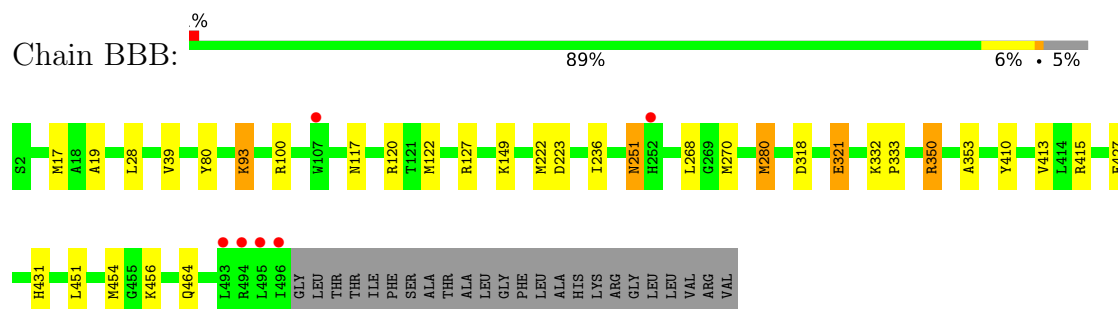
### 3 Residue-property plots [i](#)

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: Amine oxidase [flavin-containing] B



- Molecule 1: Amine oxidase [flavin-containing] B



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	132.34Å 223.93Å 86.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.68 – 1.80 47.68 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.68-1.80) 100.0 (47.68-1.80)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.52 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.144 , 0.173 0.158 , 0.185	Depositor DCC
$R_{free}$ test set	3000 reflections (2.52%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	16.4	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 39.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.007 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.011 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8968	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, A1JCO, C15

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	AAA	1.18	6/4091 (0.1%)	1.21	0/5553
1	BBB	1.17	4/4049 (0.1%)	1.21	5/5495 (0.1%)
All	All	1.18	10/8140 (0.1%)	1.21	5/11048 (0.0%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	142	GLU	C-O	6.26	1.31	1.24
1	BBB	19	ALA	C-O	5.73	1.30	1.24
1	BBB	223	ASP	C-O	5.66	1.30	1.24
1	AAA	38	ARG	C-O	5.53	1.30	1.23
1	AAA	160	SER	C-O	5.51	1.30	1.24
1	AAA	189	VAL	C-O	5.46	1.30	1.24
1	AAA	376	GLU	CD-OE2	5.44	1.35	1.25
1	BBB	80	TYR	C-O	5.38	1.30	1.23
1	BBB	236	ILE	N-CA	5.23	1.51	1.46
1	AAA	226	GLY	C-O	5.09	1.30	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	127	ARG	CG-CD-NE	-5.39	100.14	112.00
1	BBB	431	HIS	CA-CB-CG	5.36	119.16	113.80
1	BBB	318	ASP	CA-CB-CG	5.26	117.86	112.60
1	BBB	427	GLU	CB-CG-CD	5.05	121.19	112.60
1	BBB	321	GLU	CB-CG-CD	5.05	121.19	112.60

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	AAA	3988	0	3992	25	0
1	BBB	3949	0	3947	25	1
2	AAA	53	0	29	1	0
2	BBB	53	0	29	0	0
3	AAA	15	0	21	0	0
3	BBB	11	0	13	0	0
4	AAA	23	0	0	0	0
4	BBB	23	0	0	0	0
5	AAA	408	0	0	3	0
5	BBB	445	0	0	4	1
All	All	8968	0	8031	44	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 3.

All (44) 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:AAA:28:LEU:HD11	1:AAA:454:MET:HE1	1.53	0.90
1:BBB:451:LEU:HA	1:BBB:454:MET:HE2	1.68	0.75
1:AAA:251:ASN:HD22	1:AAA:251:ASN:H	1.34	0.75
1:BBB:117:ASN:HD22	1:BBB:120:ARG:HH21	1.35	0.75
1:AAA:106:VAL:HG21	1:AAA:111:THR:HG22	1.75	0.69
1:BBB:28:LEU:HD11	1:BBB:454:MET:HE1	1.74	0.68
1:AAA:117:ASN:HD22	1:AAA:120:ARG:HH21	1.43	0.66
1:AAA:353:ALA:HB1	1:BBB:280:MET:HE1	1.78	0.65
1:BBB:251:ASN:HD22	1:BBB:251:ASN:H	1.45	0.64
1:AAA:106:VAL:HG21	1:AAA:111:THR:CG2	2.30	0.62
1:BBB:451:LEU:HD23	1:BBB:454:MET:HE1	1.81	0.61
1:AAA:351:LYS:NZ	5:AAA:701:HOH:O	2.32	0.61
1:BBB:451:LEU:HD23	1:BBB:454:MET:CE	2.32	0.59
1:AAA:280:MET:HE1	1:BBB:353:ALA:HB1	1.83	0.59
1:AAA:451:LEU:HD23	1:AAA:454:MET:CE	2.33	0.58
1:BBB:410:TYR:O	1:BBB:413:VAL:HG13	2.04	0.57
1:BBB:100:ARG:HD3	5:BBB:863:HOH:O	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:270:MET:HE1	1:BBB:268:LEU:HD23	1.92	0.52
1:AAA:451:LEU:HA	1:AAA:454:MET:HE2	1.91	0.51
1:BBB:93:LYS:HA	1:BBB:93:LYS:CE	2.40	0.51
1:AAA:451:LEU:HD23	1:AAA:454:MET:HE1	1.94	0.50
1:AAA:108:ASN:OD1	1:AAA:108:ASN:C	2.55	0.50
1:BBB:332:LYS:HB3	1:BBB:333:PRO:HD2	1.94	0.49
5:AAA:802:HOH:O	1:BBB:149:LYS:HD3	2.12	0.48
1:BBB:17:MET:CE	1:BBB:39:VAL:HG11	2.43	0.48
1:AAA:163:GLN:HB3	1:AAA:318:ASP:OD2	2.13	0.48
1:AAA:353:ALA:CB	1:BBB:280:MET:HE1	2.44	0.48
1:AAA:410:TYR:OH	1:BBB:350:ARG:HD2	2.14	0.48
5:AAA:713:HOH:O	1:BBB:280:MET:CE	2.61	0.47
1:BBB:451:LEU:HA	1:BBB:454:MET:CE	2.42	0.45
1:AAA:149:LYS:HD3	5:BBB:983:HOH:O	2.17	0.45
1:BBB:464:GLN:NE2	5:BBB:702:HOH:O	2.31	0.44
1:AAA:280:MET:HE3	1:AAA:280:MET:HB3	1.76	0.44
1:BBB:415:ARG:NH1	5:BBB:707:HOH:O	2.51	0.43
1:BBB:122:MET:HE3	1:BBB:122:MET:HB3	1.93	0.42
1:AAA:106:VAL:CG1	1:AAA:112:TYR:HA	2.50	0.42
1:BBB:321:GLU:H	1:BBB:321:GLU:CD	2.27	0.42
1:AAA:58:GLY:HA2	2:AAA:601:FAD:C4X	2.50	0.41
1:AAA:121:THR:HG21	1:AAA:155:LEU:HD21	2.01	0.41
1:AAA:167:LEU:C	1:AAA:167:LEU:HD23	2.45	0.41
1:AAA:268:LEU:HD23	1:BBB:270:MET:HE1	2.02	0.41
1:AAA:106:VAL:HG11	1:AAA:112:TYR:HA	2.03	0.41
1:BBB:454:MET:HE3	1:BBB:456:LYS:HD2	2.03	0.40
1:AAA:251:ASN:HD22	1:AAA:251:ASN:N	2.12	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)
1:BBB:222:MET:CE	5:BBB:766:HOH:O[4_565]	2.10	0.10

## 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	AAA	500/519 (96%)	488 (98%)	12 (2%)	0	100	100
1	BBB	494/519 (95%)	479 (97%)	15 (3%)	0	100	100
All	All	994/1038 (96%)	967 (97%)	27 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	AAA	431/443 (97%)	425 (99%)	6 (1%)	62	56
1	BBB	426/443 (96%)	422 (99%)	4 (1%)	75	72
All	All	857/886 (97%)	847 (99%)	10 (1%)	67	62

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	251	ASN
1	AAA	280	MET
1	AAA	350	ARG
1	AAA	412	ARG
1	AAA	495	LEU
1	AAA	501	ILE
1	BBB	93	LYS
1	BBB	251	ASN
1	BBB	280	MET
1	BBB	350	ARG

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

6 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	C15	AAA	602	-	14,14,21	0.80	1 (7%)	18,19,26	0.62	0
4	A1JCO	BBB	603	-	24,24,24	2.46	7 (29%)	34,34,34	1.14	3 (8%)
4	A1JCO	AAA	603	-	24,24,24	2.11	6 (25%)	34,34,34	1.49	6 (17%)
2	FAD	AAA	601	1	53,58,58	1.23	6 (11%)	68,89,89	1.65	14 (20%)
3	C15	BBB	602	-	10,10,21	0.97	1 (10%)	14,15,26	1.00	1 (7%)
2	FAD	BBB	601	1	53,58,58	1.41	6 (11%)	68,89,89	1.49	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
3	C15	AAA	602	-	-	0/14/14/21	-
4	A1JCO	BBB	603	-	-	0/19/19/19	0/2/2/2
4	A1JCO	AAA	603	-	-	0/19/19/19	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	AAA	601	1	-	1/30/50/50	0/6/6/6
3	C15	BBB	602	-	-	4/8/8/21	-
2	FAD	BBB	601	1	-	1/30/50/50	0/6/6/6

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	BBB	603	A1JCO	O3-N1	8.40	1.42	1.22
4	AAA	603	A1JCO	O3-N1	5.86	1.36	1.22
4	BBB	603	A1JCO	C9-C1	4.46	1.55	1.47
4	AAA	603	A1JCO	C2-C1	-4.38	1.42	1.49
4	BBB	603	A1JCO	C2-C1	-4.33	1.42	1.49
2	BBB	601	FAD	C4X-N5	4.22	1.39	1.30
2	BBB	601	FAD	C9-C9A	3.86	1.45	1.39
2	AAA	601	FAD	C10-N1	3.85	1.41	1.33
2	AAA	601	FAD	C4X-N5	3.29	1.37	1.30
2	BBB	601	FAD	C10-N1	3.15	1.39	1.33
4	AAA	603	A1JCO	O2-N1	2.93	1.29	1.22
2	AAA	601	FAD	O4B-C1B	2.89	1.45	1.41
4	AAA	603	A1JCO	C9-C1	2.78	1.52	1.47
4	AAA	603	A1JCO	C14-N1	-2.75	1.40	1.47
3	BBB	602	C15	O3S-S1	2.70	1.57	1.47
2	BBB	601	FAD	C2B-C1B	-2.64	1.49	1.53
3	AAA	602	C15	O3S-S1	2.58	1.56	1.47
2	BBB	601	FAD	O4-C4	2.53	1.28	1.23
4	BBB	603	A1JCO	F2-C7	-2.51	1.23	1.32
4	AAA	603	A1JCO	F3-C7	-2.34	1.24	1.32
2	AAA	601	FAD	C2A-N3A	2.31	1.35	1.32
4	BBB	603	A1JCO	C13-C12	2.29	1.42	1.38
4	BBB	603	A1JCO	C14-N1	-2.21	1.41	1.47
2	BBB	601	FAD	C9A-N10	-2.14	1.37	1.41
2	AAA	601	FAD	O2-C2	2.13	1.28	1.24
4	BBB	603	A1JCO	C7-C6	-2.11	1.45	1.49
2	AAA	601	FAD	C2-N1	2.10	1.41	1.36

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	601	FAD	C9A-C5X-N5	-4.49	117.55	122.43
2	BBB	601	FAD	N3A-C2A-N1A	-4.41	121.78	128.68
4	AAA	603	A1JCO	C13-C12-C11	-3.76	116.35	121.25
2	AAA	601	FAD	C4-N3-C2	-3.74	118.73	125.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	601	FAD	O4B-C1B-C2B	-3.69	101.54	106.93
4	AAA	603	A1JCO	C12-C13-C14	3.29	122.40	118.73
2	AAA	601	FAD	C8M-C8-C7	3.24	127.39	120.74
2	AAA	601	FAD	C8M-C8-C9	-3.20	113.58	119.49
2	AAA	601	FAD	C4X-C10-N10	3.19	121.14	116.48
2	AAA	601	FAD	C10-C4X-N5	-3.08	118.32	124.86
2	AAA	601	FAD	C4X-C4-N3	2.90	120.55	113.19
2	BBB	601	FAD	C3B-C2B-C1B	2.85	105.26	100.98
2	AAA	601	FAD	C5A-C6A-N6A	2.84	124.67	120.35
2	BBB	601	FAD	C9-C9A-N10	2.81	125.63	121.84
4	AAA	603	A1JCO	O1-C1-C9	-2.80	116.40	120.76
2	BBB	601	FAD	C8M-C8-C7	2.74	126.35	120.74
2	BBB	601	FAD	C4X-C10-N10	2.71	120.44	116.48
2	AAA	601	FAD	C4-C4X-N5	2.66	122.02	118.23
2	BBB	601	FAD	C10-C4X-N5	-2.48	119.60	124.86
4	AAA	603	A1JCO	C2-C1-C9	2.47	122.01	119.23
4	AAA	603	A1JCO	C15-C14-C13	-2.43	119.21	122.22
2	BBB	601	FAD	C9A-C5X-N5	-2.41	119.81	122.43
2	BBB	601	FAD	C8M-C8-C9	-2.35	115.14	119.49
2	AAA	601	FAD	O4B-C1B-C2B	-2.35	103.49	106.93
2	BBB	601	FAD	C10-N1-C2	2.33	121.56	116.90
4	BBB	603	A1JCO	C13-C12-C11	-2.33	118.22	121.25
2	AAA	601	FAD	C9A-N10-C10	-2.31	117.17	120.77
2	BBB	601	FAD	C4X-C4-N3	2.31	119.05	113.19
3	BBB	602	C15	O1S-S1-C1	-2.19	104.28	106.92
4	AAA	603	A1JCO	C13-C14-N1	2.16	122.10	118.92
2	BBB	601	FAD	C4X-C10-N1	-2.14	119.75	124.73
2	AAA	601	FAD	C1'-N10-C9A	2.12	124.05	120.51
2	AAA	601	FAD	O3'-C3'-C4'	-2.11	103.73	108.81
4	BBB	603	A1JCO	F3-C7-C6	-2.10	108.31	112.93
4	BBB	603	A1JCO	C8-C6-C7	2.07	122.37	119.58
2	AAA	601	FAD	O4B-C4B-C3B	2.06	109.20	105.11
2	BBB	601	FAD	C4-C4X-N5	2.05	121.14	118.23

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	BBB	602	C15	C2-C1-S1-O2S
3	BBB	602	C15	S1-C1-C2-C3
3	BBB	602	C15	C2-C1-S1-O1S
3	BBB	602	C15	C2-C1-S1-O3S

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Mol	Chain	Res	Type	Atoms
2	BBB	601	FAD	O4B-C4B-C5B-O5B
2	AAA	601	FAD	O4B-C4B-C5B-O5B

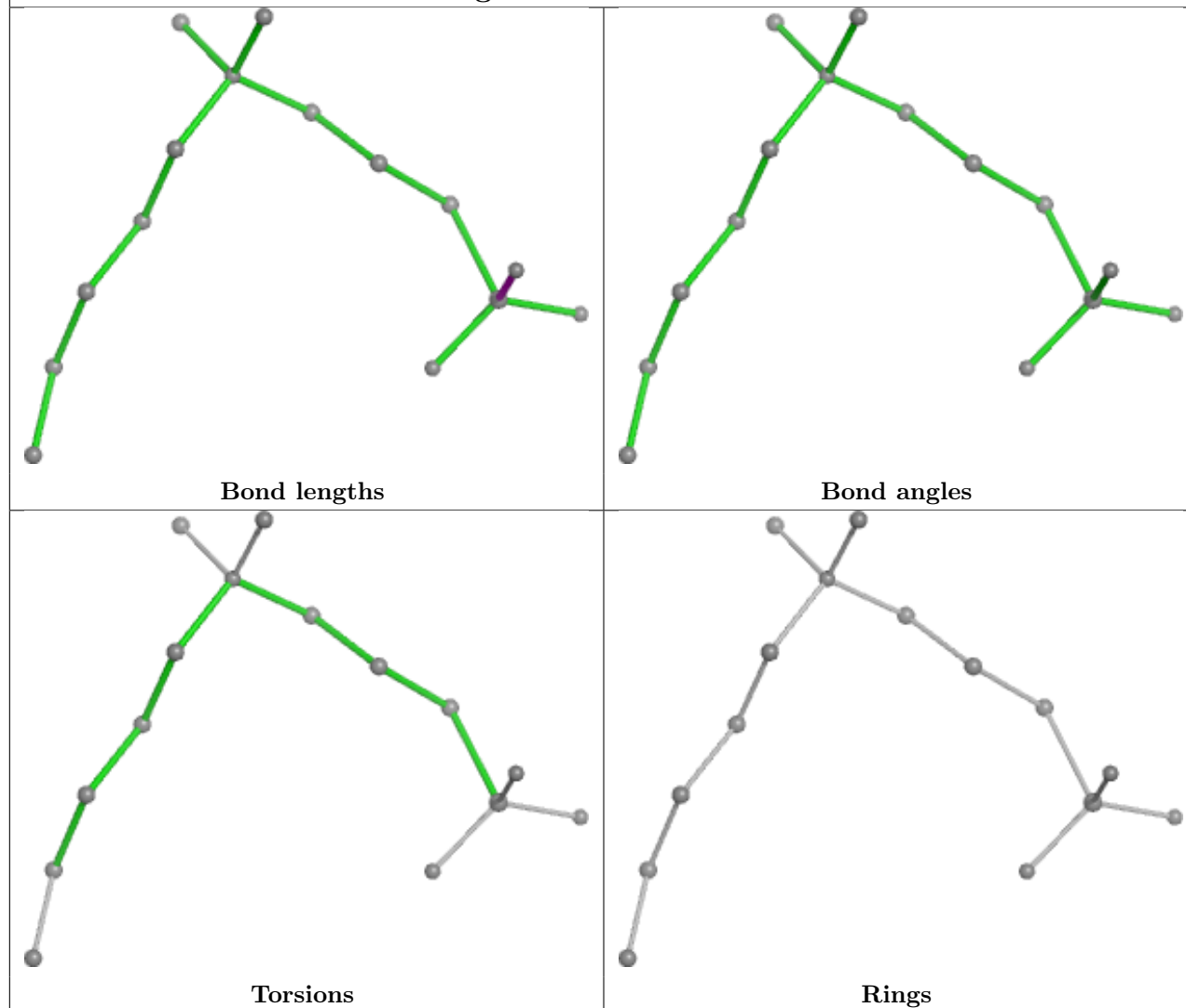
There are no ring outliers.

1 monomer is involved in 1 short contact:

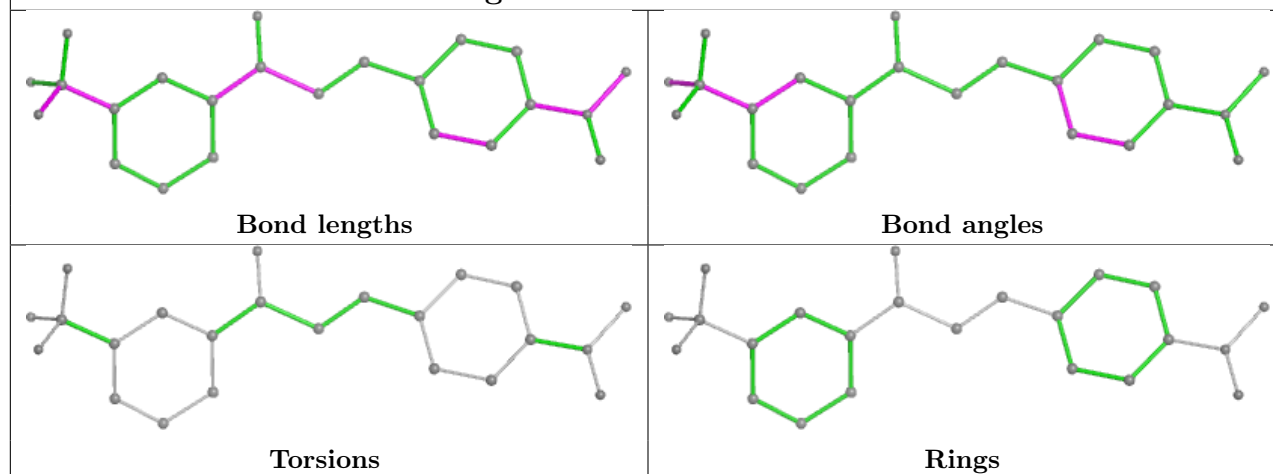
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	AAA	601	FAD	1	0

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 C15 AAA 602

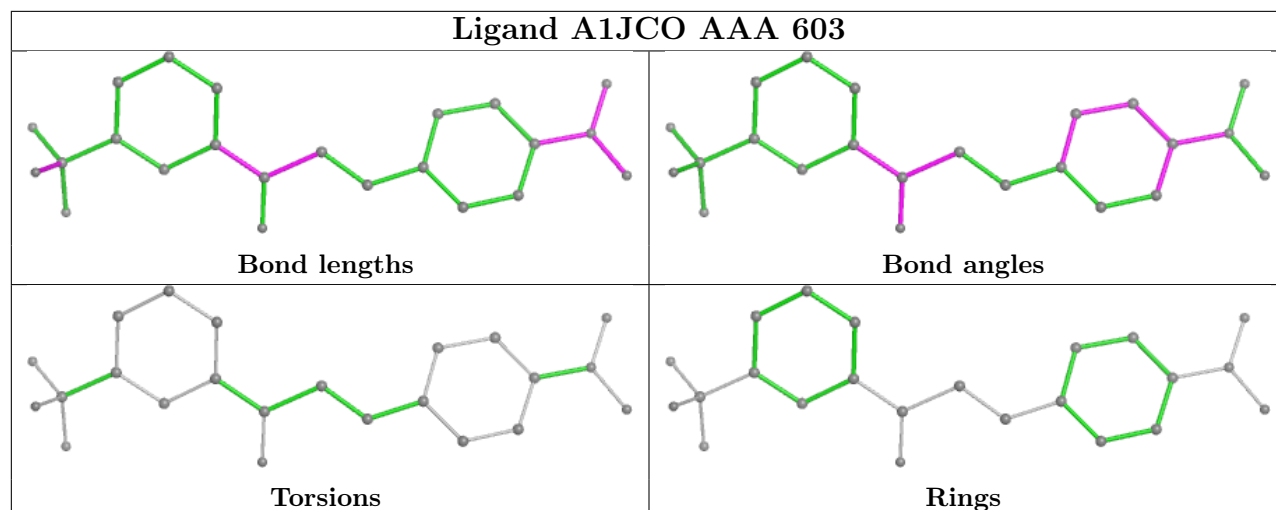


## Ligand A1JCO BBB 603

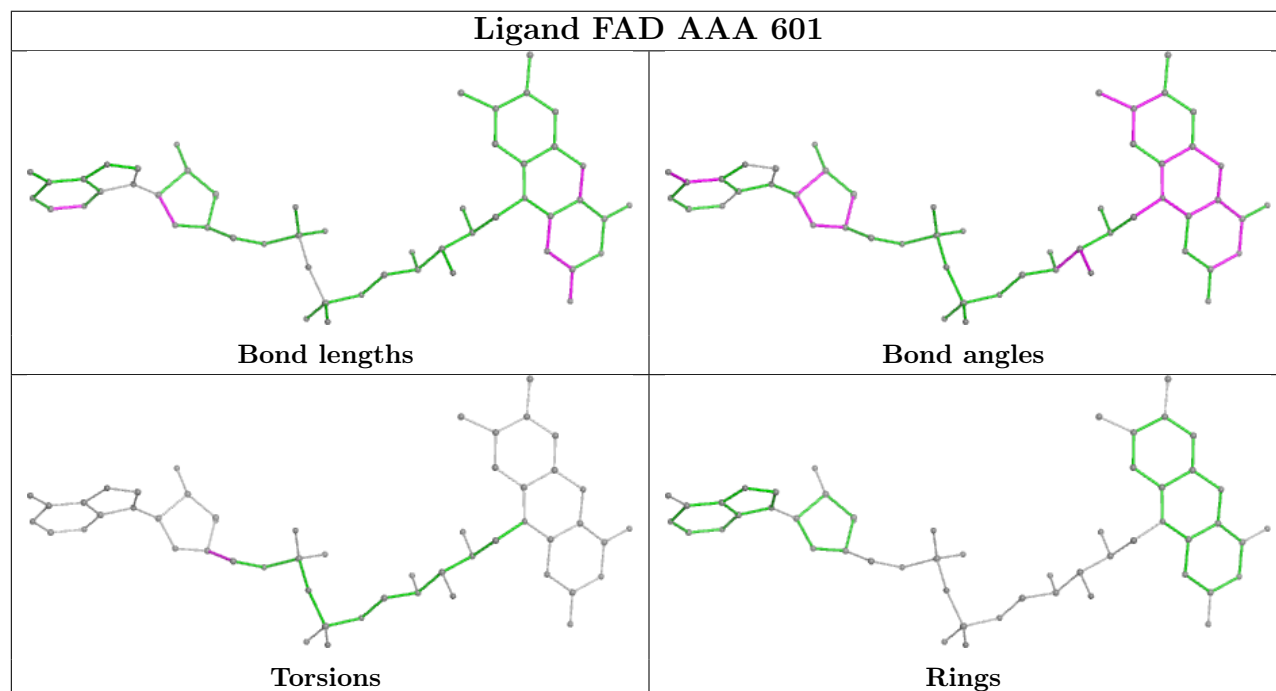


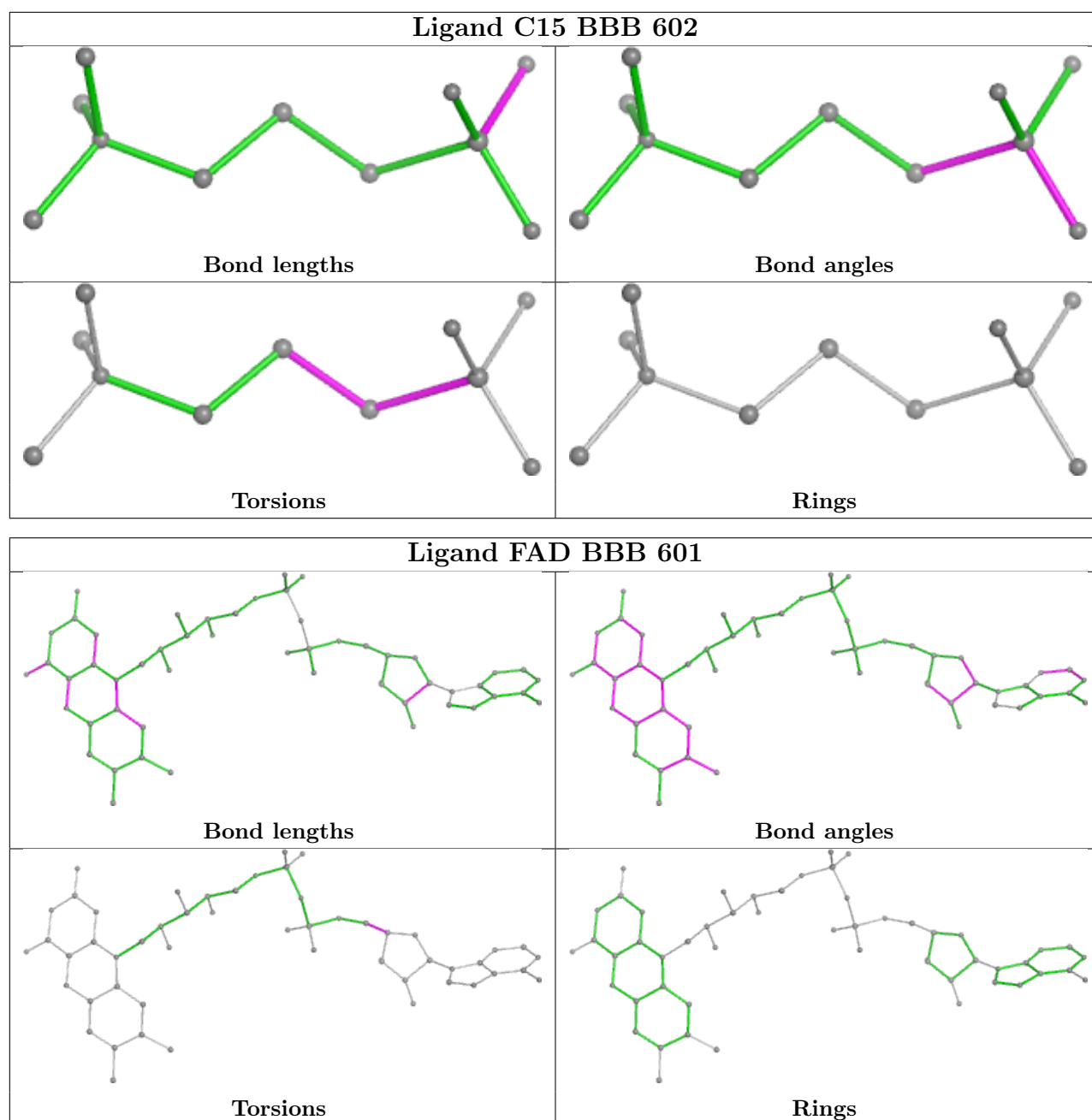


## Ligand A1JCO AAA 603



## Ligand FAD AAA 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	AAA	500/519 (96%)	-0.41	8 (1%) 70 69	9, 17, 34, 73	2 (0%)
1	BBB	495/519 (95%)	-0.52	6 (1%) 76 76	10, 15, 33, 69	1 (0%)
All	All	995/1038 (95%)	-0.46	14 (1%) 73 72	9, 16, 34, 73	3 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	501	ILE	6.6
1	BBB	496	ILE	4.3
1	BBB	495	LEU	4.3
1	BBB	494	ARG	3.3
1	AAA	498	LEU	3.2
1	AAA	252	HIS	3.1
1	AAA	500	THR	2.8
1	BBB	252	HIS	2.7
1	BBB	107	TRP	2.6
1	AAA	107	TRP	2.3
1	BBB	493	LEU	2.3
1	AAA	499	THR	2.2
1	AAA	3	ASN	2.2
1	AAA	495	LEU	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

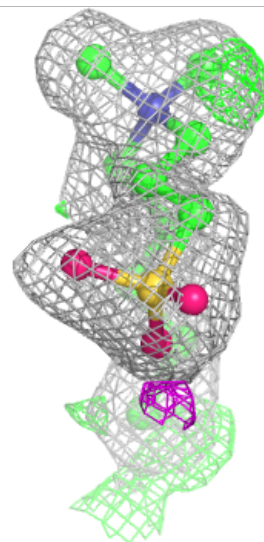
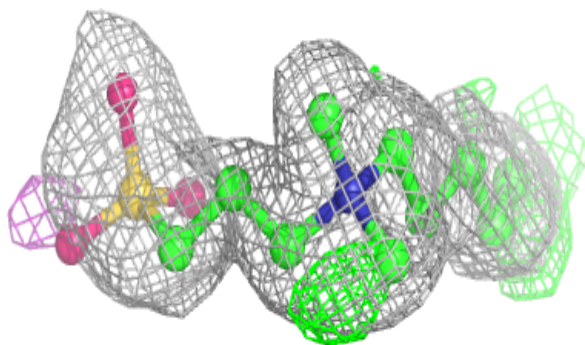
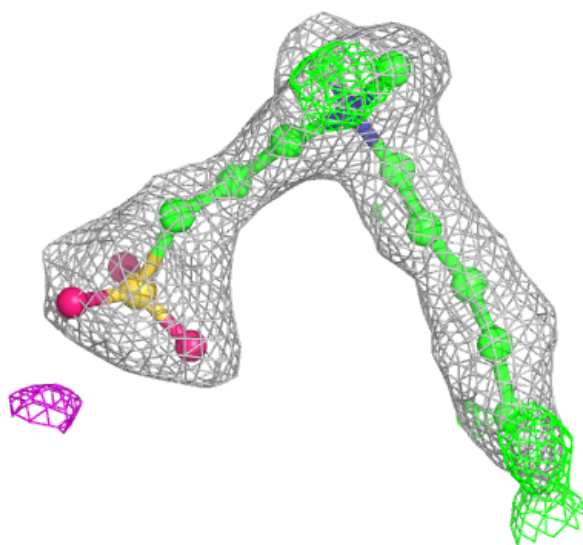
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	C15	AAA	602	15/22	0.77	0.16	28,36,66,73	0
3	C15	BBB	602	11/22	0.80	0.14	32,40,66,71	0
4	A1JCO	AAA	603	23/23	0.95	0.06	16,18,28,36	0
4	A1JCO	BBB	603	23/23	0.96	0.06	15,18,27,33	0
2	FAD	AAA	601	53/53	0.99	0.03	10,11,13,13	0
2	FAD	BBB	601	53/53	0.99	0.03	9,11,12,12	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.

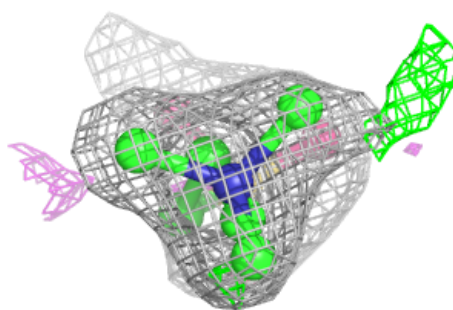
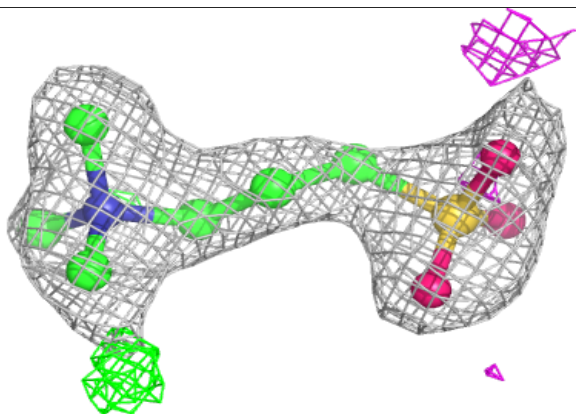
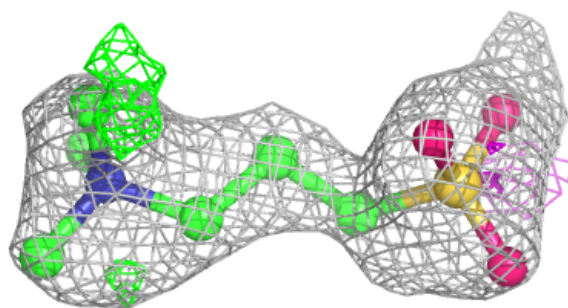
**Electron density around C15 AAA 602:**

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

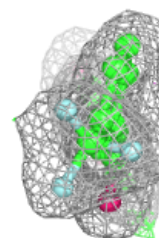
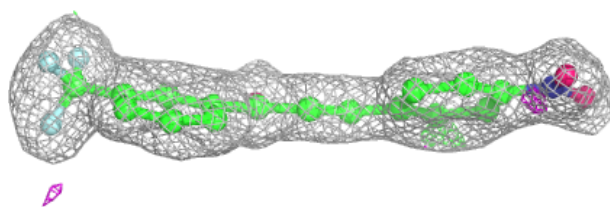
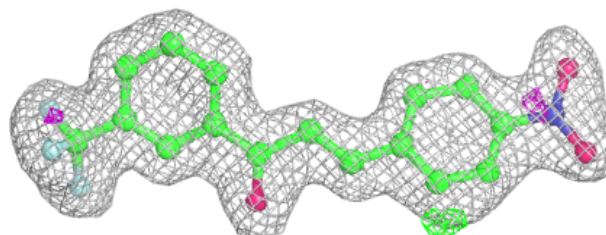


**Electron density around C15 BBB 602:**

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

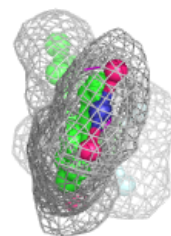
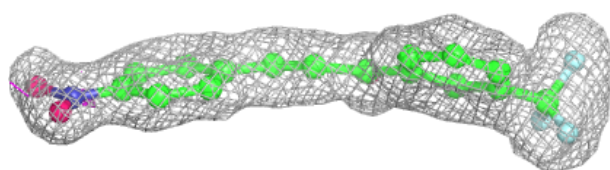
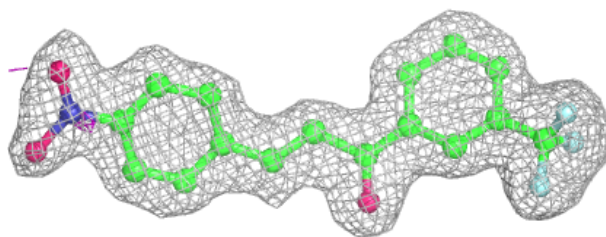
**Electron density around A1JCO AAA 603:**

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

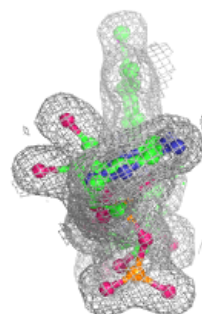
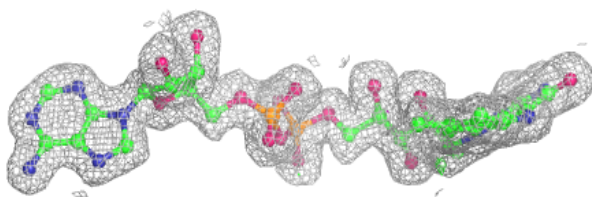
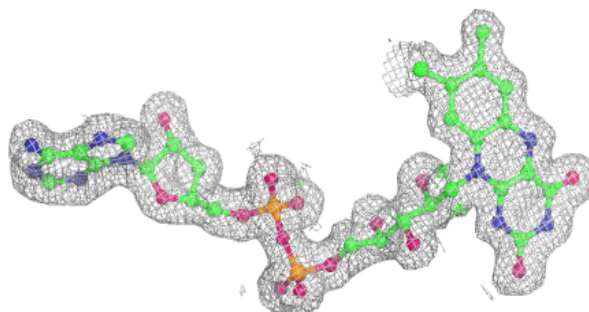


**Electron density around A1JCO BBB 603:**

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

**Electron density around FAD AAA 601:**

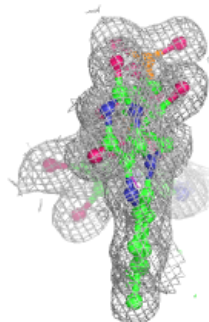
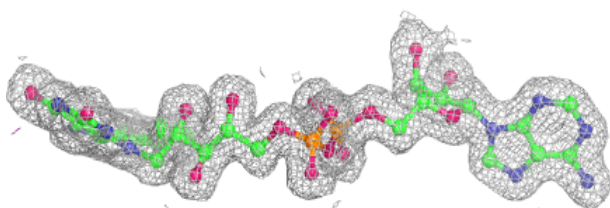
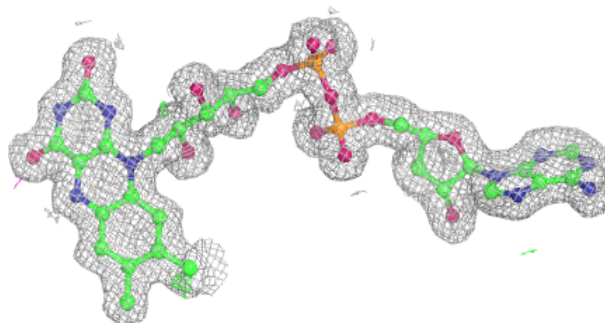
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around FAD BBB 601:**

$2mF_o - DF_c$  (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.