



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:31 AM UTC

PDB ID : 9FJT / pdb_00009fjt
Title : Human Monoamine Oxidase B in complex with MC4762 inhibitor (9a) at 1.4 Å resolution
Authors : Gottinger, A.; Binda, C.
Deposited on : 2024-05-31
Resolution : 1.40 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

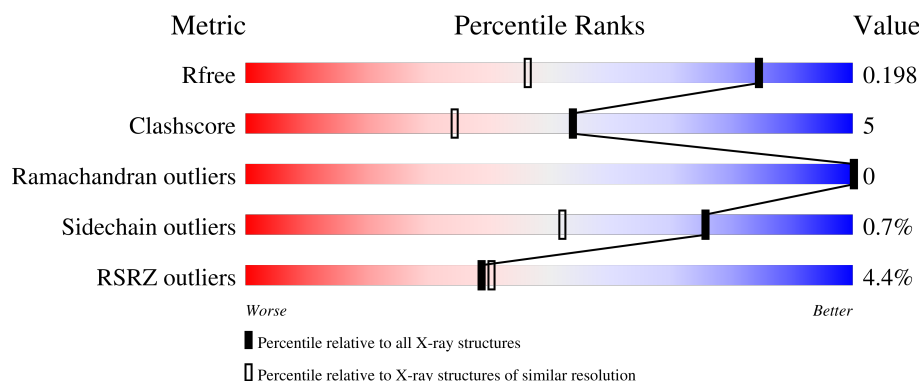
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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	2563 (1.40-1.40)
Clashscore	190562	2660 (1.40-1.40)
Ramachandran outliers	187476	2611 (1.40-1.40)
Sidechain outliers	187428	2610 (1.40-1.40)
RSRZ outliers	180081	2561 (1.40-1.40)

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	AAA	520	5% 89% 7% .
1	BBB	520	3% 88% 7% 5%

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	A1IDI	AAA	603	-	X	-	-
4	A1IDI	BBB	603	-	-	X	-

2 Entry composition [i](#)

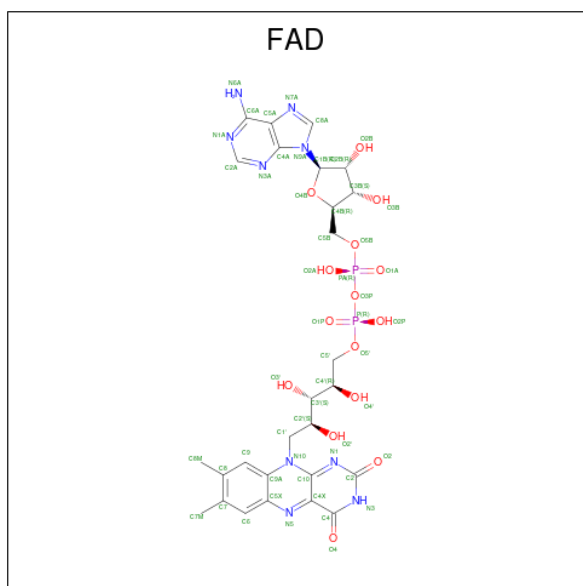
There are 5 unique types of molecules in this entry. The entry contains 9144 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

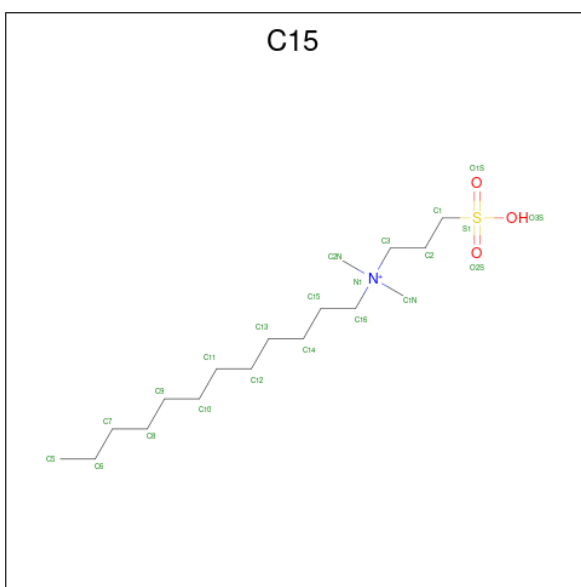
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	498	Total	C	N	O	S	0	8	0
			4001	2562	681	730	28			
1	BBB	495	Total	C	N	O	S	0	7	0
			3983	2552	678	724	29			

- 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).



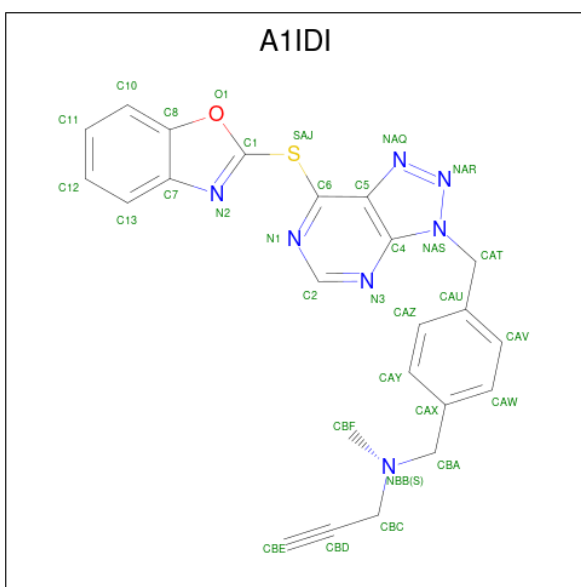
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	BBB	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is N-DODECYL-N,N-DIMETHYL-3-AMMONIO-1-PROPANESULFONATE (CCD ID: C15) (formula: $C_{17}H_{38}NO_3S$) (labeled as "Ligand of Interest" by depositor).



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 {N}-[[4-[[7-(1,3-benzoxazol-2-ylsulfanyl)-[1,2,3]triazolo[4,5-d]pyrimidin-3-yl]methyl]phenyl]methyl]- {N}-methyl-prop-2-yn-1-amine (CCD ID: A1IDI) (formula: C₂₃H₁₉N₇OS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	AAA	1	Total	C	N	S	0	0
			23	16	6	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	BBB	1	Total	C	N	O	S	0	0
			32	23	7	1	1		

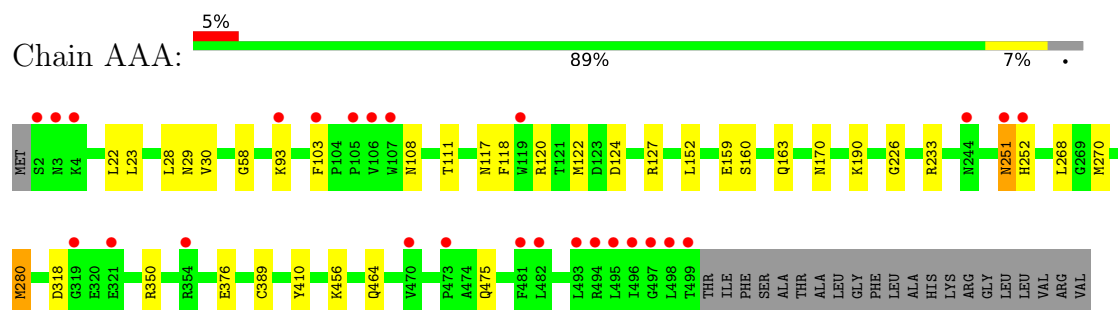
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	458	Total	O	0	0
			458	458		
5	BBB	515	Total	O	0	0
			515	515		

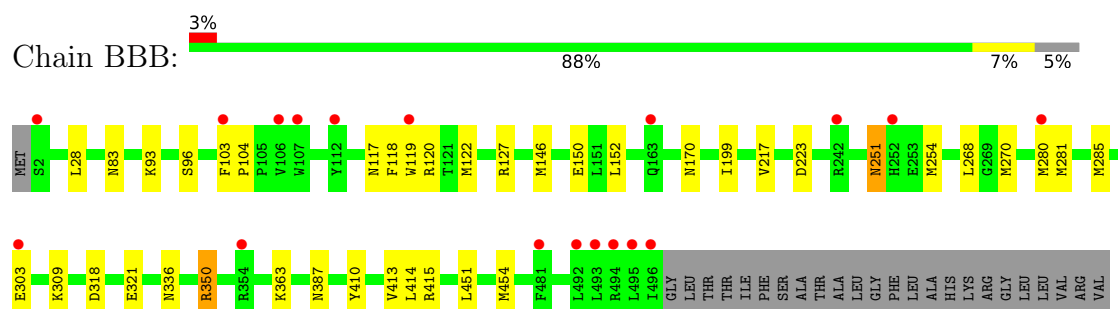
3 Residue-property plots [i](#)

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, α , β , γ	130.92Å 222.03Å 85.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.18 – 1.40 47.18 – 1.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.18-1.40) 99.7 (47.18-1.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.167 , 0.190 0.176 , 0.198	Depositor DCC
R_{free} test set	6176 reflections (2.53%)	wwPDB-VP
Wilson B-factor (Å ²)	11.8	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 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	9144	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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: C15, FAD, A1IDI

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	1.04	3/4122 (0.1%)	1.17	4/5591 (0.1%)
1	BBB	1.05	2/4103 (0.0%)	1.18	8/5566 (0.1%)
All	All	1.05	5/8225 (0.1%)	1.18	12/11157 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	376	GLU	CD-OE2	6.73	1.38	1.25
1	AAA	226	GLY	C-O	5.57	1.31	1.24
1	BBB	96	SER	CA-CB	-5.53	1.45	1.53
1	BBB	217	VAL	C-O	5.25	1.30	1.24
1	AAA	464	GLN	C-O	5.08	1.29	1.24

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BBB	415	ARG	CB-CA-C	8.96	125.86	110.72
1	AAA	252	HIS	CA-CB-CG	-7.39	106.41	113.80
1	AAA	127	ARG	CG-CD-NE	-6.74	97.18	112.00
1	AAA	170	ASN	CA-CB-CG	6.05	118.65	112.60
1	BBB	415	ARG	CG-CD-NE	-6.03	98.74	112.00
1	BBB	336	ASN	CA-CB-CG	-5.96	106.64	112.60
1	BBB	350	ARG	CG-CD-NE	5.44	123.97	112.00
1	BBB	127	ARG	CG-CD-NE	-5.43	100.06	112.00
1	BBB	170	ASN	CA-CB-CG	5.32	117.92	112.60
1	BBB	83	ASN	CA-CB-CG	-5.25	107.35	112.60
1	BBB	223	ASP	N-CA-C	-5.23	105.47	111.07
1	AAA	124	ASP	CA-CB-CG	5.11	117.71	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	4001	0	4026	30	2
1	BBB	3983	0	3998	56	0
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	32	0	0	17	0
5	AAA	458	0	0	3	0
5	BBB	515	0	0	11	0
All	All	9144	0	8116	79	2

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 5.

All (79) 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)
4:BBB:603:A1IDI:C13	5:BBB:1015:HOH:O	1.84	1.25
1:BBB:303:GLU:HG2	5:BBB:1117:HOH:O	1.60	1.02
1:BBB:103:PHE:CD1	4:BBB:603:A1IDI:C10	2.44	1.00
1:BBB:103:PHE:CE1	4:BBB:603:A1IDI:C10	2.48	0.96
1:BBB:285[B]:MET:SD	1:BBB:414:LEU:HD12	2.05	0.95
1:AAA:280:MET:CE	5:BBB:1118:HOH:O	2.23	0.86
1:AAA:280:MET:HE3	5:BBB:1118:HOH:O	1.80	0.80
1:AAA:29[B]:ASN:OD1	5:AAA:701:HOH:O	2.00	0.79
1:BBB:285[B]:MET:SD	1:BBB:414:LEU:CD1	2.72	0.78
1:BBB:119[A]:TRP:CE2	4:BBB:603:A1IDI:C8	2.68	0.77
1:BBB:103:PHE:CE1	1:BBB:199:ILE:HD11	2.21	0.76
1:BBB:451:LEU:HA	1:BBB:454:MET:HE2	1.67	0.74
1:BBB:28:LEU:HD11	1:BBB:454:MET:HE1	1.68	0.74
1:AAA:117:ASN:HD22	1:AAA:120:ARG:HH21	1.35	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:117:ASN:HD22	1:BBB:120:ARG:HH21	1.35	0.72
1:BBB:119[A]:TRP:CZ2	4:BBB:603:A1IDI:C1	2.72	0.71
1:AAA:251:ASN:HD22	1:AAA:251:ASN:H	1.39	0.71
1:BBB:451:LEU:HD23	1:BBB:454:MET:HE1	1.75	0.68
1:BBB:103:PHE:CD1	4:BBB:603:A1IDI:C11	2.77	0.67
1:BBB:451:LEU:HD23	1:BBB:454:MET:CE	2.24	0.67
1:AAA:103:PHE:HZ	5:AAA:888:HOH:O	1.77	0.67
1:AAA:103:PHE:CZ	5:AAA:888:HOH:O	2.48	0.65
1:AAA:268:LEU:HD23	1:BBB:270[A]:MET:HE1	1.78	0.65
1:BBB:251:ASN:HD22	1:BBB:251:ASN:H	1.45	0.64
1:BBB:118:PHE:CE1	1:BBB:152:LEU:CD2	2.80	0.63
1:AAA:111:THR:HG23	1:AAA:160[B]:SER:OG	1.98	0.63
1:BBB:103:PHE:CD1	4:BBB:603:A1IDI:C8	2.82	0.61
1:BBB:119[A]:TRP:CG	4:BBB:603:A1IDI:C10	2.84	0.61
1:AAA:270[A]:MET:HE1	1:BBB:268:LEU:HD23	1.83	0.60
1:BBB:119[A]:TRP:CD1	4:BBB:603:A1IDI:C11	2.85	0.60
1:BBB:118:PHE:CE1	1:BBB:152:LEU:HD22	2.38	0.58
1:BBB:280:MET:HE3	1:BBB:281:MET:HG2	1.86	0.58
1:BBB:118:PHE:HE1	1:BBB:152:LEU:CD2	2.17	0.57
1:BBB:451:LEU:HA	1:BBB:454:MET:CE	2.35	0.57
1:BBB:119[A]:TRP:CD2	4:BBB:603:A1IDI:C10	2.86	0.57
1:BBB:103:PHE:HD1	4:BBB:603:A1IDI:C8	2.19	0.56
1:BBB:103:PHE:HZ	5:BBB:801:HOH:O	1.89	0.56
1:BBB:103:PHE:CZ	5:BBB:801:HOH:O	2.53	0.54
1:BBB:321:GLU:H	1:BBB:321:GLU:CD	2.16	0.54
1:AAA:163:GLN:HB3	1:AAA:318:ASP:OD2	2.08	0.54
1:BBB:119[A]:TRP:CZ2	4:BBB:603:A1IDI:O1	2.61	0.54
1:BBB:118:PHE:HD2	1:BBB:119[B]:TRP:CD2	2.25	0.54
1:AAA:280:MET:HE1	1:BBB:387:ASN:ND2	2.24	0.53
1:AAA:280:MET:HE1	5:BBB:1118:HOH:O	2.00	0.51
1:BBB:318:ASP:CG	5:BBB:727:HOH:O	2.53	0.51
1:AAA:280:MET:CE	1:BBB:387:ASN:ND2	2.75	0.49
1:BBB:93:LYS:HE3	5:BBB:814:HOH:O	2.11	0.49
1:AAA:111:THR:CG2	1:AAA:160[B]:SER:OG	2.60	0.48
1:AAA:108:ASN:OD1	1:AAA:108:ASN:C	2.57	0.48
1:BBB:309:LYS:HG3	5:BBB:915:HOH:O	2.13	0.47
1:BBB:103:PHE:CE1	1:BBB:199:ILE:CD1	2.95	0.47
1:BBB:119[A]:TRP:CG	4:BBB:603:A1IDI:C11	2.98	0.47
1:BBB:410:TYR:O	1:BBB:413:VAL:HG13	2.15	0.47
1:AAA:58:GLY:HA2	2:AAA:601:FAD:C4X	2.46	0.46
1:BBB:117:ASN:HD22	1:BBB:120:ARG:NH2	2.10	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:119[A]:TRP:CZ2	4:BBB:603:A1IDI:C8	2.99	0.45
1:AAA:28:LEU:HD21	1:AAA:456:LYS:HE3	1.98	0.45
1:BBB:122[A]:MET:HE3	1:BBB:122[A]:MET:HB3	1.63	0.45
1:AAA:22:LEU:C	1:AAA:22:LEU:HD23	2.43	0.44
1:BBB:285[B]:MET:SD	1:BBB:413:VAL:HG22	2.58	0.43
1:AAA:117:ASN:HD22	1:AAA:120:ARG:NH2	2.09	0.43
1:AAA:270[B]:MET:SD	1:BBB:270[B]:MET:SD	3.16	0.43
1:AAA:122[A]:MET:HE3	1:AAA:122[A]:MET:HB3	1.89	0.42
1:AAA:233:ARG:HG3	1:AAA:251:ASN:HD21	1.84	0.42
1:BBB:118:PHE:HD2	1:BBB:119[B]:TRP:CE3	2.36	0.42
1:BBB:119[A]:TRP:CH2	4:BBB:603:A1IDI:O1	2.72	0.42
1:AAA:389:CYS:CB	1:BBB:280:MET:HG3	2.50	0.42
1:BBB:93:LYS:HA	1:BBB:93:LYS:CE	2.49	0.42
1:AAA:410:TYR:OH	1:BBB:350:ARG:HD2	2.20	0.42
1:AAA:270[A]:MET:HE1	1:BBB:268:LEU:CD2	2.50	0.41
1:BBB:363:LYS:HE2	5:BBB:1128:HOH:O	2.19	0.41
1:AAA:117:ASN:ND2	1:AAA:120:ARG:HH21	2.12	0.41
1:AAA:280:MET:HE3	1:AAA:280:MET:HB3	1.67	0.41
1:BBB:103:PHE:CE1	4:BBB:603:A1IDI:C11	3.01	0.41
1:BBB:146[B]:MET:HG2	1:BBB:150:GLU:HB2	2.03	0.41
1:BBB:254:MET:HE2	1:BBB:254:MET:HB2	1.75	0.41
1:AAA:23:LEU:HB2	1:AAA:30:VAL:HG11	2.03	0.40
1:BBB:119[A]:TRP:CD2	4:BBB:603:A1IDI:C8	3.05	0.40
1:AAA:118:PHE:CE1	1:AAA:152:LEU:CD2	3.04	0.40

All (2) 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:AAA:475:GLN:NE2	1:AAA:475:GLN:NE2[3_656]	1.87	0.33
1:AAA:93:LYS:NZ	1:AAA:159:GLU:OE1[6_565]	2.13	0.07

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	504/520 (97%)	492 (98%)	12 (2%)	0	100	100
1	BBB	500/520 (96%)	488 (98%)	12 (2%)	0	100	100
All	All	1004/1040 (96%)	980 (98%)	24 (2%)	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	435/444 (98%)	431 (99%)	4 (1%)	70	44
1	BBB	432/444 (97%)	430 (100%)	2 (0%)	81	61
All	All	867/888 (98%)	861 (99%)	6 (1%)	76	52

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	190	LYS
1	AAA	251	ASN
1	AAA	280	MET
1	AAA	350	ARG
1	BBB	104	PRO
1	BBB	251	ASN

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 [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](#)

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$
4	A1IDI	AAA	603	2	23,25,36	3.28	10 (43%)	31,34,50	7.79	15 (48%)
2	FAD	AAA	601	1,4	58,58,58	1.06	4 (6%)	85,89,89	1.51	12 (14%)
4	A1IDI	BBB	603	2	35,36,36	2.53	13 (37%)	46,50,50	6.31	14 (30%)
2	FAD	BBB	601	1,4	58,58,58	1.25	5 (8%)	85,89,89	1.58	17 (20%)
3	C15	AAA	602	-	14,14,21	0.79	1 (7%)	19,19,26	0.87	1 (5%)
3	C15	BBB	602	-	10,10,21	0.85	1 (10%)	15,15,26	0.66	0

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	A1IDI	AAA	603	2	-	5/10/11/15	0/3/3/5
2	FAD	AAA	601	1,4	-	0/34/50/50	0/6/6/6
4	A1IDI	BBB	603	2	-	7/14/15/15	0/5/5/5
2	FAD	BBB	601	1,4	-	2/34/50/50	0/6/6/6
3	C15	AAA	602	-	-	2/14/14/21	-
3	C15	BBB	602	-	-	1/8/8/21	-

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	AAA	603	A1IDI	CBD-CBE	9.50	1.46	1.18
4	AAA	603	A1IDI	NAQ-NAR	7.11	1.41	1.30
4	BBB	603	A1IDI	CBD-CBE	6.88	1.38	1.18
4	BBB	603	A1IDI	C2-N1	4.99	1.42	1.33
4	BBB	603	A1IDI	C2-N3	4.82	1.42	1.33
4	AAA	603	A1IDI	C2-N3	4.39	1.41	1.33
4	BBB	603	A1IDI	CBA-CAX	-4.18	1.43	1.51
2	AAA	601	FAD	C4X-N5	3.97	1.39	1.30
4	AAA	603	A1IDI	CBA-CAX	-3.94	1.44	1.51
2	BBB	601	FAD	C4X-N5	3.77	1.38	1.30
4	BBB	603	A1IDI	CBA-NBB	-3.74	1.38	1.47
4	BBB	603	A1IDI	C5-C6	-3.63	1.35	1.39
4	AAA	603	A1IDI	C2-N1	3.55	1.40	1.33
4	AAA	603	A1IDI	CBA-NBB	-3.37	1.39	1.47
4	BBB	603	A1IDI	C6-SAJ	-3.35	1.68	1.76
4	BBB	603	A1IDI	C4-NAS	-3.32	1.32	1.35
4	BBB	603	A1IDI	C13-C7	-3.29	1.34	1.40
4	BBB	603	A1IDI	C7-N2	-3.23	1.33	1.39
4	BBB	603	A1IDI	C8-C7	-3.07	1.34	1.38
2	BBB	601	FAD	C6-C7	-3.05	1.35	1.39
4	BBB	603	A1IDI	C5-C4	-2.99	1.33	1.39
4	AAA	603	A1IDI	CAT-NAS	2.73	1.49	1.46
4	AAA	603	A1IDI	NAS-NAR	-2.70	1.33	1.36
3	AAA	602	C15	O3S-S1	2.68	1.57	1.47
2	BBB	601	FAD	C4X-C10	-2.63	1.36	1.44
2	AAA	601	FAD	C9-C9A	-2.63	1.35	1.39
2	AAA	601	FAD	C4X-C10	-2.56	1.36	1.44
4	AAA	603	A1IDI	C5-C6	-2.48	1.36	1.39
2	AAA	601	FAD	C8A-N7A	2.47	1.36	1.31
4	AAA	603	A1IDI	CAT-CAU	-2.39	1.47	1.51
2	BBB	601	FAD	P-O3P	2.38	1.62	1.59
3	BBB	602	C15	O3S-S1	2.36	1.56	1.47
2	BBB	601	FAD	C4X-C4	-2.29	1.36	1.44
4	BBB	603	A1IDI	C10-C8	-2.14	1.35	1.39

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	AAA	603	A1IDI	CBC-CBD-CBE	-39.97	112.80	177.63
4	BBB	603	A1IDI	CBC-CBD-CBE	-39.75	113.14	177.63
4	AAA	603	A1IDI	C5-C4-NAS	8.81	111.25	104.25
4	BBB	603	A1IDI	CAT-NAS-C4	-6.55	123.66	127.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	601	FAD	C9A-C5X-N5	-5.66	116.45	122.45
4	AAA	603	A1IDI	N3-C2-N1	-5.49	120.27	128.58
2	AAA	601	FAD	C9A-C5X-N5	-5.31	116.82	122.45
4	AAA	603	A1IDI	C4-C5-C6	5.22	121.50	116.44
4	BBB	603	A1IDI	C5-C4-NAS	5.21	108.38	104.25
4	BBB	603	A1IDI	N3-C2-N1	-5.19	120.72	128.58
4	AAA	603	A1IDI	CBC-NBB-CBA	4.91	119.06	111.42
4	AAA	603	A1IDI	C5-C4-N3	-4.80	120.10	126.72
4	BBB	603	A1IDI	CAT-NAS-NAR	4.62	128.69	120.27
4	BBB	603	A1IDI	CBC-NBB-CBA	4.39	118.25	111.42
4	AAA	603	A1IDI	C4-C5-NAQ	-4.20	106.07	109.31
4	AAA	603	A1IDI	C4-NAS-NAR	-3.96	107.04	109.93
4	BBB	603	A1IDI	C4-C5-C6	3.90	120.22	116.44
4	BBB	603	A1IDI	CAT-CAU-CAV	3.71	127.58	120.75
2	AAA	601	FAD	C10-C4X-N5	-3.69	117.28	124.81
2	BBB	601	FAD	C4X-C10-N10	3.65	121.70	116.48
4	AAA	603	A1IDI	CAT-CAU-CAV	3.63	127.42	120.75
2	AAA	601	FAD	C4-C4X-N5	3.57	123.14	118.21
2	BBB	601	FAD	C10-C4X-N5	-3.53	117.61	124.81
4	BBB	603	A1IDI	C4-C5-NAQ	-3.26	106.79	109.31
4	BBB	603	A1IDI	C4-NAS-NAR	-3.18	107.61	109.93
2	BBB	601	FAD	C9A-C9-C8	3.12	125.49	119.22
2	BBB	601	FAD	C9A-N10-C10	-3.11	116.01	120.75
2	AAA	601	FAD	C6-C5X-C9A	3.10	123.31	119.05
2	AAA	601	FAD	C9-C9A-N10	2.95	125.82	121.85
2	BBB	601	FAD	C5A-C4A-N3A	-2.93	122.68	126.72
2	AAA	601	FAD	C4-N3-C2	-2.83	120.62	125.64
2	AAA	601	FAD	C4X-C10-N10	2.75	120.42	116.48
2	AAA	601	FAD	C5'-C4'-C3'	-2.73	107.06	112.22
2	BBB	601	FAD	C9-C9A-C5X	-2.67	115.31	120.03
4	AAA	603	A1IDI	CAT-NAS-NAR	2.57	124.95	120.27
2	BBB	601	FAD	C4-C4X-N5	2.53	121.70	118.21
2	AAA	601	FAD	C9-C9A-C5X	-2.50	115.61	120.03
4	AAA	603	A1IDI	C2-N3-C4	2.50	117.94	111.83
4	BBB	603	A1IDI	CAV-CAU-CAZ	-2.46	114.58	118.23
2	BBB	601	FAD	C6-C5X-N5	2.45	122.52	118.44
2	AAA	601	FAD	C5A-C4A-N9A	2.44	108.47	105.81
4	AAA	603	A1IDI	CAT-CAU-CAZ	-2.42	116.28	120.75
2	BBB	601	FAD	C9-C8-C7	-2.41	116.15	119.69
3	AAA	602	C15	O3S-S1-C1	-2.41	101.29	106.00
2	BBB	601	FAD	C9-C9A-N10	2.38	125.05	121.85
4	BBB	603	A1IDI	CAX-CBA-NBB	2.37	117.58	113.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	AAA	603	A1IDI	CBD-CBC-NBB	2.34	118.26	113.70
4	AAA	603	A1IDI	CAY-CAZ-CAU	2.31	124.03	121.00
2	AAA	601	FAD	C9A-C9-C8	2.28	123.81	119.22
4	BBB	603	A1IDI	C10-C8-C7	-2.27	120.86	123.62
2	BBB	601	FAD	O3'-C3'-C4'	-2.17	103.99	108.93
4	AAA	603	A1IDI	CBA-CAX-CAY	-2.17	116.75	120.75
2	BBB	601	FAD	C2B-C3B-C4B	2.15	106.77	102.61
4	BBB	603	A1IDI	C12-C11-C10	-2.14	117.60	120.24
2	BBB	601	FAD	C4X-C10-N1	-2.12	119.39	124.59
2	BBB	601	FAD	C6A-C5A-C4A	2.10	120.05	117.18
2	BBB	601	FAD	C7M-C7-C6	-2.07	115.93	119.57
2	AAA	601	FAD	C8M-C8-C7	2.07	124.98	120.76
2	BBB	601	FAD	C4-C4X-C10	2.04	120.43	116.93

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	BBB	602	C15	S1-C1-C2-C3
4	AAA	603	A1IDI	CAU-CAT-NAS-NAR
4	AAA	603	A1IDI	CAX-CBA-NBB-CBC
4	AAA	603	A1IDI	CBD-CBC-NBB-CBF
4	BBB	603	A1IDI	N1-C6-SAJ-C1
4	BBB	603	A1IDI	C5-C6-SAJ-C1
4	BBB	603	A1IDI	CAU-CAT-NAS-NAR
4	BBB	603	A1IDI	CAX-CBA-NBB-CBC
4	BBB	603	A1IDI	CBD-CBC-NBB-CBF
3	AAA	602	C15	S1-C1-C2-C3
4	AAA	603	A1IDI	CBD-CBC-NBB-CBA
4	BBB	603	A1IDI	CBD-CBC-NBB-CBA
4	AAA	603	A1IDI	CAU-CAT-NAS-C4
4	BBB	603	A1IDI	CAU-CAT-NAS-C4
3	AAA	602	C15	C12-C13-C14-C15
2	BBB	601	FAD	C2'-C1'-N10-C10
2	BBB	601	FAD	O4B-C4B-C5B-O5B

There are no ring outliers.

2 monomers are involved in 18 short contacts:

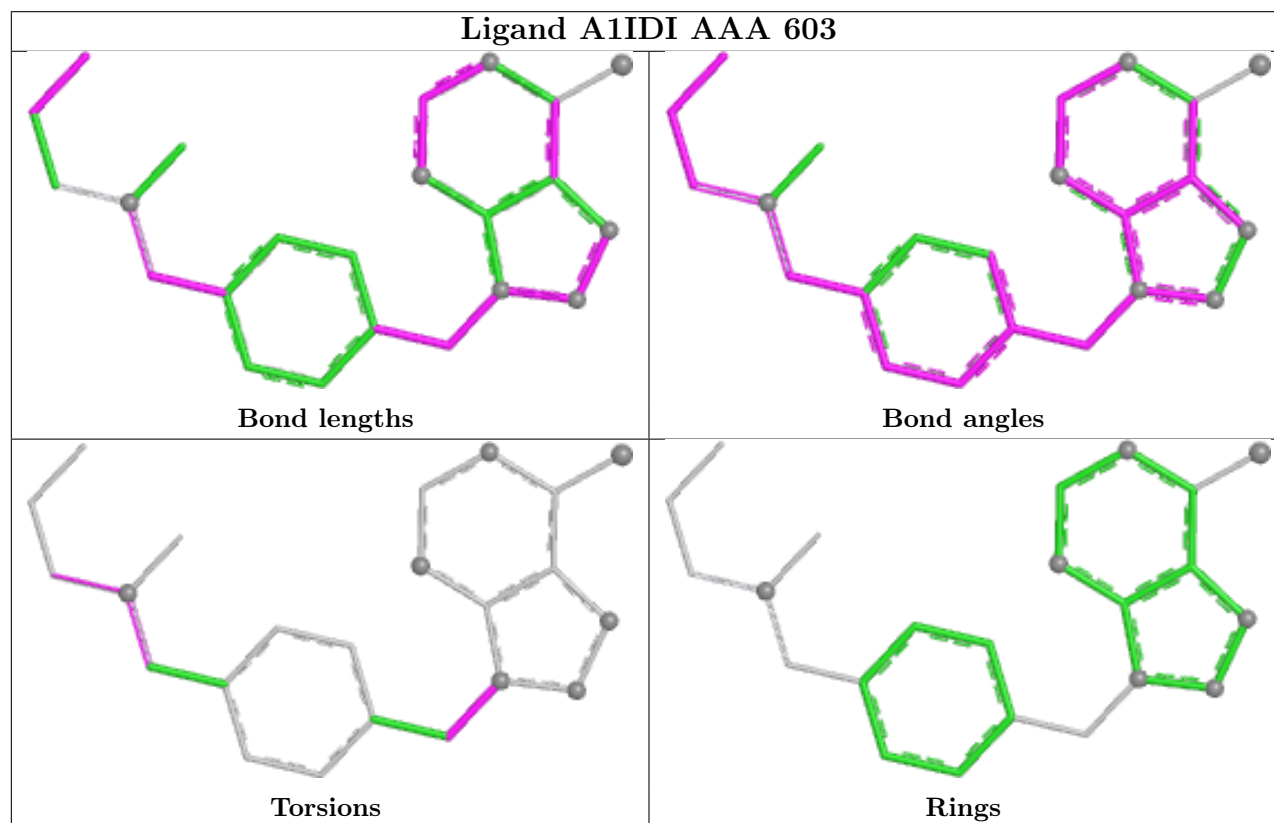
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	AAA	601	FAD	1	0

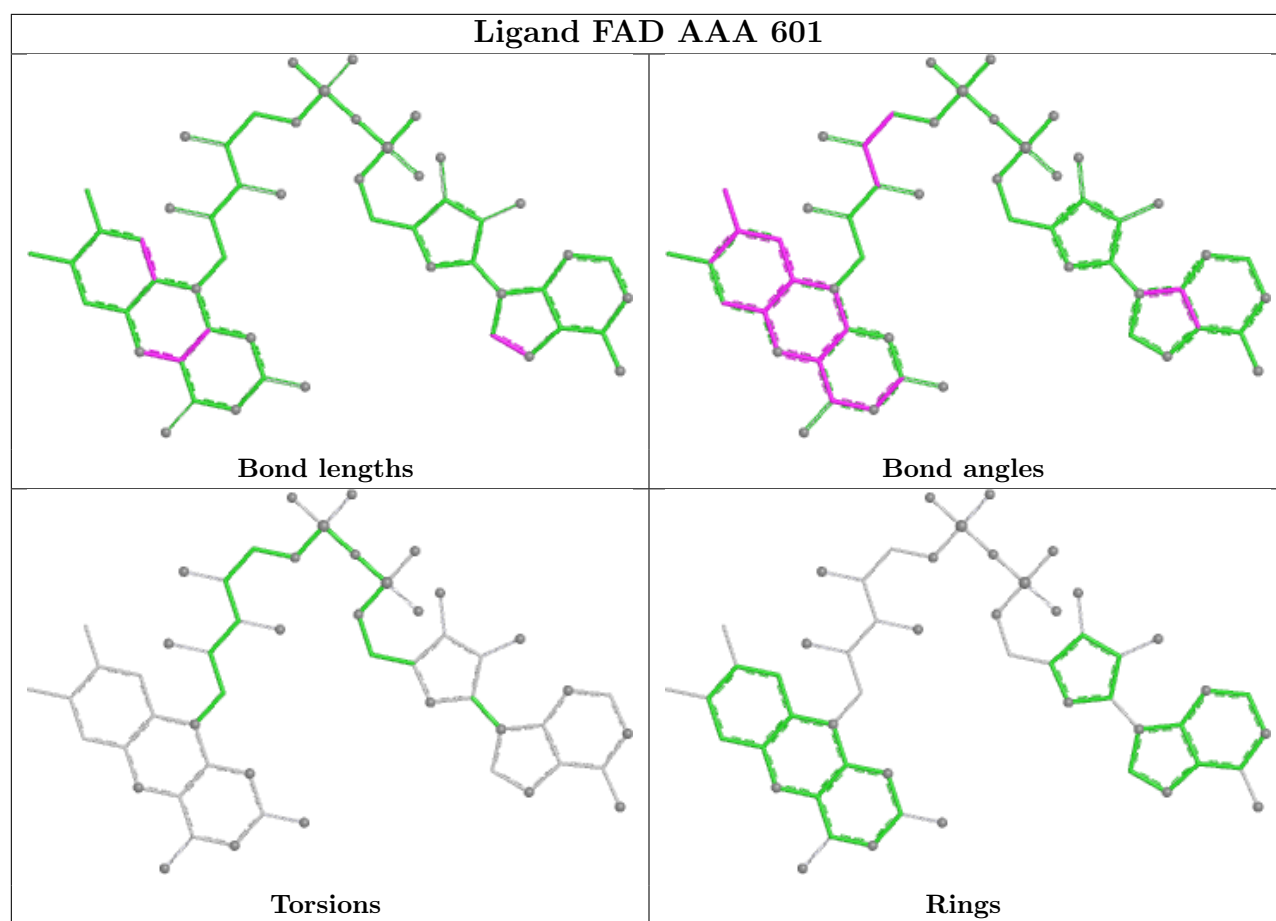
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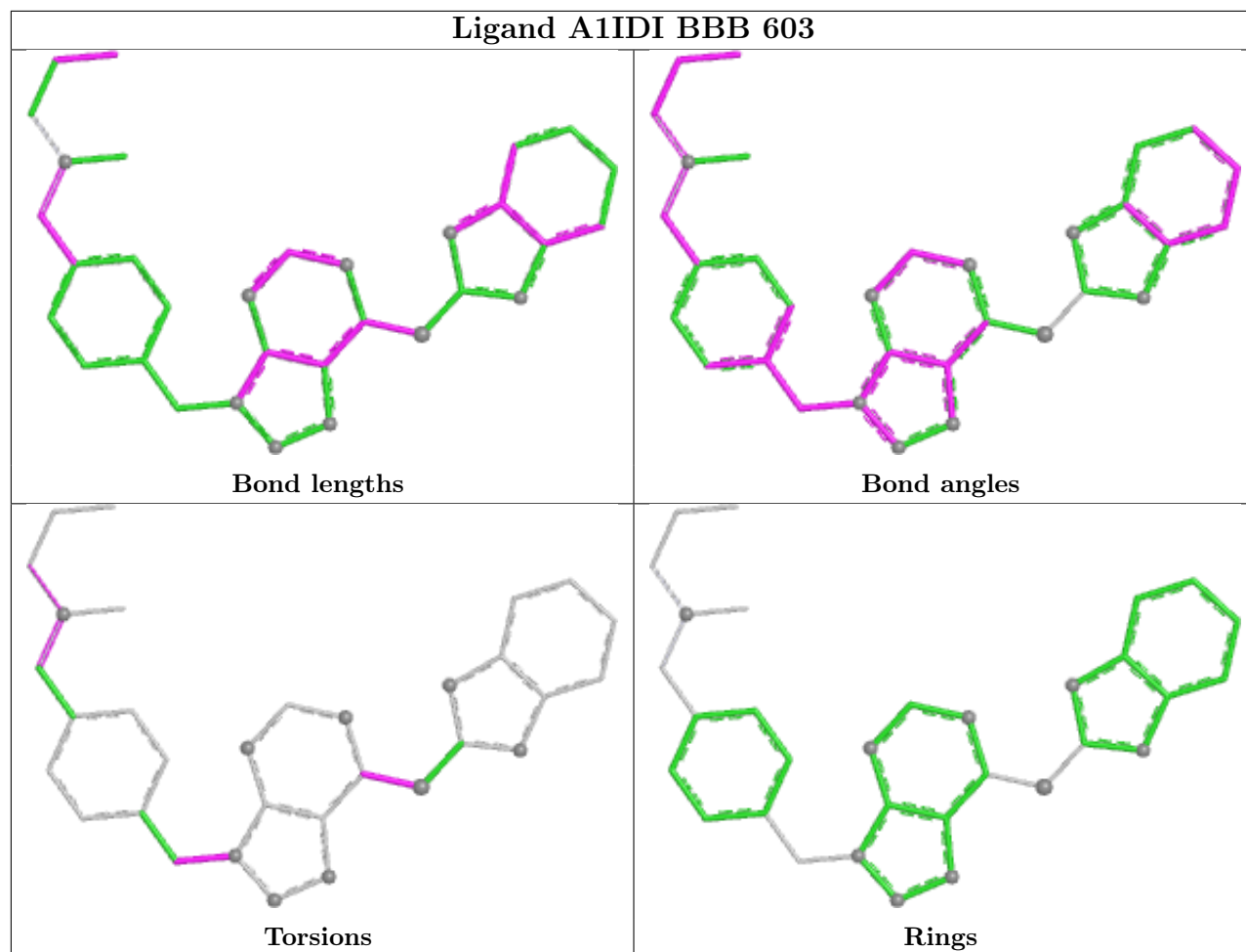
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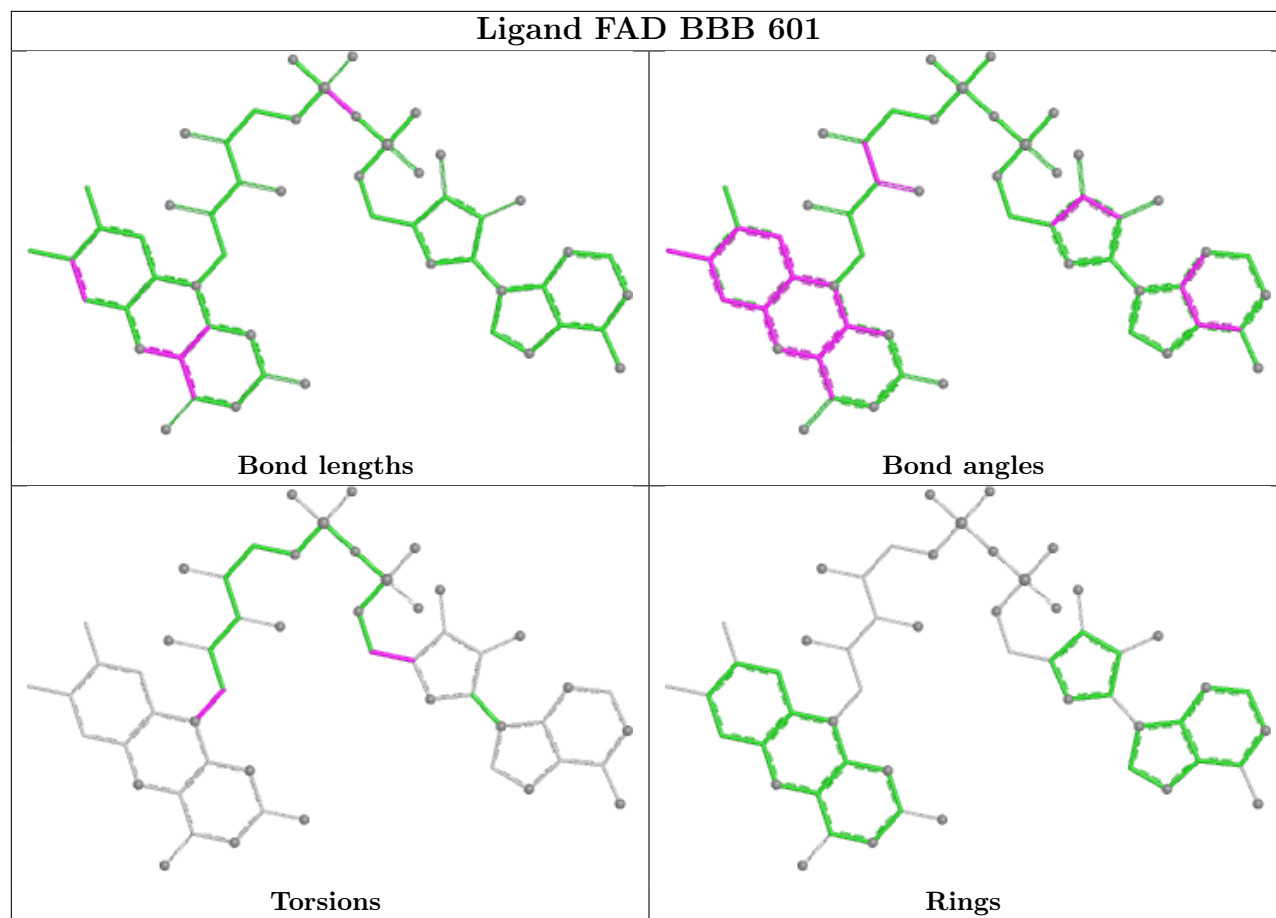
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	BBB	603	A1IDI	17	0

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.

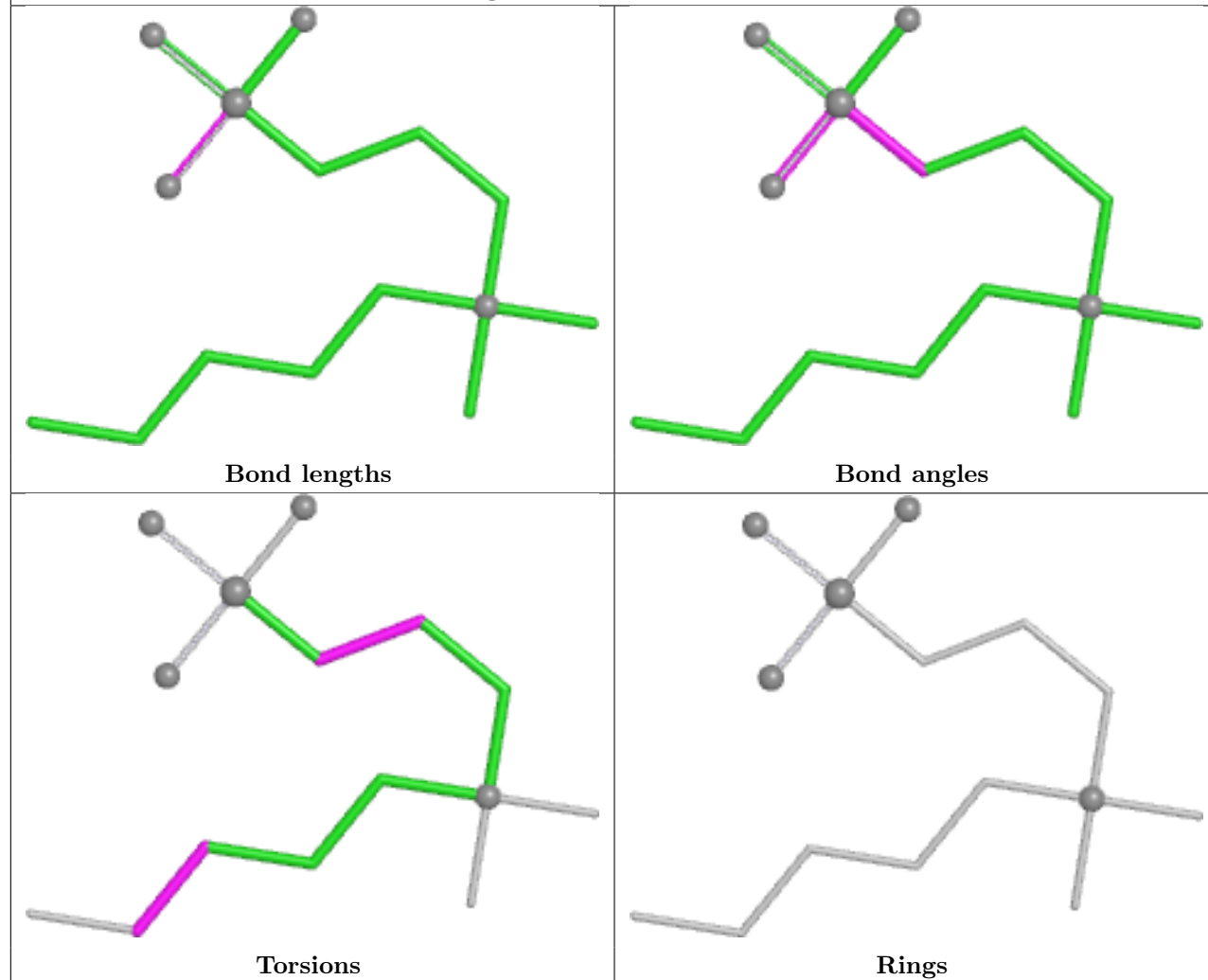




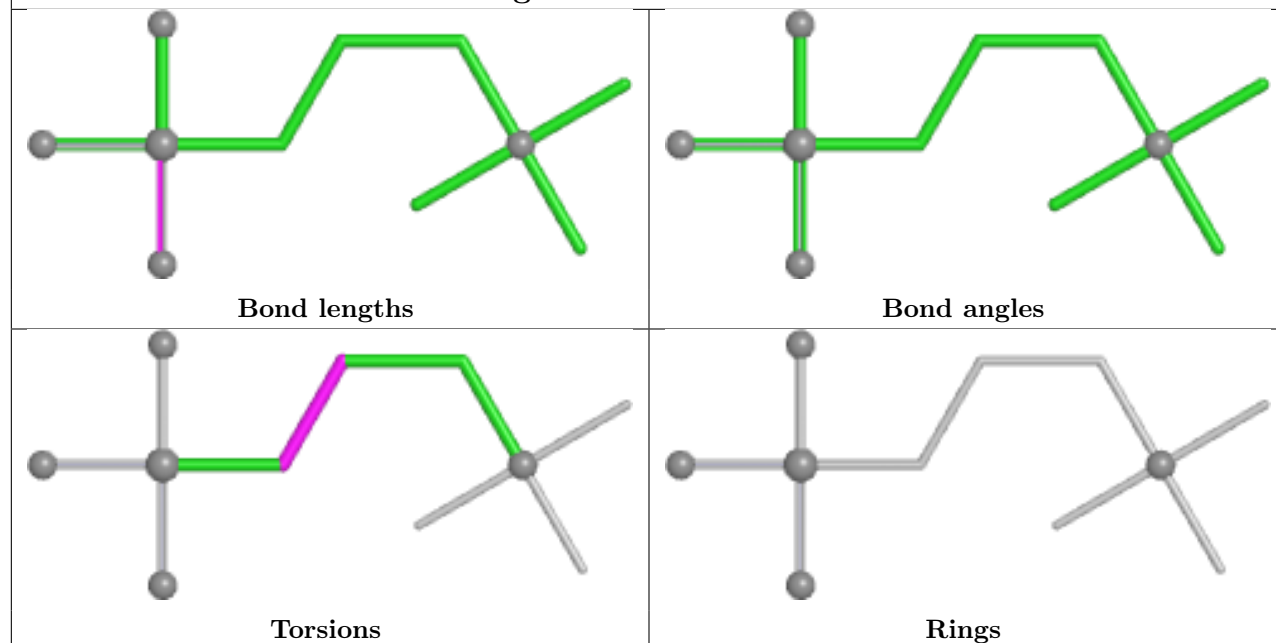




Ligand C15 AAA 602



Ligand C15 BBB 602



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	AAA	498/520 (95%)	-0.02	26 (5%)	33 35	6, 13, 29, 63	8 (1%)
1	BBB	495/520 (95%)	-0.16	18 (3%)	46 48	6, 11, 27, 66	7 (1%)
All	All	993/1040 (95%)	-0.09	44 (4%)	39 40	6, 12, 28, 66	15 (1%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	495	LEU	6.8
1	BBB	496	ILE	6.3
1	AAA	498	LEU	5.8
1	AAA	499	THR	5.3
1	BBB	103	PHE	5.0
1	BBB	494	ARG	4.6
1	AAA	107	TRP	4.6
1	BBB	107	TRP	4.6
1	AAA	252	HIS	4.2
1	AAA	103	PHE	4.0
1	BBB	493	LEU	4.0
1	AAA	495	LEU	3.8
1	BBB	119[A]	TRP	3.4
1	AAA	481	PHE	3.1
1	AAA	321	GLU	3.1
1	AAA	497	GLY	3.0
1	AAA	494	ARG	2.9
1	BBB	163	GLN	2.9
1	BBB	252	HIS	2.9
1	BBB	2	SER	2.8
1	BBB	303	GLU	2.8
1	AAA	482	LEU	2.8
1	BBB	242	ARG	2.6
1	AAA	496	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	AAA	354	ARG	2.6
1	AAA	106	VAL	2.5
1	AAA	119	TRP	2.5
1	AAA	105	PRO	2.5
1	BBB	492	LEU	2.4
1	BBB	354	ARG	2.4
1	AAA	93	LYS	2.4
1	AAA	251	ASN	2.4
1	AAA	319	GLY	2.3
1	BBB	481	PHE	2.3
1	AAA	473	PRO	2.3
1	BBB	280	MET	2.2
1	AAA	2	SER	2.2
1	AAA	3	ASN	2.2
1	AAA	470	VAL	2.2
1	AAA	4	LYS	2.2
1	AAA	493	LEU	2.1
1	AAA	244	ASN	2.1
1	BBB	106	VAL	2.1
1	BBB	112	TYR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	C15	AAA	602	15/22	0.81	0.16	22,30,52,55	0
3	C15	BBB	602	11/22	0.82	0.15	24,33,59,66	0
4	A1IDI	AAA	603	23/32	0.88	0.12	19,24,33,42	0

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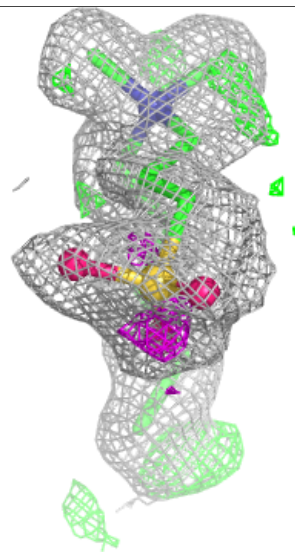
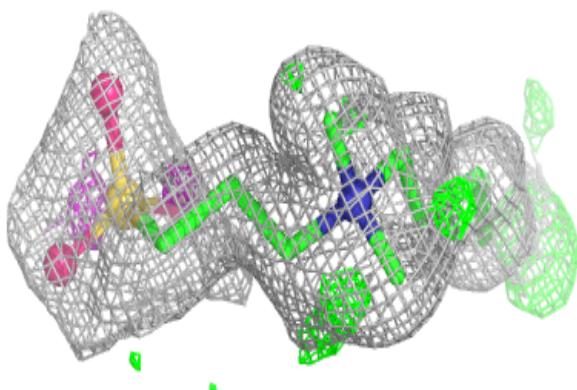
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	A1IDI	BBB	603	32/32	0.91	0.12	17,23,48,53	0
2	FAD	AAA	601	53/53	0.99	0.03	7,9,10,10	0
2	FAD	BBB	601	53/53	0.99	0.03	7,8,9,10	0

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.

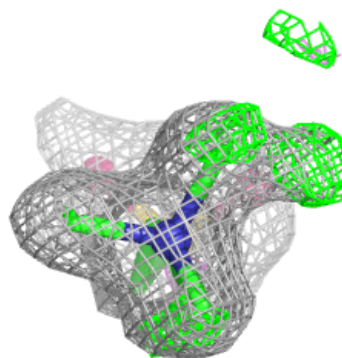
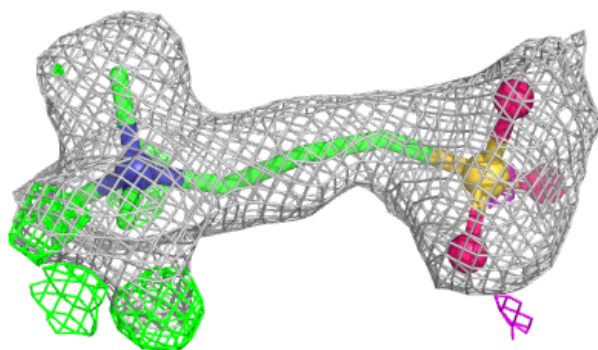
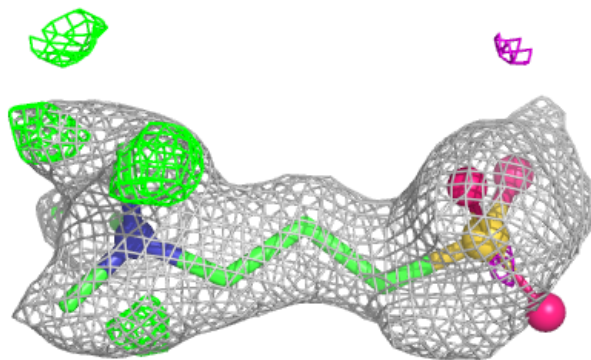
Electron density around C15 AAA 602:

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

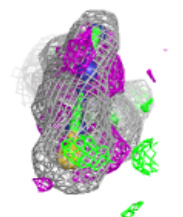
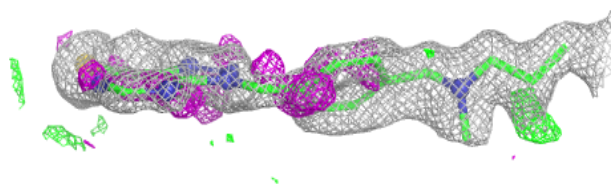
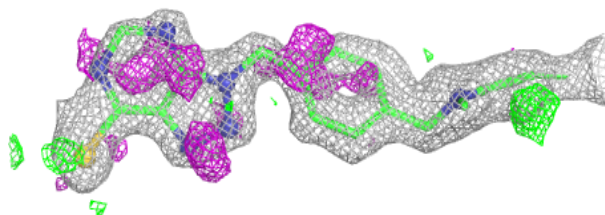


Electron density around C15 BBB 602:

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

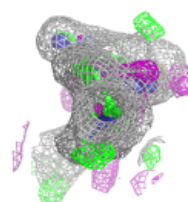
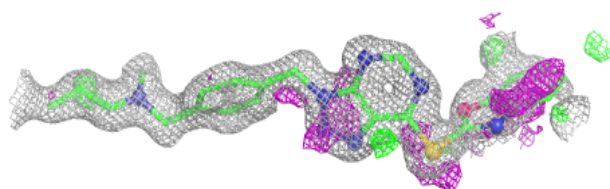
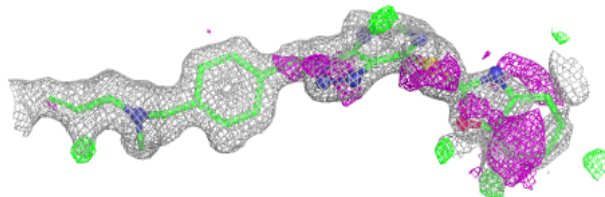
**Electron density around A1IDI AAA 603:**

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

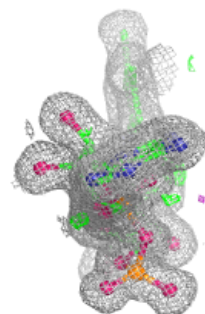
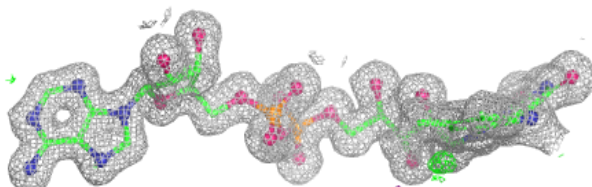
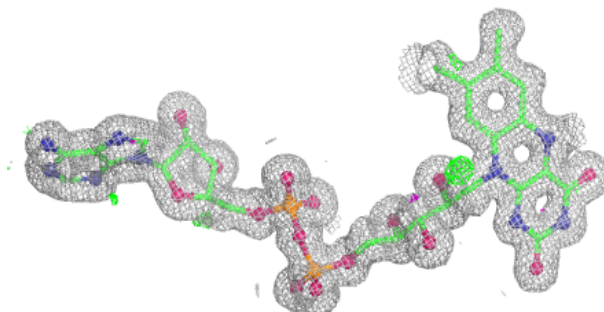


Electron density around A1IDI BBB 603:

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

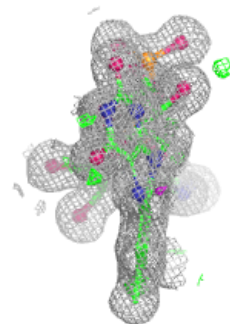
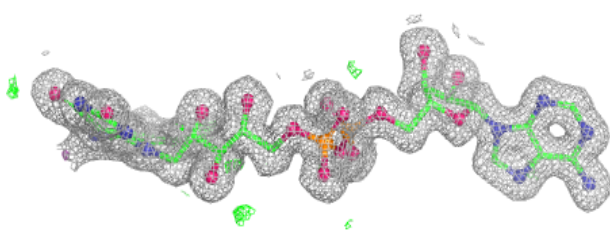
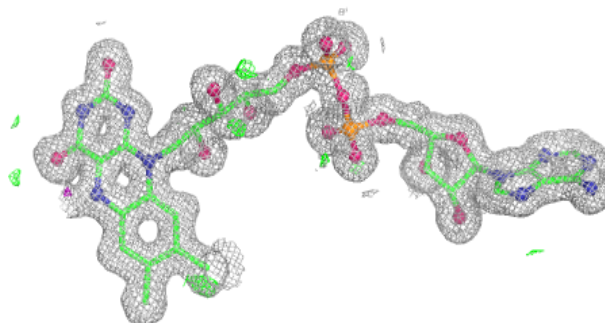
**Electron density around FAD AAA 601:**

$2mF_o-DF_c$ (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.