



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 18, 2020 – 10:09 AM JST

PDB ID : 7CU1
Title : CRYSTAL STRUCTURE OF STREPTOMYCES ALBOGRISEOLUS
FLAVIN-DEPENDENT TRYPTOPHAN 6-HALOGENASE (THAL) IN
COMPLEX WITH FAD and AMP
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Deposited on : 2020-08-20
Resolution : 1.91 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 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 : 2.14.6
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) : 2.14.6

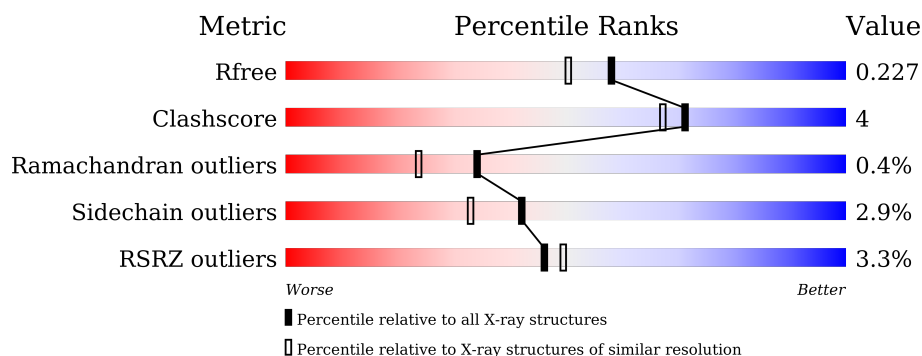
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.91 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 on 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	551	<div> <div>3%</div> <div>88%</div> <div>8%</div> <div>• •</div> </div>
1	B	551	<div> <div>3%</div> <div>84%</div> <div>10%</div> <div>• 5%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9245 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 Tryptophan 6-halogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	0	0	0
			4263	2723	735	784	21			
1	B	524	Total	C	N	O	S	0	0	0
			4175	2666	719	770	20			

There are 40 discrepancies between the modelled and reference sequences:

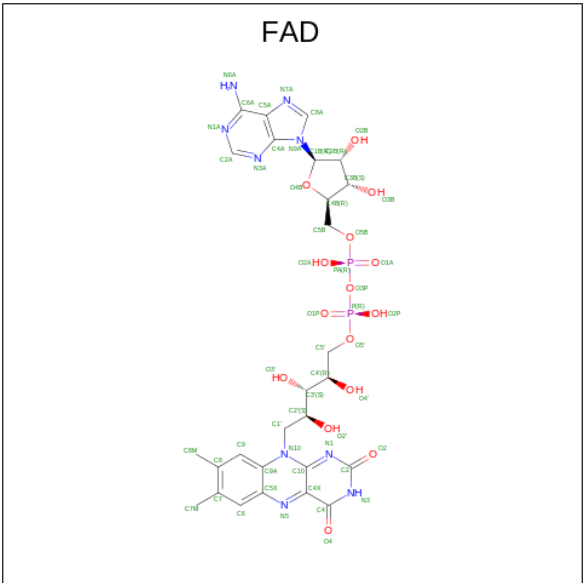
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP A1E280
A	-18	GLY	-	expression tag	UNP A1E280
A	-17	SER	-	expression tag	UNP A1E280
A	-16	SER	-	expression tag	UNP A1E280
A	-15	HIS	-	expression tag	UNP A1E280
A	-14	HIS	-	expression tag	UNP A1E280
A	-13	HIS	-	expression tag	UNP A1E280
A	-12	HIS	-	expression tag	UNP A1E280
A	-11	HIS	-	expression tag	UNP A1E280
A	-10	HIS	-	expression tag	UNP A1E280
A	-9	SER	-	expression tag	UNP A1E280
A	-8	SER	-	expression tag	UNP A1E280
A	-7	GLY	-	expression tag	UNP A1E280
A	-6	LEU	-	expression tag	UNP A1E280
A	-5	VAL	-	expression tag	UNP A1E280
A	-4	PRO	-	expression tag	UNP A1E280
A	-3	ARG	-	expression tag	UNP A1E280
A	-2	GLY	-	expression tag	UNP A1E280
A	-1	SER	-	expression tag	UNP A1E280
A	0	HIS	-	expression tag	UNP A1E280
B	-19	MET	-	expression tag	UNP A1E280
B	-18	GLY	-	expression tag	UNP A1E280
B	-17	SER	-	expression tag	UNP A1E280
B	-16	SER	-	expression tag	UNP A1E280
B	-15	HIS	-	expression tag	UNP A1E280

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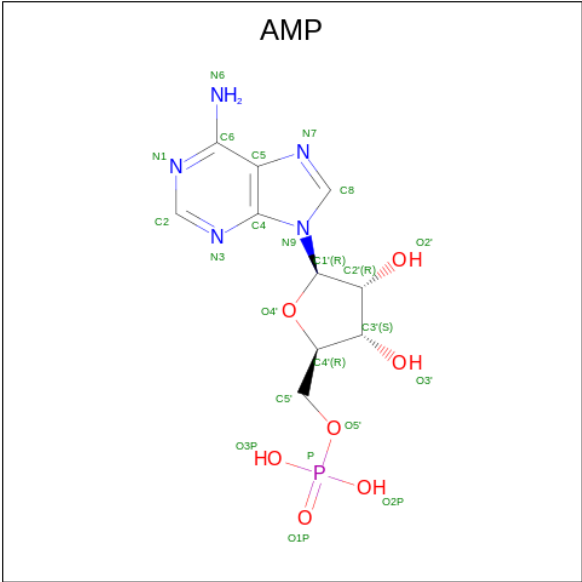
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP A1E280
B	-13	HIS	-	expression tag	UNP A1E280
B	-12	HIS	-	expression tag	UNP A1E280
B	-11	HIS	-	expression tag	UNP A1E280
B	-10	HIS	-	expression tag	UNP A1E280
B	-9	SER	-	expression tag	UNP A1E280
B	-8	SER	-	expression tag	UNP A1E280
B	-7	GLY	-	expression tag	UNP A1E280
B	-6	LEU	-	expression tag	UNP A1E280
B	-5	VAL	-	expression tag	UNP A1E280
B	-4	PRO	-	expression tag	UNP A1E280
B	-3	ARG	-	expression tag	UNP A1E280
B	-2	GLY	-	expression tag	UNP A1E280
B	-1	SER	-	expression tag	UNP A1E280
B	0	HIS	-	expression tag	UNP A1E280

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



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

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

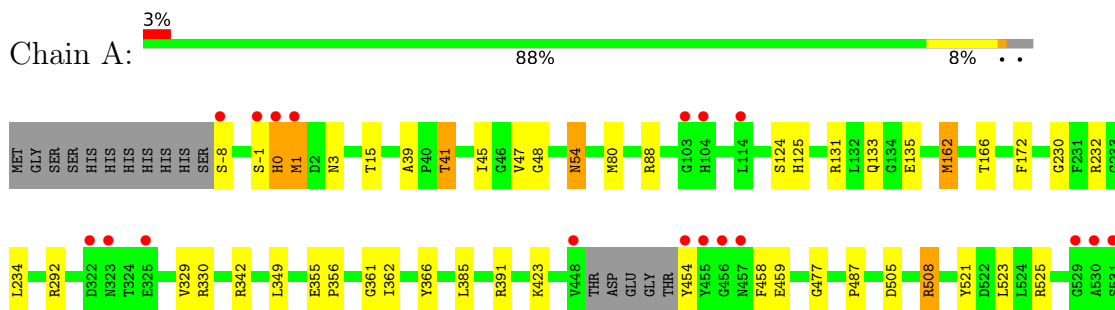
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	326	Total	O	0	0
			326	326		
4	B	405	Total	O	0	0
			405	405		

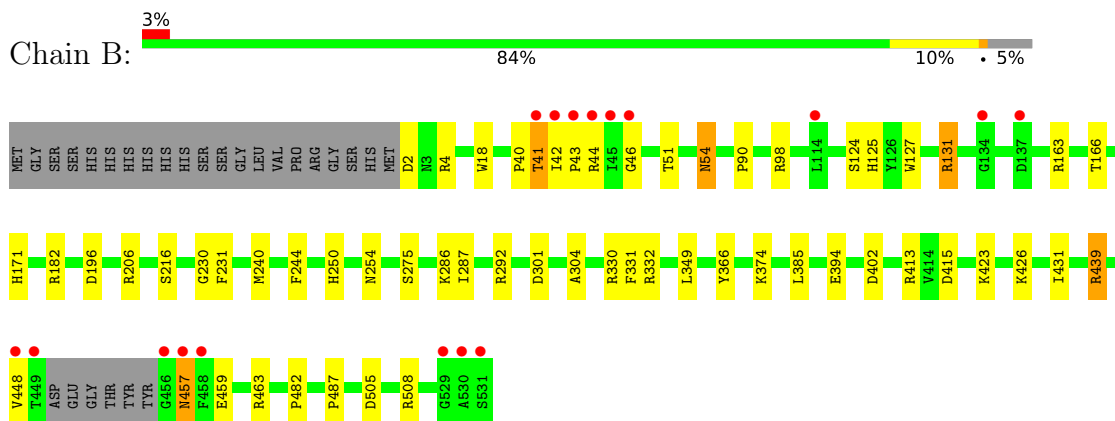
3 Residue-property plots [i](#)

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: Tryptophan 6-halogenase



• Molecule 1: Tryptophan 6-halogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.55Å 118.15Å 82.86Å 90.00° 100.88° 90.00°	Depositor
Resolution (Å)	20.64 – 1.91 20.64 – 1.91	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.64-1.91) 99.1 (20.64-1.91)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.174 , 0.220 0.183 , 0.227	Depositor DCC
R_{free} test set	3807 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.326	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9245	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AMP, FAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.50	0/4379	0.68	0/5942
1	B	0.53	1/4287 (0.0%)	0.70	2/5818 (0.0%)
All	All	0.51	1/8666 (0.0%)	0.69	2/11760 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	394	GLU	CD-OE2	-5.39	1.19	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	331	PHE	CB-CA-C	-5.68	99.03	110.40
1	B	413	ARG	NE-CZ-NH1	-5.13	117.74	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	206	ARG	Sidechain
1	B	439	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	457	ASN	Peptide
1	B	98	ARG	Sidechain

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	4263	0	4128	32	0
1	B	4175	0	4042	29	0
2	A	53	0	31	7	0
3	B	23	0	12	0	0
4	A	326	0	0	3	1
4	B	405	0	0	5	1
All	All	9245	0	8213	59	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 4.

All (59) 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:45:ILE:HG22	1:A:47:VAL:HG22	1.68	0.73
1:A:162:MET:HE2	1:A:162:MET:HA	1.73	0.69
1:A:342:ARG:NH2	4:A:701:HOH:O	2.26	0.68
1:B:505:ASP:OD1	1:B:508:ARG:NH2	2.28	0.66
1:B:127:TRP:CZ2	1:B:131:ARG:HG3	2.32	0.65
1:B:231:PHE:CZ	1:B:349:LEU:HD13	2.33	0.64
1:A:162:MET:CE	1:A:523:LEU:HD11	2.28	0.64
1:B:54:ASN:ND2	1:B:366:TYR:OH	2.32	0.63
1:B:2:ASP:OD1	1:B:4:ARG:HD3	2.00	0.62
1:A:124:SER:OG	1:A:125:HIS:HD2	1.83	0.61
1:A:0:HIS:HA	1:A:1:MET:O	2.01	0.59
1:A:162:MET:HE3	1:A:523:LEU:HD11	1.84	0.59
1:A:162:MET:CE	1:A:162:MET:HA	2.32	0.59
1:B:374:LYS:NZ	1:B:459:GLU:OE2	2.28	0.58
1:A:133:GLN:NE2	1:A:135:GLU:OE1	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:ARG:CZ	4:A:701:HOH:O	2.51	0.58
1:A:385:LEU:HD11	1:B:487:PRO:HG2	1.87	0.56
1:A:361:GLY:HA3	2:A:601:FAD:H1'2	1.88	0.56
1:B:124:SER:OG	1:B:125:HIS:HD2	1.88	0.56
1:A:232:ARG:HG3	1:A:234:LEU:HG	1.87	0.55
1:B:42:ILE:N	1:B:43:PRO:CD	2.70	0.55
1:B:40:PRO:O	1:B:41:THR:C	2.46	0.54
1:A:15:THR:HG21	4:A:951:HOH:O	2.09	0.53
1:A:47:VAL:HG21	2:A:601:FAD:C7	2.38	0.53
1:B:402:ASP:HB3	1:B:431:ILE:HG12	1.90	0.52
1:A:505:ASP:OD1	1:A:508:ARG:NH2	2.42	0.52
1:A:54:ASN:ND2	1:A:366:TYR:OH	2.43	0.52
1:A:362:ILE:HG12	2:A:601:FAD:O2	2.08	0.52
1:B:250:HIS:HD2	4:B:715:HOH:O	1.94	0.51
1:A:45:ILE:HG13	1:A:349:LEU:HD12	1.93	0.50
1:B:254:ASN:O	1:B:330:ARG:HD2	2.12	0.49
1:A:39:ALA:HB1	1:A:41:THR:HG22	1.96	0.48
1:B:244:PHE:CE2	1:B:332:ARG:HG2	2.49	0.48
1:B:196:ASP:OD2	1:B:216:SER:HB3	2.15	0.47
1:B:301:ASP:HA	1:B:304:ALA:O	2.15	0.47
1:A:487:PRO:HG2	1:B:385:LEU:HD21	1.96	0.47
1:A:356:PRO:HB3	2:A:601:FAD:C5X	2.45	0.47
1:A:45:ILE:CG2	1:A:47:VAL:HG22	2.41	0.46
1:B:171:HIS:CG	1:B:287:ILE:HG23	2.50	0.46
1:B:2:ASP:OD1	1:B:4:ARG:CD	2.64	0.46
1:B:51:THR:HB	4:B:928:HOH:O	2.16	0.45
1:A:88:ARG:HD3	1:A:477:GLY:O	2.16	0.45
1:B:463:ARG:NH2	4:B:706:HOH:O	2.42	0.44
1:A:-1:SER:O	1:A:0:HIS:C	2.55	0.44
1:B:439:ARG:HA	1:B:482:PRO:HA	1.98	0.44
1:B:275:SER:HB2	1:B:286:LYS:HB3	2.00	0.44
1:A:48:GLY:HA2	2:A:601:FAD:C4X	2.47	0.44
1:A:15:THR:OG1	2:A:601:FAD:H5'2	2.18	0.44
1:B:90:PRO:HB3	1:B:415:ASP:HB3	1.99	0.43
1:A:356:PRO:HG3	2:A:601:FAD:C6	2.49	0.42
1:B:230:GLY:HA2	1:B:349:LEU:HB2	2.01	0.42
1:A:0:HIS:HA	1:A:1:MET:C	2.39	0.42
1:B:42:ILE:N	1:B:43:PRO:HD2	2.34	0.42
1:A:230:GLY:HA2	1:A:349:LEU:HB2	2.02	0.42
1:A:521:TYR:CZ	1:A:525:ARG:HD2	2.55	0.42
1:A:80:MET:CE	1:A:166:THR:HG21	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:THR:HB	4:B:755:HOH:O	2.20	0.41
1:B:18:TRP:CE3	1:B:182:ARG:HA	2.55	0.41
1:B:240:MET:HE2	4:B:736:HOH:O	2.20	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)
4:A:934:HOH:O	4:B:730:HOH:O[1_655]	2.17	0.03

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	531/551 (96%)	513 (97%)	16 (3%)	2 (0%)	34	24
1	B	520/551 (94%)	504 (97%)	14 (3%)	2 (0%)	34	24
All	All	1051/1102 (95%)	1017 (97%)	30 (3%)	4 (0%)	34	24

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	0	HIS
1	B	41	THR
1	B	46	GLY

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	448/462 (97%)	431 (96%)	17 (4%)	33	22
1	B	439/462 (95%)	430 (98%)	9 (2%)	53	46
All	All	887/924 (96%)	861 (97%)	26 (3%)	42	33

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-8	SER
1	A	3	ASN
1	A	41	THR
1	A	54	ASN
1	A	131	ARG
1	A	162	MET
1	A	172	PHE
1	A	292	ARG
1	A	329	VAL
1	A	330	ARG
1	A	355	GLU
1	A	391	ARG
1	A	423	LYS
1	A	454	TYR
1	A	458	PHE
1	A	459	GLU
1	A	508	ARG
1	B	44	ARG
1	B	54	ASN
1	B	131	ARG
1	B	163	ARG
1	B	292	ARG
1	B	423	LYS
1	B	426	LYS
1	B	448	VAL
1	B	457	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN

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Mol	Chain	Res	Type
1	A	125	HIS
1	A	189	GLN
1	A	221	GLN
1	A	457	ASN
1	A	494	GLN
1	B	3	ASN
1	B	54	ASN
1	B	125	HIS
1	B	221	GLN
1	B	250	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	AMP	B	601	-	22,25,25	1.13	2 (9%)	25,38,38	1.71	5 (20%)
2	FAD	A	601	-	51,58,58	2.12	8 (15%)	60,89,89	2.26	20 (33%)

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	AMP	B	601	-	-	3/6/26/26	0/3/3/3
2	FAD	A	601	-	-	2/30/50/50	0/6/6/6

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	C4X-C10	11.71	1.50	1.38
2	A	601	FAD	C4-C4X	4.12	1.48	1.41
2	A	601	FAD	C9A-C5X	3.23	1.49	1.42
2	A	601	FAD	C9A-N10	3.06	1.42	1.38
3	B	601	AMP	C5-C4	2.79	1.48	1.40
2	A	601	FAD	C10-N1	2.72	1.36	1.33
2	A	601	FAD	C8-C7	2.72	1.47	1.40
2	A	601	FAD	C4X-N5	2.52	1.36	1.33
2	A	601	FAD	C5A-C4A	2.17	1.46	1.40
3	B	601	AMP	C2-N3	2.13	1.35	1.32

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	C4-N3-C2	10.47	123.98	115.14
2	A	601	FAD	C4-C4X-C10	-4.66	116.86	119.95
2	A	601	FAD	C4X-C4-N3	-4.46	117.33	123.43
2	A	601	FAD	C4X-N5-C5X	4.29	121.06	116.77
2	A	601	FAD	N3A-C2A-N1A	-4.06	122.33	128.68
3	B	601	AMP	N3-C2-N1	-3.63	123.01	128.68
3	B	601	AMP	C1'-N9-C4	-3.34	120.77	126.64
2	A	601	FAD	C2A-N1A-C6A	3.20	124.23	118.75
3	B	601	AMP	C2-N1-C6	3.13	124.10	118.75
2	A	601	FAD	C1'-N10-C9A	3.07	120.71	118.29
2	A	601	FAD	N6A-C6A-N1A	2.87	124.54	118.57
2	A	601	FAD	C4-C4X-N5	2.78	121.77	118.60
3	B	601	AMP	N6-C6-N1	2.73	124.24	118.57
2	A	601	FAD	O4'-C4'-C3'	2.61	115.45	109.10
2	A	601	FAD	O2A-PA-O1A	2.60	125.11	112.24
2	A	601	FAD	O4B-C1B-C2B	-2.59	103.13	106.93
2	A	601	FAD	O2'-C2'-C3'	-2.57	102.84	109.10
2	A	601	FAD	C4X-C10-N10	-2.57	117.66	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	O2P-P-O1P	2.49	124.56	112.24
2	A	601	FAD	C1B-N9A-C4A	-2.24	122.70	126.64
2	A	601	FAD	O2'-C2'-C1'	2.23	114.95	109.59
2	A	601	FAD	O4'-C4'-C5'	-2.18	105.03	109.92
2	A	601	FAD	C1'-C2'-C3'	2.11	115.69	109.79
2	A	601	FAD	C4'-C3'-C2'	2.01	117.54	113.36
3	B	601	AMP	O3P-P-O2P	2.00	115.30	107.64

There are no chirality outliers.

All (5) torsion outliers are listed below:

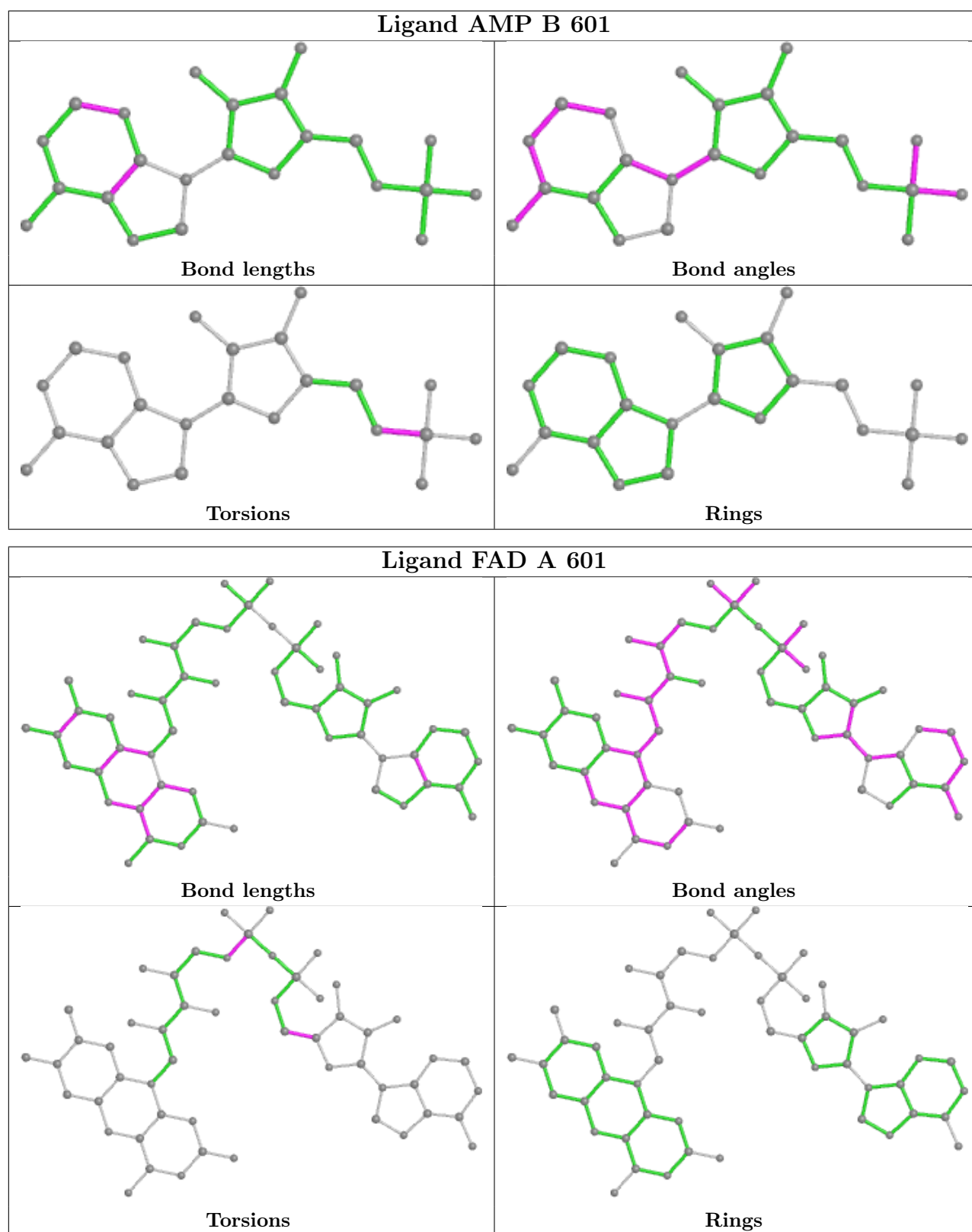
Mol	Chain	Res	Type	Atoms
3	B	601	AMP	C5'-O5'-P-O3P
2	A	601	FAD	C5'-O5'-P-O1P
3	B	601	AMP	C5'-O5'-P-O2P
3	B	601	AMP	C5'-O5'-P-O1P
2	A	601	FAD	O4B-C4B-C5B-O5B

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	7	0

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

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	535/551 (97%)	0.11	18 (3%) 45 48	8, 18, 37, 85	0
1	B	524/551 (95%)	-0.07	17 (3%) 47 50	7, 13, 30, 89	0
All	All	1059/1102 (96%)	0.02	35 (3%) 46 49	7, 15, 35, 89	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	531	SER	11.0
1	A	-1	SER	8.8
1	B	44	ARG	7.5
1	A	0	HIS	6.6
1	B	42	ILE	6.0
1	A	454	TYR	5.8
1	A	530	ALA	5.7
1	B	457	ASN	5.6
1	B	43	PRO	4.8
1	B	45	ILE	4.5
1	B	530	ALA	4.4
1	A	1	MET	4.3
1	A	529	GLY	3.9
1	A	322	ASP	3.8
1	A	323	ASN	3.8
1	B	114	LEU	3.4
1	B	456	GLY	3.4
1	B	41	THR	3.2
1	A	104	HIS	3.1
1	B	448	VAL	3.1
1	B	531	SER	3.0
1	B	458	PHE	2.9
1	A	448	VAL	2.5
1	B	46	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	-8	SER	2.3
1	A	457	ASN	2.3
1	B	449	THR	2.3
1	B	134	GLY	2.3
1	A	114	LEU	2.3
1	A	455	TYR	2.2
1	A	325	GLU	2.2
1	B	137	ASP	2.1
1	A	103	GLY	2.1
1	B	529	GLY	2.1
1	A	456	GLY	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](#)

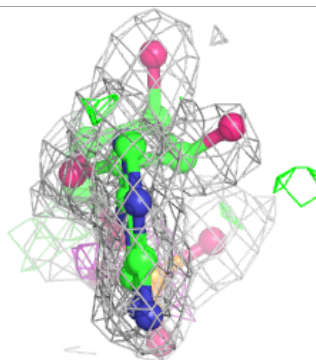
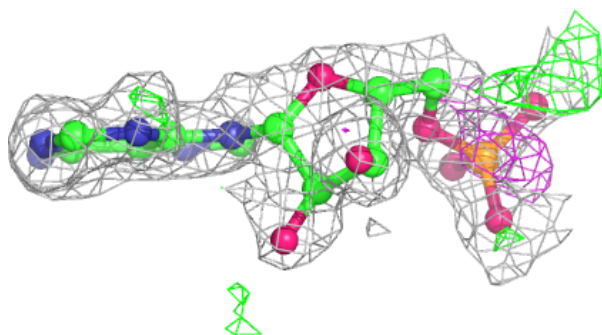
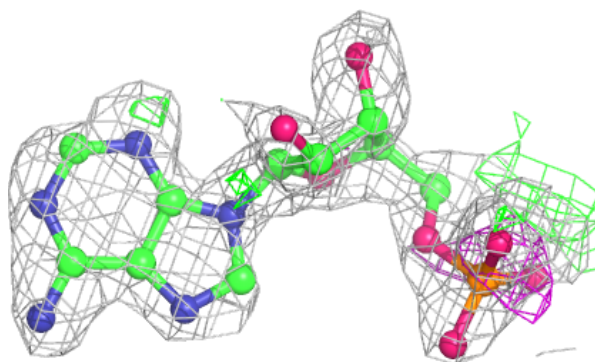
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	AMP	B	601	23/23	0.72	0.25	28,38,57,60	0
2	FAD	A	601	53/53	0.96	0.09	10,16,22,23	0

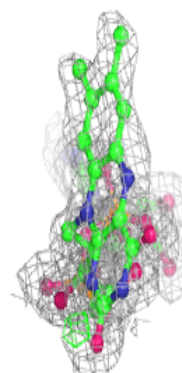
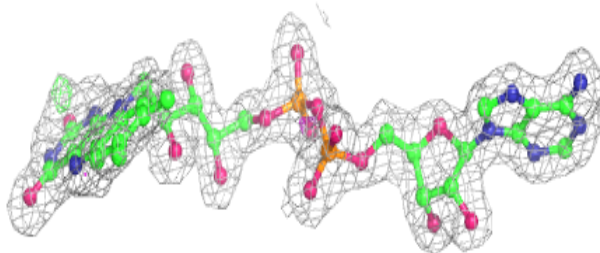
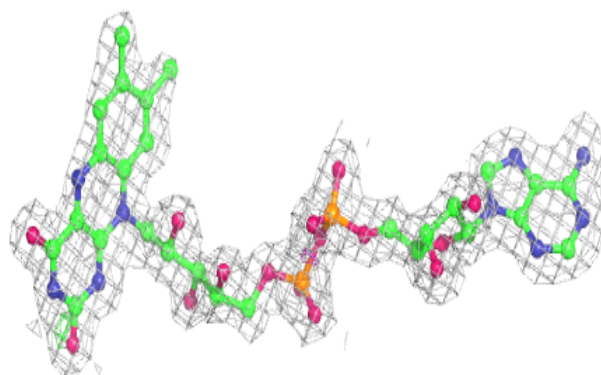
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around AMP B 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 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.