



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 05:06 am BST

PDB ID : 5JZX
Title : Crystal Structure of UDP-N-acetylenolpyruvoylglucosamine reductase (MurB) from *Mycobacterium tuberculosis*
Authors : Dharavath, S.; Eniyan, K.; Bajpai, U.; Gourinath, S.
Deposited on : 2016-05-17
Resolution : 2.20 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 : 2.11
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) : 2.11

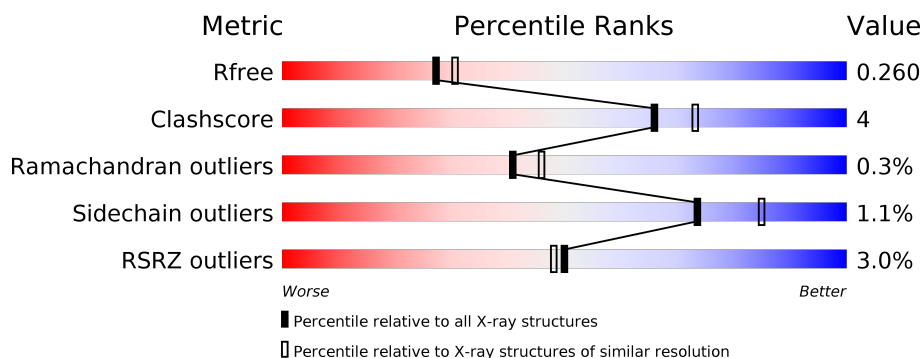
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	405	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 80%; height: 10px; background-color: green;"></div> <div style="width: 9%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: orange;"></div> <div style="width: 10%; height: 10px; background-color: grey;"></div> </div>
1	B	405	<div> <div style="width: 4%; height: 10px; background-color: red;"></div> <div style="width: 71%; height: 10px; background-color: green;"></div> <div style="width: 8%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: orange;"></div> <div style="width: 19%; height: 10px; background-color: grey;"></div> </div>
1	C	405	<div> <div style="width: 78%; height: 10px; background-color: green;"></div> <div style="width: 9%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: orange;"></div> <div style="width: 12%; height: 10px; background-color: grey;"></div> </div>
1	D	405	<div> <div style="width: 81%; height: 10px; background-color: green;"></div> <div style="width: 6%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: orange;"></div> <div style="width: 13%; height: 10px; background-color: grey;"></div> </div>
1	E	405	<div> <div style="width: 4%; height: 10px; background-color: red;"></div> <div style="width: 70%; height: 10px; background-color: green;"></div> <div style="width: 7%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: orange;"></div> <div style="width: 23%; height: 10px; background-color: grey;"></div> </div>
1	F	405	<div> <div style="width: 5%; height: 10px; background-color: red;"></div> <div style="width: 75%; height: 10px; background-color: green;"></div> <div style="width: 8%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: orange;"></div> <div style="width: 16%; height: 10px; background-color: grey;"></div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15461 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 UDP-N-acetylenolpyruvoylglucosamine reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	0	0	0
			2633	1650	478	498	7			
1	B	327	Total	C	N	O	S	0	0	0
			2351	1480	422	442	7			
1	C	355	Total	C	N	O	S	0	0	0
			2599	1631	470	491	7			
1	D	353	Total	C	N	O	S	0	0	0
			2579	1615	468	489	7			
1	E	313	Total	C	N	O	S	0	0	0
			2236	1404	401	424	7			
1	F	339	Total	C	N	O	S	0	0	0
			2379	1498	425	449	7			

There are 216 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	MET	-	expression tag	UNP P9WJL9
A	-34	GLY	-	expression tag	UNP P9WJL9
A	-33	SER	-	expression tag	UNP P9WJL9
A	-32	SER	-	expression tag	UNP P9WJL9
A	-31	HIS	-	expression tag	UNP P9WJL9
A	-30	HIS	-	expression tag	UNP P9WJL9
A	-29	HIS	-	expression tag	UNP P9WJL9
A	-28	HIS	-	expression tag	UNP P9WJL9
A	-27	HIS	-	expression tag	UNP P9WJL9
A	-26	HIS	-	expression tag	UNP P9WJL9
A	-25	SER	-	expression tag	UNP P9WJL9
A	-24	SER	-	expression tag	UNP P9WJL9
A	-23	GLY	-	expression tag	UNP P9WJL9
A	-22	LEU	-	expression tag	UNP P9WJL9
A	-21	VAL	-	expression tag	UNP P9WJL9
A	-20	PRO	-	expression tag	UNP P9WJL9
A	-19	ARG	-	expression tag	UNP P9WJL9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	GLY	-	expression tag	UNP P9WJL9
A	-17	SER	-	expression tag	UNP P9WJL9
A	-16	HIS	-	expression tag	UNP P9WJL9
A	-15	MET	-	expression tag	UNP P9WJL9
A	-14	ALA	-	expression tag	UNP P9WJL9
A	-13	SER	-	expression tag	UNP P9WJL9
A	-12	MET	-	expression tag	UNP P9WJL9
A	-11	THR	-	expression tag	UNP P9WJL9
A	-10	GLY	-	expression tag	UNP P9WJL9
A	-9	GLY	-	expression tag	UNP P9WJL9
A	-8	GLN	-	expression tag	UNP P9WJL9
A	-7	GLN	-	expression tag	UNP P9WJL9
A	-6	MET	-	expression tag	UNP P9WJL9
A	-5	GLY	-	expression tag	UNP P9WJL9
A	-4	ARG	-	expression tag	UNP P9WJL9
A	-3	GLY	-	expression tag	UNP P9WJL9
A	-2	SER	-	expression tag	UNP P9WJL9
A	-1	GLU	-	expression tag	UNP P9WJL9
A	0	PHE	-	expression tag	UNP P9WJL9
B	-35	MET	-	expression tag	UNP P9WJL9
B	-34	GLY	-	expression tag	UNP P9WJL9
B	-33	SER	-	expression tag	UNP P9WJL9
B	-32	SER	-	expression tag	UNP P9WJL9
B	-31	HIS	-	expression tag	UNP P9WJL9
B	-30	HIS	-	expression tag	UNP P9WJL9
B	-29	HIS	-	expression tag	UNP P9WJL9
B	-28	HIS	-	expression tag	UNP P9WJL9
B	-27	HIS	-	expression tag	UNP P9WJL9
B	-26	HIS	-	expression tag	UNP P9WJL9
B	-25	SER	-	expression tag	UNP P9WJL9
B	-24	SER	-	expression tag	UNP P9WJL9
B	-23	GLY	-	expression tag	UNP P9WJL9
B	-22	LEU	-	expression tag	UNP P9WJL9
B	-21	VAL	-	expression tag	UNP P9WJL9
B	-20	PRO	-	expression tag	UNP P9WJL9
B	-19	ARG	-	expression tag	UNP P9WJL9
B	-18	GLY	-	expression tag	UNP P9WJL9
B	-17	SER	-	expression tag	UNP P9WJL9
B	-16	HIS	-	expression tag	UNP P9WJL9
B	-15	MET	-	expression tag	UNP P9WJL9
B	-14	ALA	-	expression tag	UNP P9WJL9
B	-13	SER	-	expression tag	UNP P9WJL9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	MET	-	expression tag	UNP P9WJL9
B	-11	THR	-	expression tag	UNP P9WJL9
B	-10	GLY	-	expression tag	UNP P9WJL9
B	-9	GLY	-	expression tag	UNP P9WJL9
B	-8	GLN	-	expression tag	UNP P9WJL9
B	-7	GLN	-	expression tag	UNP P9WJL9
B	-6	MET	-	expression tag	UNP P9WJL9
B	-5	GLY	-	expression tag	UNP P9WJL9
B	-4	ARG	-	expression tag	UNP P9WJL9
B	-3	GLY	-	expression tag	UNP P9WJL9
B	-2	SER	-	expression tag	UNP P9WJL9
B	-1	GLU	-	expression tag	UNP P9WJL9
B	0	PHE	-	expression tag	UNP P9WJL9
C	-35	MET	-	expression tag	UNP P9WJL9
C	-34	GLY	-	expression tag	UNP P9WJL9
C	-33	SER	-	expression tag	UNP P9WJL9
C	-32	SER	-	expression tag	UNP P9WJL9
C	-31	HIS	-	expression tag	UNP P9WJL9
C	-30	HIS	-	expression tag	UNP P9WJL9
C	-29	HIS	-	expression tag	UNP P9WJL9
C	-28	HIS	-	expression tag	UNP P9WJL9
C	-27	HIS	-	expression tag	UNP P9WJL9
C	-26	HIS	-	expression tag	UNP P9WJL9
C	-25	SER	-	expression tag	UNP P9WJL9
C	-24	SER	-	expression tag	UNP P9WJL9
C	-23	GLY	-	expression tag	UNP P9WJL9
C	-22	LEU	-	expression tag	UNP P9WJL9
C	-21	VAL	-	expression tag	UNP P9WJL9
C	-20	PRO	-	expression tag	UNP P9WJL9
C	-19	ARG	-	expression tag	UNP P9WJL9
C	-18	GLY	-	expression tag	UNP P9WJL9
C	-17	SER	-	expression tag	UNP P9WJL9
C	-16	HIS	-	expression tag	UNP P9WJL9
C	-15	MET	-	expression tag	UNP P9WJL9
C	-14	ALA	-	expression tag	UNP P9WJL9
C	-13	SER	-	expression tag	UNP P9WJL9
C	-12	MET	-	expression tag	UNP P9WJL9
C	-11	THR	-	expression tag	UNP P9WJL9
C	-10	GLY	-	expression tag	UNP P9WJL9
C	-9	GLY	-	expression tag	UNP P9WJL9
C	-8	GLN	-	expression tag	UNP P9WJL9
C	-7	GLN	-	expression tag	UNP P9WJL9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	MET	-	expression tag	UNP P9WJL9
C	-5	GLY	-	expression tag	UNP P9WJL9
C	-4	ARG	-	expression tag	UNP P9WJL9
C	-3	GLY	-	expression tag	UNP P9WJL9
C	-2	SER	-	expression tag	UNP P9WJL9
C	-1	GLU	-	expression tag	UNP P9WJL9
C	0	PHE	-	expression tag	UNP P9WJL9
D	-35	MET	-	expression tag	UNP P9WJL9
D	-34	GLY	-	expression tag	UNP P9WJL9
D	-33	SER	-	expression tag	UNP P9WJL9
D	-32	SER	-	expression tag	UNP P9WJL9
D	-31	HIS	-	expression tag	UNP P9WJL9
D	-30	HIS	-	expression tag	UNP P9WJL9
D	-29	HIS	-	expression tag	UNP P9WJL9
D	-28	HIS	-	expression tag	UNP P9WJL9
D	-27	HIS	-	expression tag	UNP P9WJL9
D	-26	HIS	-	expression tag	UNP P9WJL9
D	-25	SER	-	expression tag	UNP P9WJL9
D	-24	SER	-	expression tag	UNP P9WJL9
D	-23	GLY	-	expression tag	UNP P9WJL9
D	-22	LEU	-	expression tag	UNP P9WJL9
D	-21	VAL	-	expression tag	UNP P9WJL9
D	-20	PRO	-	expression tag	UNP P9WJL9
D	-19	ARG	-	expression tag	UNP P9WJL9
D	-18	GLY	-	expression tag	UNP P9WJL9
D	-17	SER	-	expression tag	UNP P9WJL9
D	-16	HIS	-	expression tag	UNP P9WJL9
D	-15	MET	-	expression tag	UNP P9WJL9
D	-14	ALA	-	expression tag	UNP P9WJL9
D	-13	SER	-	expression tag	UNP P9WJL9
D	-12	MET	-	expression tag	UNP P9WJL9
D	-11	THR	-	expression tag	UNP P9WJL9
D	-10	GLY	-	expression tag	UNP P9WJL9
D	-9	GLY	-	expression tag	UNP P9WJL9
D	-8	GLN	-	expression tag	UNP P9WJL9
D	-7	GLN	-	expression tag	UNP P9WJL9
D	-6	MET	-	expression tag	UNP P9WJL9
D	-5	GLY	-	expression tag	UNP P9WJL9
D	-4	ARG	-	expression tag	UNP P9WJL9
D	-3	GLY	-	expression tag	UNP P9WJL9
D	-2	SER	-	expression tag	UNP P9WJL9
D	-1	GLU	-	expression tag	UNP P9WJL9

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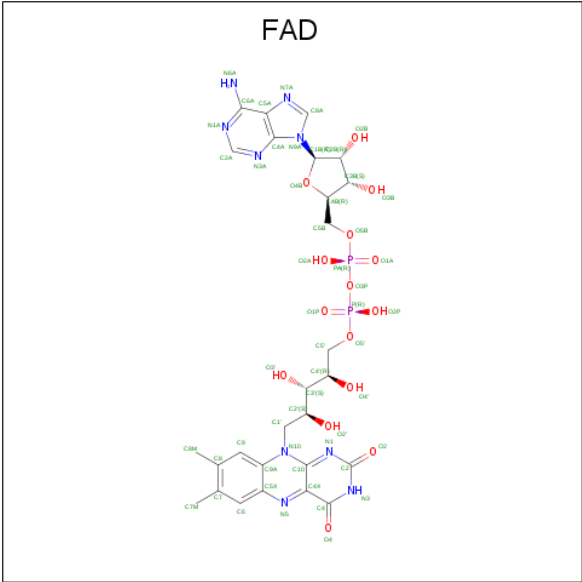
Chain	Residue	Modelled	Actual	Comment	Reference
D	0	PHE	-	expression tag	UNP P9WJL9
E	-35	MET	-	expression tag	UNP P9WJL9
E	-34	GLY	-	expression tag	UNP P9WJL9
E	-33	SER	-	expression tag	UNP P9WJL9
E	-32	SER	-	expression tag	UNP P9WJL9
E	-31	HIS	-	expression tag	UNP P9WJL9
E	-30	HIS	-	expression tag	UNP P9WJL9
E	-29	HIS	-	expression tag	UNP P9WJL9
E	-28	HIS	-	expression tag	UNP P9WJL9
E	-27	HIS	-	expression tag	UNP P9WJL9
E	-26	HIS	-	expression tag	UNP P9WJL9
E	-25	SER	-	expression tag	UNP P9WJL9
E	-24	SER	-	expression tag	UNP P9WJL9
E	-23	GLY	-	expression tag	UNP P9WJL9
E	-22	LEU	-	expression tag	UNP P9WJL9
E	-21	VAL	-	expression tag	UNP P9WJL9
E	-20	PRO	-	expression tag	UNP P9WJL9
E	-19	ARG	-	expression tag	UNP P9WJL9
E	-18	GLY	-	expression tag	UNP P9WJL9
E	-17	SER	-	expression tag	UNP P9WJL9
E	-16	HIS	-	expression tag	UNP P9WJL9
E	-15	MET	-	expression tag	UNP P9WJL9
E	-14	ALA	-	expression tag	UNP P9WJL9
E	-13	SER	-	expression tag	UNP P9WJL9
E	-12	MET	-	expression tag	UNP P9WJL9
E	-11	THR	-	expression tag	UNP P9WJL9
E	-10	GLY	-	expression tag	UNP P9WJL9
E	-9	GLY	-	expression tag	UNP P9WJL9
E	-8	GLN	-	expression tag	UNP P9WJL9
E	-7	GLN	-	expression tag	UNP P9WJL9
E	-6	MET	-	expression tag	UNP P9WJL9
E	-5	GLY	-	expression tag	UNP P9WJL9
E	-4	ARG	-	expression tag	UNP P9WJL9
E	-3	GLY	-	expression tag	UNP P9WJL9
E	-2	SER	-	expression tag	UNP P9WJL9
E	-1	GLU	-	expression tag	UNP P9WJL9
E	0	PHE	-	expression tag	UNP P9WJL9
F	-35	MET	-	expression tag	UNP P9WJL9
F	-34	GLY	-	expression tag	UNP P9WJL9
F	-33	SER	-	expression tag	UNP P9WJL9
F	-32	SER	-	expression tag	UNP P9WJL9
F	-31	HIS	-	expression tag	UNP P9WJL9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-30	HIS	-	expression tag	UNP P9WJL9
F	-29	HIS	-	expression tag	UNP P9WJL9
F	-28	HIS	-	expression tag	UNP P9WJL9
F	-27	HIS	-	expression tag	UNP P9WJL9
F	-26	HIS	-	expression tag	UNP P9WJL9
F	-25	SER	-	expression tag	UNP P9WJL9
F	-24	SER	-	expression tag	UNP P9WJL9
F	-23	GLY	-	expression tag	UNP P9WJL9
F	-22	LEU	-	expression tag	UNP P9WJL9
F	-21	VAL	-	expression tag	UNP P9WJL9
F	-20	PRO	-	expression tag	UNP P9WJL9
F	-19	ARG	-	expression tag	UNP P9WJL9
F	-18	GLY	-	expression tag	UNP P9WJL9
F	-17	SER	-	expression tag	UNP P9WJL9
F	-16	HIS	-	expression tag	UNP P9WJL9
F	-15	MET	-	expression tag	UNP P9WJL9
F	-14	ALA	-	expression tag	UNP P9WJL9
F	-13	SER	-	expression tag	UNP P9WJL9
F	-12	MET	-	expression tag	UNP P9WJL9
F	-11	THR	-	expression tag	UNP P9WJL9
F	-10	GLY	-	expression tag	UNP P9WJL9
F	-9	GLY	-	expression tag	UNP P9WJL9
F	-8	GLN	-	expression tag	UNP P9WJL9
F	-7	GLN	-	expression tag	UNP P9WJL9
F	-6	MET	-	expression tag	UNP P9WJL9
F	-5	GLY	-	expression tag	UNP P9WJL9
F	-4	ARG	-	expression tag	UNP P9WJL9
F	-3	GLY	-	expression tag	UNP P9WJL9
F	-2	SER	-	expression tag	UNP P9WJL9
F	-1	GLU	-	expression tag	UNP P9WJL9
F	0	PHE	-	expression tag	UNP P9WJL9

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	K	0	0
			1	1		
3	A	1	Total	K	0	0
			1	1		
3	C	1	Total	K	0	0
			1	1		

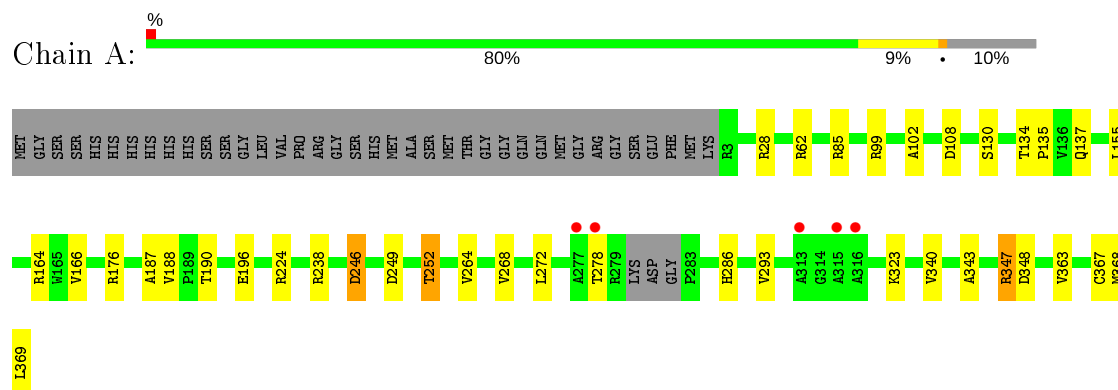
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	109	Total 109	O 109	0	0
4	B	34	Total 34	O 34	0	0
4	C	94	Total 94	O 94	0	0
4	D	91	Total 91	O 91	0	0
4	E	21	Total 21	O 21	0	0
4	F	14	Total 14	O 14	0	0

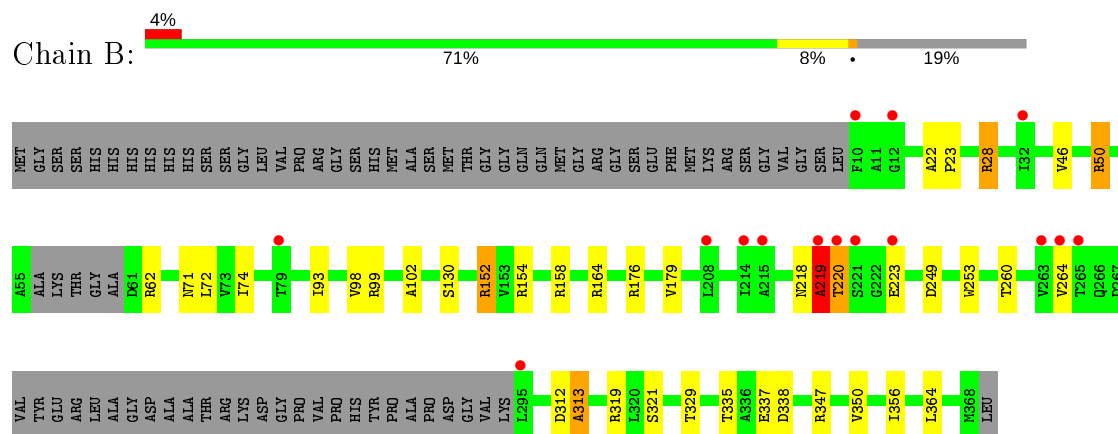
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

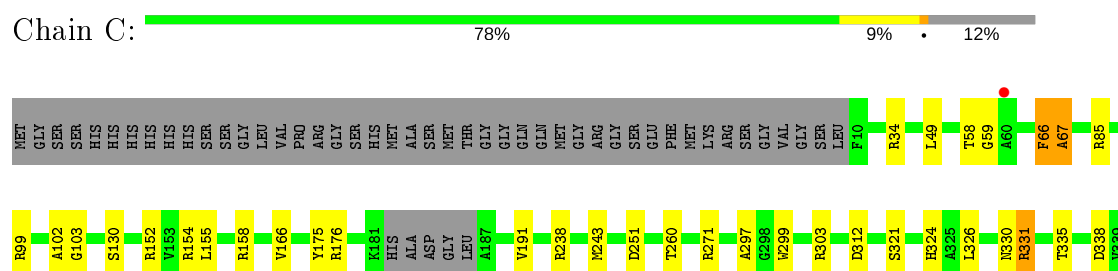
- Molecule 1: UDP-N-acetylenolpyruvoylglucosamine reductase



- Molecule 1: UDP-N-acetylenolpyruvoylglucosamine reductase



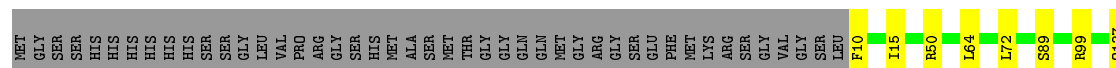
- Molecule 1: UDP-N-acetylenolpyruvoylglucosamine reductase





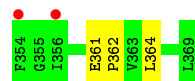
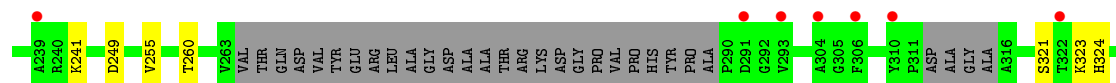
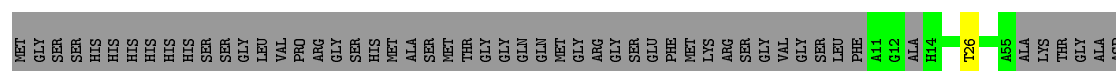
- Molecule 1: UDP-N-acetylenolpyruvoylglucosamine reductase

Chain D: 81% 6% 13%



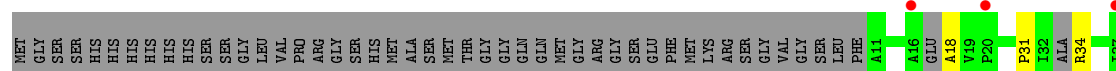
- Molecule 1: UDP-N-acetylenolpyruvoylglucosamine reductase

Chain E: 4% 70% 7% 23%



- Molecule 1: UDP-N-acetylenolpyruvoylglucosamine reductase

Chain F: 5% 75% 8% 16%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	274.92Å 79.97Å 139.00Å 90.00° 112.44° 90.00°	Depositor
Resolution (Å)	38.43 – 2.20 38.43 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (38.43-2.20) 99.0 (38.43-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.207 , 0.259 0.210 , 0.260	Depositor DCC
R_{free} test set	6985 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15461	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: K, 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	0.93	0/2678	1.06	14/3657 (0.4%)
1	B	0.78	1/2389 (0.0%)	1.04	13/3265 (0.4%)
1	C	1.08	6/2644 (0.2%)	1.13	20/3609 (0.6%)
1	D	0.91	1/2623 (0.0%)	1.03	9/3580 (0.3%)
1	E	0.75	0/2269	0.91	3/3097 (0.1%)
1	F	0.70	0/2417	0.86	5/3305 (0.2%)
All	All	0.87	8/15020 (0.1%)	1.01	64/20513 (0.3%)

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	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	67	ALA	N-CA	12.37	1.71	1.46
1	C	34	ARG	C-N	-9.01	1.13	1.34
1	D	331	ARG	CD-NE	-6.09	1.36	1.46
1	C	103	GLY	C-O	-5.96	1.14	1.23
1	C	331	ARG	CD-NE	-5.80	1.36	1.46
1	C	175	TYR	CB-CG	5.29	1.59	1.51
1	C	66	PHE	C-N	-5.21	1.22	1.34
1	B	219	ALA	C-N	-5.11	1.22	1.34

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	ARG	NE-CZ-NH2	-16.49	112.06	120.30
1	D	99	ARG	NE-CZ-NH2	-12.67	113.97	120.30
1	C	99	ARG	NE-CZ-NH2	-11.34	114.63	120.30
1	A	99	ARG	NE-CZ-NH1	10.97	125.79	120.30
1	C	66	PHE	O-C-N	-10.75	105.50	122.70
1	B	220	THR	O-C-N	-10.49	105.91	122.70
1	C	34	ARG	O-C-N	-9.29	107.84	122.70
1	C	66	PHE	C-N-CA	-8.39	100.73	121.70
1	B	50	ARG	NE-CZ-NH1	8.00	124.30	120.30
1	B	50	ARG	NE-CZ-NH2	-7.87	116.36	120.30
1	D	99	ARG	CG-CD-NE	-7.76	95.50	111.80
1	D	331	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	B	220	THR	CA-C-N	7.53	133.76	117.20
1	C	99	ARG	NE-CZ-NH1	7.49	124.05	120.30
1	A	108	ASP	CB-CG-OD1	7.48	125.03	118.30
1	A	99	ARG	CG-CD-NE	-7.40	96.26	111.80
1	C	331	ARG	NE-CZ-NH2	-7.20	116.70	120.30
1	C	238	ARG	NE-CZ-NH1	7.14	123.87	120.30
1	A	108	ASP	CB-CG-OD2	-7.12	111.89	118.30
1	A	28	ARG	NE-CZ-NH2	6.87	123.74	120.30
1	C	67	ALA	N-CA-C	6.77	129.27	111.00
1	A	85	ARG	NE-CZ-NH2	-6.68	116.96	120.30
1	C	34	ARG	C-N-CA	6.62	138.26	121.70
1	C	34	ARG	CA-C-N	6.01	130.42	117.20
1	C	99	ARG	CG-CD-NE	-5.99	99.22	111.80
1	C	154	ARG	NE-CZ-NH1	5.85	123.23	120.30
1	B	164	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	C	271	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	B	158	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	C	85	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	A	246	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	C	238	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	F	281	ASP	CB-CG-OD1	5.52	123.27	118.30
1	A	62	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	B	152	ARG	NE-CZ-NH1	-5.47	117.56	120.30
1	A	28	ARG	NE-CZ-NH1	-5.46	117.57	120.30
1	C	271	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	F	76	GLU	N-CA-C	5.35	125.45	111.00
1	A	246	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	238	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	B	99	ARG	NE-CZ-NH2	-5.27	117.66	120.30
1	D	279	ARG	NE-CZ-NH1	5.27	122.93	120.30
1	C	331	ARG	NE-CZ-NH1	5.27	122.93	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	76	GLU	C-N-CA	5.24	134.80	121.70
1	D	331	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	B	179	VAL	CB-CA-C	-5.21	101.50	111.40
1	A	99	ARG	CD-NE-CZ	5.18	130.86	123.60
1	E	152	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	A	164	ARG	NE-CZ-NH2	-5.15	117.72	120.30
1	D	279	ARG	N-CA-C	5.13	124.84	111.00
1	D	267	ASP	CB-CG-OD1	5.12	122.91	118.30
1	C	99	ARG	CD-NE-CZ	5.10	130.74	123.60
1	B	28	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	D	50	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	F	238	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	B	338	ASP	CB-CG-OD1	5.05	122.85	118.30
1	C	347	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	E	85	ARG	NE-CZ-NH1	5.05	122.82	120.30
1	B	158	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	B	62	ARG	NE-CZ-NH1	5.03	122.82	120.30
1	E	169	ARG	NE-CZ-NH2	-5.02	117.79	120.30
1	F	152	ARG	NE-CZ-NH2	-5.02	117.79	120.30
1	C	66	PHE	CA-C-N	5.01	128.23	117.20
1	D	99	ARG	NE-CZ-NH1	5.00	122.80	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	313	ALA	Peptide

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	2633	0	2638	22	0
1	B	2351	0	2344	20	0
1	C	2599	0	2617	34	0
1	D	2579	0	2591	12	0
1	E	2236	0	2205	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2379	0	2296	17	0
2	A	53	0	31	1	0
2	B	53	0	31	1	0
2	C	53	0	31	0	0
2	D	53	0	31	0	0
2	E	53	0	31	0	0
2	F	53	0	31	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	109	0	0	1	0
4	B	34	0	0	1	0
4	C	94	0	0	1	0
4	D	91	0	0	1	0
4	E	21	0	0	0	0
4	F	14	0	0	0	0
All	All	15461	0	14877	119	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 (119) 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:C:67:ALA:CA	1:C:67:ALA:N	1.71	1.50
1:C:66:PHE:O	1:C:67:ALA:CA	1.93	1.16
1:A:196:GLU:OE1	1:C:152:ARG:NH1	1.83	1.10
1:C:66:PHE:C	1:C:67:ALA:CA	2.27	1.02
1:A:347:ARG:NH1	1:A:348:ASP:OD1	1.97	0.98
1:C:344:ARG:NH1	1:C:348:ASP:OD2	2.06	0.88
1:C:66:PHE:O	1:C:67:ALA:CB	2.27	0.81
1:C:66:PHE:O	1:C:67:ALA:HB3	1.85	0.77
1:F:155:LEU:HD12	1:F:166:VAL:HG21	1.67	0.74
1:C:67:ALA:N	1:C:67:ALA:HA	1.97	0.70
1:A:196:GLU:CD	1:C:152:ARG:NH1	2.43	0.70
1:C:243:MET:HE1	1:C:324:HIS:CD2	2.26	0.70
1:A:246:ASP:H	1:A:252:THR:HG21	1.58	0.69
1:A:224:ARG:NH1	4:A:501:HOH:O	2.26	0.68
1:C:243:MET:CE	1:C:324:HIS:CD2	2.78	0.66
1:C:335:THR:HG21	4:C:581:HOH:O	1.95	0.66
1:A:155:LEU:HD13	1:A:190:THR:HG21	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:134:THR:HG23	1:E:146:VAL:HG11	1.78	0.65
1:C:176:ARG:NH1	1:C:361:GLU:OE1	2.32	0.62
1:C:243:MET:HE1	1:C:324:HIS:HD2	1.65	0.62
1:F:71:ASN:ND2	1:F:137:GLN:OE1	2.30	0.61
1:E:134:THR:HG23	1:E:146:VAL:HG21	1.82	0.61
1:C:335:THR:HG23	1:C:338:ASP:H	1.65	0.61
1:B:154:ARG:NH1	4:B:501:HOH:O	2.35	0.60
1:C:251:ASP:OD1	1:C:331:ARG:NH2	2.32	0.59
1:B:28:ARG:HG3	1:B:253:TRP:CE3	2.39	0.58
1:E:155:LEU:HD13	1:E:190:THR:HG21	1.85	0.58
1:F:137:GLN:O	1:F:176:ARG:HA	2.04	0.58
1:C:49:LEU:HD13	1:C:191:VAL:HG21	1.85	0.58
1:A:343:ALA:HB3	1:A:369:LEU:HD11	1.85	0.57
1:A:196:GLU:CD	1:C:152:ARG:HH12	2.09	0.55
1:F:18:ALA:N	1:F:34:ARG:O	2.39	0.55
1:A:264:VAL:HB	1:A:268:VAL:HG22	1.89	0.55
1:E:249:ASP:OD2	1:E:323:LYS:NZ	2.34	0.54
1:C:158:ARG:HD2	1:C:191:VAL:HG23	1.90	0.54
1:F:246:ASP:H	1:F:252:THR:HG21	1.71	0.54
1:F:246:ASP:N	1:F:252:THR:HG21	2.23	0.54
1:A:137:GLN:O	1:A:176:ARG:HA	2.08	0.53
1:B:152:ARG:NH1	1:D:196:GLU:OE1	2.40	0.53
1:D:10:PHE:CD2	1:D:15:ILE:HD11	2.43	0.53
1:D:72:LEU:HD11	1:D:365:ILE:HD11	1.90	0.53
1:F:155:LEU:HD13	1:F:190:THR:HG21	1.89	0.53
1:B:72:LEU:HD13	1:B:74:ILE:HD11	1.91	0.53
1:A:249:ASP:O	1:A:252:THR:HB	2.10	0.52
1:E:134:THR:HG22	1:E:135:PRO:HD3	1.91	0.52
1:E:134:THR:CG2	1:E:146:VAL:HG11	2.39	0.52
1:D:351:HIS:O	1:D:354:PHE:O	2.28	0.52
1:B:176:ARG:HG3	1:B:260:THR:CG2	2.39	0.51
1:D:89:SER:O	4:D:501:HOH:O	2.19	0.51
1:A:368:MET:O	1:A:369:LEU:HB2	2.11	0.51
1:C:49:LEU:CD1	1:C:191:VAL:HG21	2.41	0.51
1:E:155:LEU:HD22	1:E:190:THR:HG22	1.93	0.50
1:C:66:PHE:O	1:C:67:ALA:C	2.49	0.50
1:B:220:THR:HB	1:B:223:GLU:HG3	1.94	0.50
1:B:46:VAL:O	1:B:50:ARG:HG3	2.12	0.50
1:C:67:ALA:CB	1:C:67:ALA:N	2.65	0.50
1:E:255:VAL:HG21	1:E:362:PRO:HG3	1.93	0.50
1:E:241:LYS:O	1:E:324:HIS:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:321:SER:HB3	1:E:324:HIS:O	2.12	0.49
1:B:220:THR:HB	1:B:223:GLU:CG	2.42	0.49
1:C:340:VAL:HG13	1:C:368:MET:O	2.13	0.49
1:C:243:MET:HE3	1:C:324:HIS:CD2	2.46	0.48
1:E:131:ALA:O	1:E:134:THR:HB	2.12	0.48
1:C:155:LEU:HD12	1:C:166:VAL:HG21	1.96	0.48
1:B:319:ARG:O	1:B:329:THR:OG1	2.27	0.47
1:F:251:ASP:OD2	1:F:322:THR:OG1	2.29	0.47
1:D:249:ASP:OD2	1:D:323:LYS:NZ	2.44	0.47
1:B:22:ALA:HB3	1:B:23:PRO:HD3	1.97	0.47
1:D:72:LEU:HD11	1:D:365:ILE:CD1	2.45	0.46
1:F:76:GLU:N	1:F:77:ASN:CB	2.79	0.46
1:C:297:ALA:HB3	1:C:326:LEU:HD11	1.98	0.46
1:A:155:LEU:HD12	1:A:166:VAL:HG21	1.97	0.46
1:A:340:VAL:HG21	1:A:367:CYS:HB2	1.98	0.46
1:B:176:ARG:HG3	1:B:260:THR:HG22	1.96	0.46
1:A:196:GLU:OE2	1:C:152:ARG:NH1	2.49	0.46
1:F:75:ALA:HB2	1:F:366:GLY:O	2.16	0.46
1:B:347:ARG:O	1:B:350:VAL:HG12	2.16	0.45
1:C:260:THR:HG22	1:C:361:GLU:OE2	2.16	0.45
1:E:137:GLN:O	1:E:176:ARG:HA	2.15	0.45
1:A:249:ASP:OD2	1:A:323:LYS:NZ	2.42	0.45
1:E:72:LEU:HD13	1:E:74:ILE:HD11	1.99	0.45
1:A:246:ASP:N	1:A:252:THR:HG21	2.28	0.45
1:B:93:ILE:CD1	1:B:98:VAL:HG22	2.46	0.45
1:D:64:LEU:HD11	1:D:365:ILE:HD13	1.97	0.45
1:B:249:ASP:OD1	1:B:249:ASP:C	2.56	0.45
1:C:58:THR:HG23	1:C:59:GLY:N	2.32	0.44
1:E:155:LEU:HD13	1:E:190:THR:CG2	2.47	0.44
1:B:102:ALA:O	1:B:130:SER:HB2	2.17	0.44
1:C:299:TRP:CE2	1:C:303:ARG:HG3	2.53	0.44
1:C:321:SER:HB3	1:C:324:HIS:O	2.17	0.44
1:E:155:LEU:HD12	1:E:166:VAL:HG21	2.00	0.44
1:F:72:LEU:HD23	1:F:74:ILE:HD11	2.00	0.44
1:A:363:VAL:HG21	2:A:401:FAD:H4B	1.99	0.43
1:B:335:THR:HG22	1:B:337:GLU:N	2.32	0.43
1:D:10:PHE:HD2	1:D:15:ILE:HD11	1.83	0.43
1:A:134:THR:OG1	1:A:135:PRO:HD3	2.18	0.43
1:F:320:LEU:HD13	1:F:325:ALA:O	2.18	0.43
1:F:230:VAL:O	1:F:234:VAL:HG23	2.19	0.43
1:A:286:HIS:HB2	1:A:293:VAL:CG1	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:260:THR:HG22	1:E:361:GLU:OE1	2.19	0.43
1:B:264:VAL:HG11	1:B:356:ILE:HD12	2.01	0.43
1:C:340:VAL:CG1	1:C:368:MET:O	2.67	0.43
1:F:67:ALA:O	2:F:401:FAD:O1P	2.38	0.42
1:E:26:THR:OG1	1:E:70:SER:HA	2.19	0.42
1:B:312:ASP:HA	1:B:313:ALA:HA	1.80	0.42
1:C:102:ALA:O	1:C:130:SER:HB2	2.20	0.42
1:E:137:GLN:NE2	1:E:361:GLU:O	2.50	0.42
1:E:153:VAL:HG13	1:E:192:VAL:HG13	2.02	0.41
1:A:102:ALA:O	1:A:130:SER:HB2	2.20	0.41
1:D:321:SER:HB3	1:D:324:HIS:O	2.21	0.41
1:F:208:LEU:O	1:F:214:ILE:HD11	2.21	0.41
1:D:137:GLN:O	1:D:176:ARG:HA	2.20	0.41
1:F:261:ASN:HB3	1:F:295:LEU:O	2.20	0.41
1:B:218:ASN:O	1:B:219:ALA:O	2.38	0.40
1:A:187:ALA:O	1:A:188:VAL:C	2.60	0.40
1:C:330:ASN:OD1	1:C:330:ASN:C	2.60	0.40
1:B:71:ASN:HB2	2:B:401:FAD:H2'	2.03	0.40
1:D:153:VAL:HG13	1:D:192:VAL:HG13	2.04	0.40
1:F:134:THR:HB	1:F:146:VAL:HG11	2.04	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	360/405 (89%)	345 (96%)	15 (4%)	0	100	100
1	B	321/405 (79%)	305 (95%)	15 (5%)	1 (0%)	41	46
1	C	351/405 (87%)	338 (96%)	13 (4%)	0	100	100
1	D	349/405 (86%)	339 (97%)	10 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	301/405 (74%)	288 (96%)	11 (4%)	2 (1%)	22	22
1	F	325/405 (80%)	308 (95%)	14 (4%)	3 (1%)	17	16
All	All	2007/2430 (83%)	1923 (96%)	78 (4%)	6 (0%)	41	46

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	219	ALA
1	F	176	ARG
1	E	77	ASN
1	E	205	SER
1	F	31	PRO
1	F	332	GLY

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	A	266/306 (87%)	262 (98%)	4 (2%)	65	78
1	B	235/306 (77%)	233 (99%)	2 (1%)	78	88
1	C	265/306 (87%)	263 (99%)	2 (1%)	81	90
1	D	263/306 (86%)	261 (99%)	2 (1%)	81	90
1	E	223/306 (73%)	221 (99%)	2 (1%)	78	88
1	F	226/306 (74%)	222 (98%)	4 (2%)	59	72
All	All	1478/1836 (80%)	1462 (99%)	16 (1%)	73	85

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	252	THR
1	A	272	LEU
1	A	278	THR
1	A	347	ARG

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Mol	Chain	Res	Type
1	B	321	SER
1	B	364	LEU
1	C	312	ASP
1	C	340	VAL
1	D	220	THR
1	D	340	VAL
1	E	134	THR
1	E	364	LEU
1	F	130	SER
1	F	232	GLU
1	F	252	THR
1	F	367	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	324	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FAD	F	401	-	51,58,58	2.16	9 (17%)	60,89,89	2.03	15 (25%)
2	FAD	D	401	-	51,58,58	1.98	8 (15%)	60,89,89	2.35	14 (23%)
2	FAD	B	401	3	51,58,58	2.02	7 (13%)	60,89,89	2.08	17 (28%)
2	FAD	E	401	-	51,58,58	2.05	8 (15%)	60,89,89	2.06	14 (23%)
2	FAD	C	401	-	51,58,58	2.23	8 (15%)	60,89,89	2.12	12 (20%)
2	FAD	A	401	-	51,58,58	2.16	13 (25%)	60,89,89	2.51	18 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	F	401	-	-	6/30/50/50	0/6/6/6
2	FAD	D	401	-	-	3/30/50/50	0/6/6/6
2	FAD	B	401	3	-	3/30/50/50	0/6/6/6
2	FAD	E	401	-	-	4/30/50/50	0/6/6/6
2	FAD	C	401	-	-	2/30/50/50	0/6/6/6
2	FAD	A	401	-	-	2/30/50/50	0/6/6/6

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	FAD	C4X-C10	11.72	1.50	1.38
2	F	401	FAD	C4X-C10	11.61	1.50	1.38
2	E	401	FAD	C4X-C10	10.41	1.49	1.38
2	A	401	FAD	C4X-C10	10.41	1.49	1.38
2	D	401	FAD	C4X-C10	10.15	1.49	1.38
2	B	401	FAD	C4X-C10	9.68	1.48	1.38
2	A	401	FAD	C9A-N10	5.70	1.46	1.38
2	B	401	FAD	C4-C4X	5.33	1.50	1.41
2	D	401	FAD	C4-C4X	4.32	1.48	1.41
2	B	401	FAD	C9A-C5X	4.32	1.51	1.42
2	E	401	FAD	C4-C4X	4.21	1.48	1.41
2	C	401	FAD	C4-C4X	4.12	1.48	1.41
2	F	401	FAD	C9A-C5X	4.11	1.50	1.42
2	E	401	FAD	C9A-C5X	3.82	1.50	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	FAD	C9A-N10	3.71	1.43	1.38
2	C	401	FAD	C6-C5X	-3.67	1.36	1.41
2	E	401	FAD	C9A-N10	3.67	1.43	1.38
2	C	401	FAD	C5'-C4'	3.64	1.56	1.51
2	A	401	FAD	C4X-N5	3.47	1.38	1.33
2	F	401	FAD	C4-C4X	3.46	1.47	1.41
2	F	401	FAD	C9A-N10	3.36	1.43	1.38
2	E	401	FAD	C8-C7	3.19	1.48	1.40
2	F	401	FAD	C8-C7	3.12	1.48	1.40
2	B	401	FAD	C9A-N10	3.08	1.42	1.38
2	F	401	FAD	C2A-N3A	3.03	1.37	1.32
2	A	401	FAD	C4A-N3A	-2.99	1.31	1.35
2	C	401	FAD	C1'-N10	2.89	1.51	1.48
2	B	401	FAD	C8-C7	2.76	1.47	1.40
2	E	401	FAD	C5A-C4A	2.75	1.48	1.40
2	C	401	FAD	O4B-C1B	-2.73	1.37	1.41
2	A	401	FAD	C4-C4X	2.63	1.45	1.41
2	C	401	FAD	C9A-C5X	2.61	1.47	1.42
2	F	401	FAD	C10-N1	2.60	1.36	1.33
2	D	401	FAD	C2B-C1B	2.59	1.57	1.53
2	A	401	FAD	C2B-C1B	2.50	1.57	1.53
2	A	401	FAD	C10-N1	2.47	1.36	1.33
2	A	401	FAD	C2B-C3B	-2.47	1.46	1.53
2	D	401	FAD	C8-C7	2.44	1.47	1.40
2	F	401	FAD	C5A-C4A	2.38	1.47	1.40
2	A	401	FAD	C9A-C5X	2.29	1.47	1.42
2	E	401	FAD	C2A-N3A	2.25	1.35	1.32
2	D	401	FAD	C5X-N5	2.23	1.39	1.35
2	A	401	FAD	C2A-N3A	2.21	1.35	1.32
2	D	401	FAD	C2B-C3B	-2.15	1.47	1.53
2	A	401	FAD	C5A-C4A	2.12	1.46	1.40
2	A	401	FAD	C9-C9A	-2.12	1.36	1.40
2	D	401	FAD	C10-N1	2.11	1.36	1.33
2	B	401	FAD	C5A-N7A	-2.08	1.32	1.39
2	E	401	FAD	C10-N1	2.06	1.35	1.33
2	D	401	FAD	O3B-C3B	2.06	1.47	1.43
2	B	401	FAD	C2'-C3'	2.05	1.57	1.53
2	F	401	FAD	C1'-N10	-2.02	1.46	1.48
2	A	401	FAD	C5X-N5	2.00	1.38	1.35

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	FAD	C4-N3-C2	11.76	125.07	115.14
2	A	401	FAD	C4-N3-C2	10.53	124.04	115.14
2	F	401	FAD	C4-N3-C2	8.66	122.45	115.14
2	E	401	FAD	C4-N3-C2	7.77	121.70	115.14
2	C	401	FAD	C4-N3-C2	7.20	121.22	115.14
2	A	401	FAD	C4-C4X-C10	-7.19	115.19	119.95
2	C	401	FAD	C1'-N10-C9A	7.10	123.88	118.29
2	A	401	FAD	C1'-N10-C9A	6.54	123.44	118.29
2	B	401	FAD	C1'-N10-C9A	6.02	123.03	118.29
2	C	401	FAD	C4-C4X-C10	-5.92	116.03	119.95
2	B	401	FAD	C4-N3-C2	5.88	120.11	115.14
2	B	401	FAD	C4-C4X-C10	-5.42	116.36	119.95
2	F	401	FAD	C1'-N10-C9A	5.30	122.47	118.29
2	E	401	FAD	C1'-N10-C9A	5.29	122.45	118.29
2	E	401	FAD	C4-C4X-C10	-5.05	116.61	119.95
2	D	401	FAD	C4-C4X-C10	-4.95	116.67	119.95
2	D	401	FAD	C4X-C4-N3	-4.87	116.78	123.43
2	D	401	FAD	C2A-N1A-C6A	4.62	126.65	118.75
2	C	401	FAD	C9A-N10-C10	-4.60	115.88	121.91
2	F	401	FAD	C4X-N5-C5X	4.30	121.07	116.77
2	E	401	FAD	O4B-C1B-C2B	-4.26	100.70	106.93
2	D	401	FAD	N3A-C2A-N1A	-4.26	122.02	128.68
2	C	401	FAD	C4-C4X-N5	4.24	123.44	118.60
2	B	401	FAD	C4X-N5-C5X	4.14	120.91	116.77
2	B	401	FAD	C4-C4X-N5	4.11	123.29	118.60
2	B	401	FAD	P-O3P-PA	-3.99	119.14	132.83
2	D	401	FAD	C1'-N10-C9A	3.95	121.41	118.29
2	A	401	FAD	O3B-C3B-C2B	-3.88	99.28	111.82
2	A	401	FAD	C4-C4X-N5	3.77	122.91	118.60
2	F	401	FAD	C4-C4X-C10	-3.73	117.48	119.95
2	C	401	FAD	C4X-C4-N3	-3.65	118.44	123.43
2	A	401	FAD	C4X-N5-C5X	3.65	120.42	116.77
2	E	401	FAD	C4A-C5A-N7A	-3.64	105.60	109.40
2	B	401	FAD	N3A-C2A-N1A	-3.53	123.16	128.68
2	E	401	FAD	C4X-N5-C5X	3.51	120.28	116.77
2	F	401	FAD	P-O3P-PA	-3.50	120.83	132.83
2	A	401	FAD	C5X-C9A-N10	3.47	120.23	117.72
2	B	401	FAD	C9A-N10-C10	-3.38	117.48	121.91
2	A	401	FAD	P-O3P-PA	-3.34	121.37	132.83
2	C	401	FAD	O3B-C3B-C2B	-3.33	101.06	111.82
2	A	401	FAD	O4B-C1B-C2B	-3.30	102.11	106.93
2	F	401	FAD	C4X-C4-N3	-3.27	118.96	123.43
2	A	401	FAD	N3A-C2A-N1A	-3.19	123.70	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	FAD	N6A-C6A-N1A	3.18	125.17	118.57
2	E	401	FAD	C4X-C4-N3	-3.17	119.09	123.43
2	E	401	FAD	P-O3P-PA	-3.14	122.06	132.83
2	E	401	FAD	O2'-C2'-C3'	3.11	116.66	109.10
2	A	401	FAD	N6A-C6A-N1A	3.08	124.97	118.57
2	D	401	FAD	C4-C4X-N5	3.08	122.11	118.60
2	C	401	FAD	N3A-C2A-N1A	-3.05	123.91	128.68
2	D	401	FAD	N6A-C6A-N1A	3.01	124.81	118.57
2	A	401	FAD	C4X-C4-N3	-2.96	119.39	123.43
2	D	401	FAD	O3B-C3B-C2B	-2.96	102.26	111.82
2	E	401	FAD	N3A-C2A-N1A	-2.89	124.16	128.68
2	B	401	FAD	C4X-C4-N3	-2.86	119.51	123.43
2	A	401	FAD	O2B-C2B-C3B	-2.77	102.87	111.82
2	A	401	FAD	C9A-N10-C10	-2.75	118.31	121.91
2	E	401	FAD	C9A-N10-C10	-2.73	118.33	121.91
2	B	401	FAD	C4A-C5A-N7A	-2.70	106.59	109.40
2	C	401	FAD	C1B-N9A-C4A	-2.62	122.04	126.64
2	F	401	FAD	O3B-C3B-C2B	2.60	120.22	111.82
2	B	401	FAD	C3B-C2B-C1B	2.59	104.89	100.98
2	B	401	FAD	C1B-N9A-C4A	-2.59	122.09	126.64
2	B	401	FAD	O2A-PA-O1A	2.59	125.03	112.24
2	E	401	FAD	C4-C4X-N5	2.57	121.53	118.60
2	A	401	FAD	O2B-C2B-C1B	2.49	120.05	110.85
2	C	401	FAD	C5X-C9A-N10	2.47	119.50	117.72
2	D	401	FAD	C5A-C6A-N1A	-2.46	114.77	120.35
2	E	401	FAD	O3B-C3B-C2B	-2.42	103.98	111.82
2	F	401	FAD	C9A-N10-C10	-2.42	118.74	121.91
2	D	401	FAD	O4B-C1B-C2B	-2.35	103.49	106.93
2	F	401	FAD	O5'-C5'-C4'	-2.35	103.10	109.36
2	F	401	FAD	C2B-C3B-C4B	-2.33	98.11	102.64
2	F	401	FAD	N3A-C2A-N1A	-2.33	125.04	128.68
2	A	401	FAD	O5B-C5B-C4B	2.27	116.79	108.99
2	A	401	FAD	C5A-C6A-N6A	-2.24	116.95	120.35
2	F	401	FAD	C4A-C5A-N7A	-2.18	107.13	109.40
2	D	401	FAD	O2'-C2'-C1'	2.18	114.84	109.59
2	A	401	FAD	C6-C7-C8	-2.16	116.28	119.91
2	B	401	FAD	C4'-C3'-C2'	-2.14	108.92	113.36
2	B	401	FAD	O2B-C2B-C3B	-2.08	105.09	111.82
2	B	401	FAD	C5'-C4'-C3'	2.07	116.21	112.20
2	B	401	FAD	C2A-N1A-C6A	2.07	122.29	118.75
2	C	401	FAD	C4'-C3'-C2'	-2.06	109.07	113.36
2	D	401	FAD	P-O3P-PA	-2.05	125.78	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	FAD	C1'-N10-C10	2.05	120.24	118.41
2	D	401	FAD	C4X-N5-C5X	2.04	118.81	116.77
2	F	401	FAD	C4-C4X-N5	2.04	120.93	118.60
2	F	401	FAD	C5X-C9A-N10	2.02	119.18	117.72
2	E	401	FAD	O4'-C4'-C3'	-2.01	104.21	109.10

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	401	FAD	C5B-O5B-PA-O1A
2	F	401	FAD	N10-C1'-C2'-O2'
2	D	401	FAD	N10-C1'-C2'-O2'
2	B	401	FAD	N10-C1'-C2'-O2'
2	E	401	FAD	N10-C1'-C2'-O2'
2	C	401	FAD	N10-C1'-C2'-O2'
2	A	401	FAD	N10-C1'-C2'-O2'
2	F	401	FAD	C5B-O5B-PA-O3P
2	D	401	FAD	C5'-O5'-P-O3P
2	F	401	FAD	C5B-O5B-PA-O2A
2	E	401	FAD	C5B-O5B-PA-O1A
2	F	401	FAD	N10-C1'-C2'-C3'
2	D	401	FAD	N10-C1'-C2'-C3'
2	B	401	FAD	N10-C1'-C2'-C3'
2	E	401	FAD	N10-C1'-C2'-C3'
2	C	401	FAD	N10-C1'-C2'-C3'
2	A	401	FAD	N10-C1'-C2'-C3'
2	B	401	FAD	C5'-O5'-P-O3P
2	E	401	FAD	C5B-O5B-PA-O3P
2	F	401	FAD	C5'-O5'-P-O1P

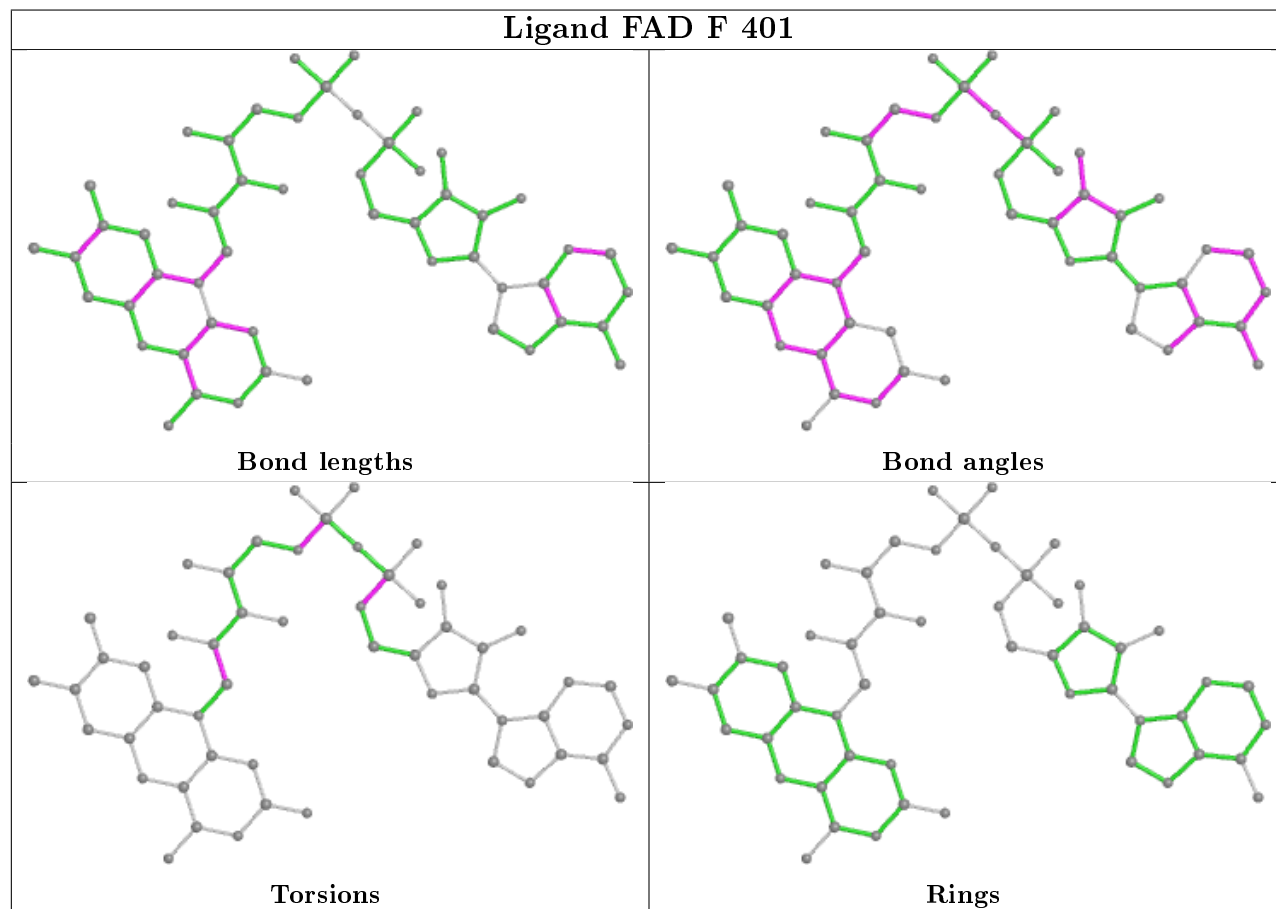
There are no ring outliers.

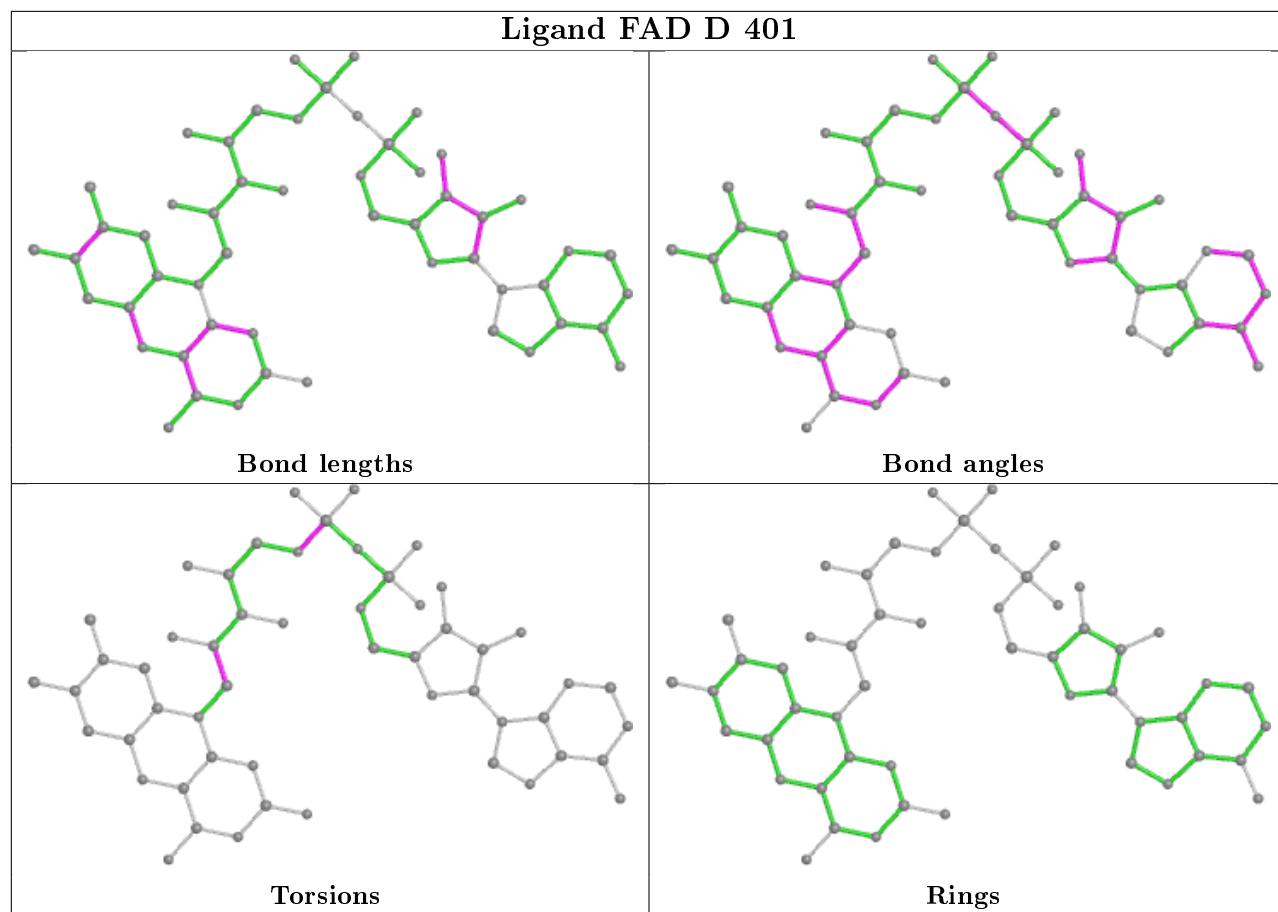
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	401	FAD	1	0
2	B	401	FAD	1	0
2	A	401	FAD	1	0

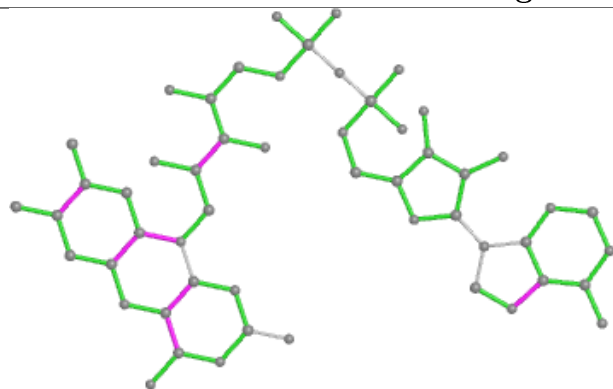
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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.

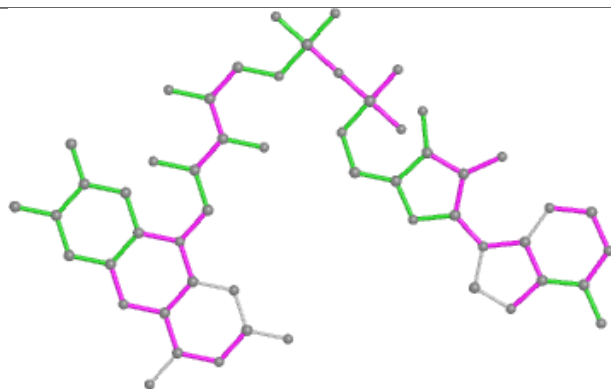




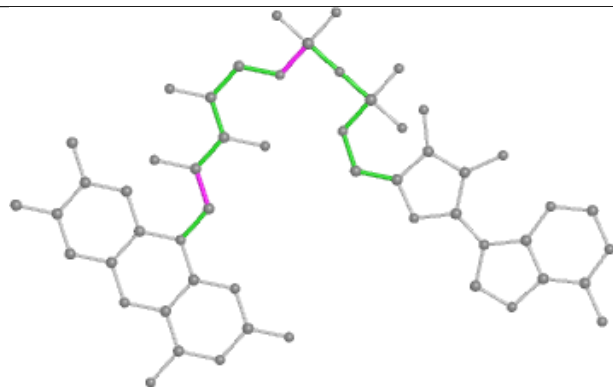
Ligand FAD B 401



Bond lengths



Bond angles

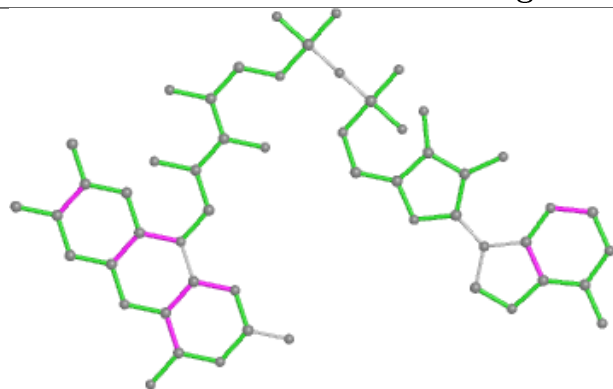


Torsions

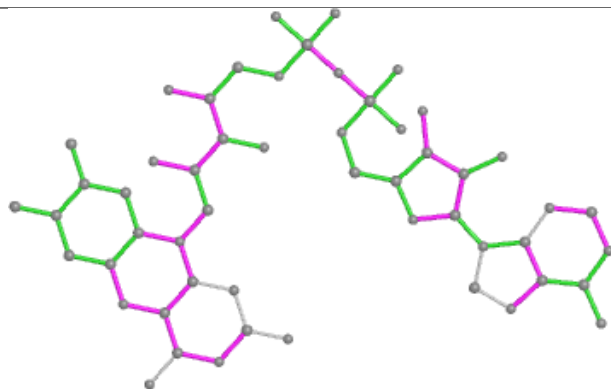


Rings

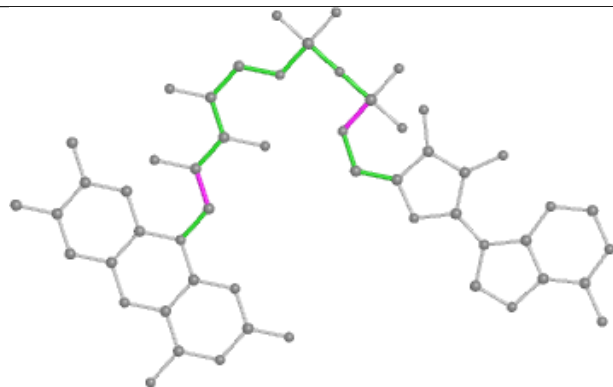
Ligand FAD E 401



Bond lengths



Bond angles

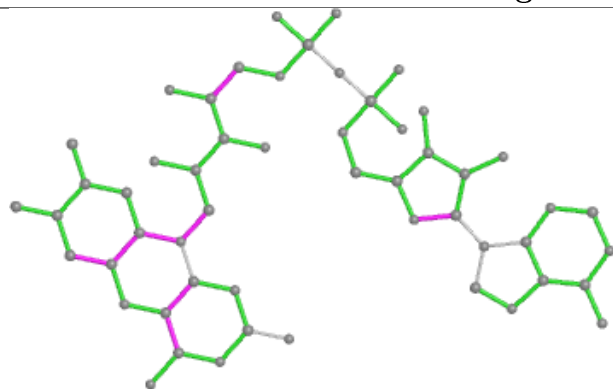


Torsions

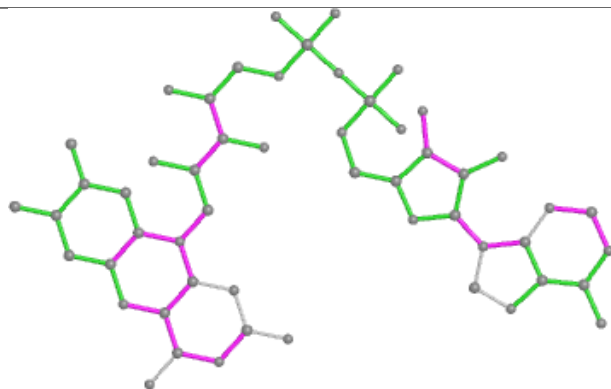


Rings

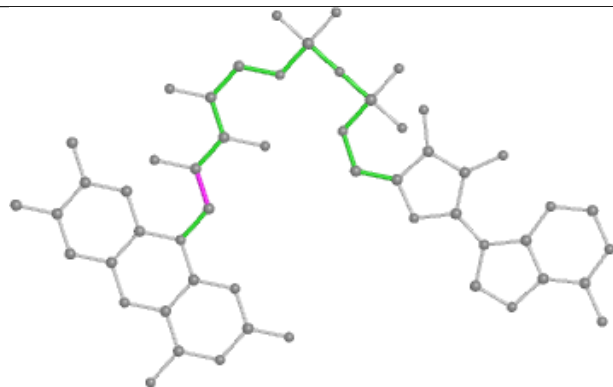
Ligand FAD C 401



Bond lengths



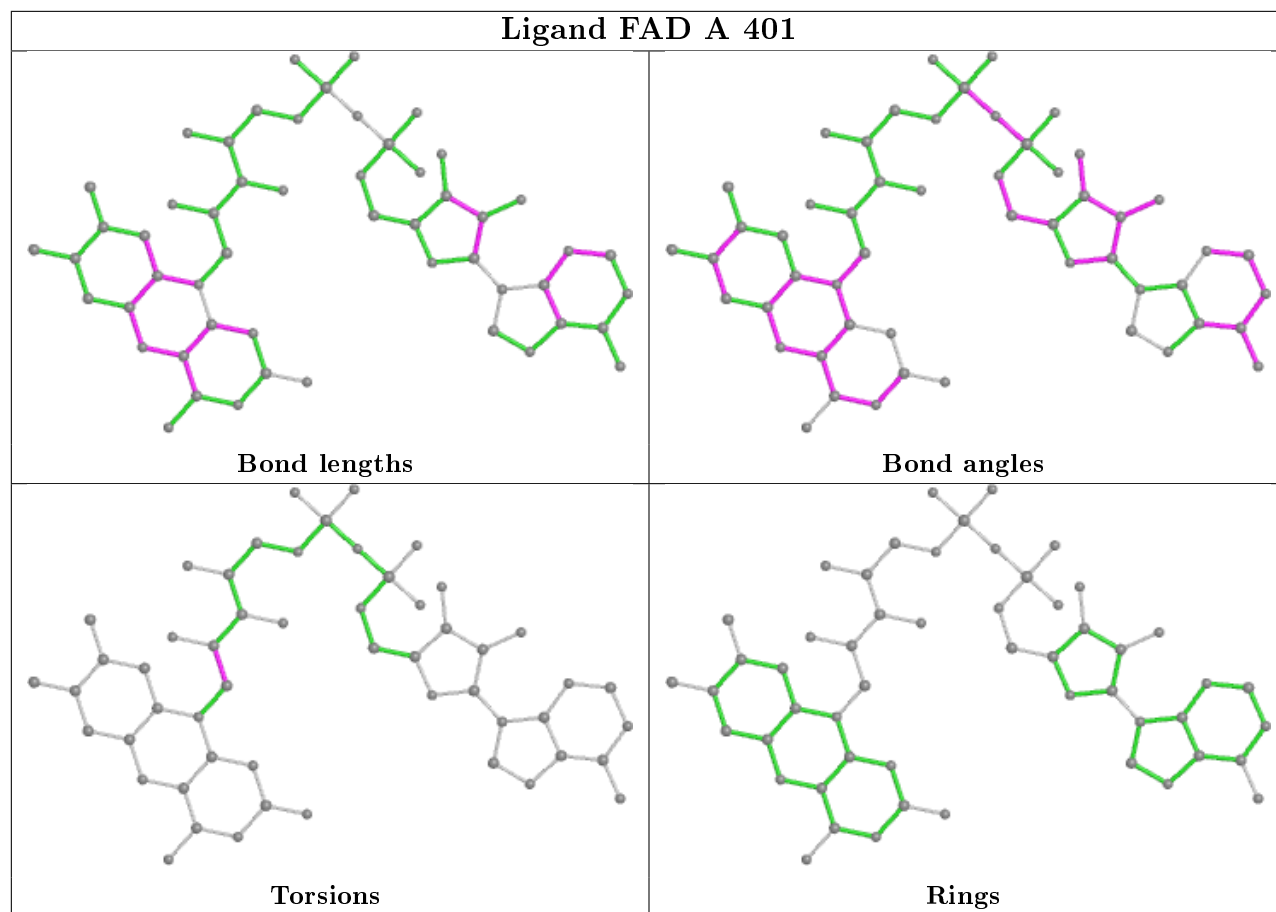
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	34:ARG	C	35:ARG	N	1.13

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	364/405 (89%)	-0.24	5 (1%) 75 73	20, 36, 74, 102	0
1	B	327/405 (80%)	-0.03	15 (4%) 32 31	28, 47, 76, 95	0
1	C	355/405 (87%)	-0.31	1 (0%) 94 93	20, 33, 58, 86	0
1	D	353/405 (87%)	-0.34	2 (0%) 89 88	22, 36, 64, 88	0
1	E	313/405 (77%)	0.20	18 (5%) 23 22	36, 55, 78, 94	0
1	F	339/405 (83%)	0.27	21 (6%) 20 19	36, 61, 80, 94	0
All	All	2051/2430 (84%)	-0.08	62 (3%) 50 48	20, 45, 76, 102	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	278	THR	5.0
1	B	263	VAL	4.3
1	B	264	VAL	4.3
1	E	214	ILE	3.6
1	B	79	THR	3.6
1	C	60	ALA	3.5
1	E	209	ARG	3.4
1	F	248	THR	3.4
1	F	20	PRO	3.3
1	F	16	ALA	3.2
1	F	283	PRO	3.2
1	E	304	ALA	3.1
1	F	276	ALA	3.1
1	B	295	LEU	3.1
1	E	291	ASP	3.1
1	B	265	THR	3.1
1	B	214	ILE	3.1
1	B	220	THR	3.0
1	F	333	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	322	THR	3.0
1	A	277	ALA	2.9
1	B	219	ALA	2.9
1	E	236	ALA	2.9
1	E	206	ALA	2.9
1	F	37	ILE	2.9
1	E	354	PHE	2.8
1	B	221	SER	2.8
1	D	290	PRO	2.7
1	E	225	ALA	2.7
1	E	306	PHE	2.7
1	F	217	LEU	2.7
1	A	315	ALA	2.7
1	F	332	GLY	2.7
1	B	32	ILE	2.7
1	E	210	TYR	2.6
1	F	281	ASP	2.6
1	F	51	HIS	2.6
1	F	282	GLY	2.6
1	E	227	PRO	2.5
1	E	310	TYR	2.5
1	B	10	PHE	2.5
1	E	239	ALA	2.4
1	D	291	ASP	2.4
1	F	291	ASP	2.4
1	B	223	GLU	2.4
1	E	213	LEU	2.3
1	B	215	ALA	2.3
1	E	293	VAL	2.3
1	E	356	ILE	2.3
1	F	266	GLN	2.2
1	A	313	ALA	2.2
1	F	274	GLY	2.2
1	A	278	THR	2.2
1	B	208	LEU	2.2
1	F	208	LEU	2.2
1	F	54	SER	2.1
1	F	275	ASP	2.1
1	B	12	GLY	2.1
1	E	208	LEU	2.1
1	F	269	TYR	2.1
1	A	316	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	273	ALA	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 carbohydrates in this entry.

6.4 Ligands [i](#)

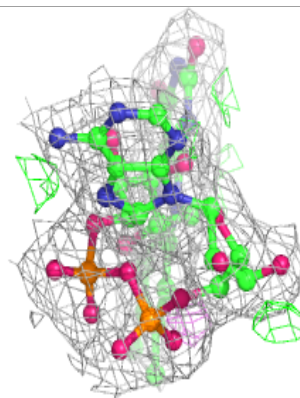
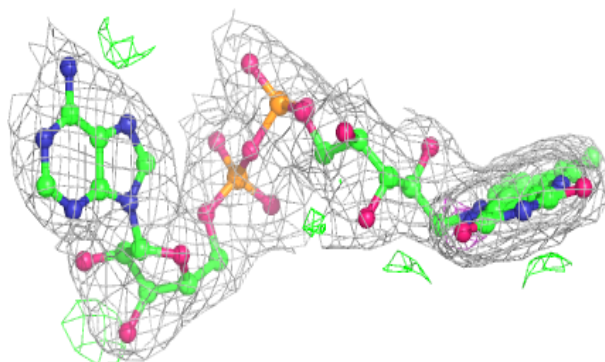
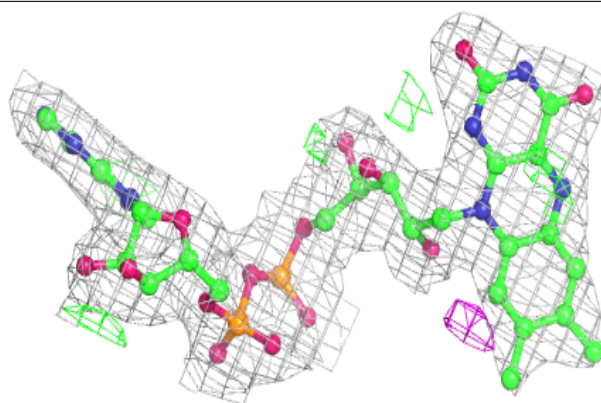
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	K	C	402	1/1	0.84	0.19	79,79,79,79	0
3	K	A	402	1/1	0.89	0.17	88,88,88,88	0
3	K	B	402	1/1	0.94	0.14	81,81,81,81	0
2	FAD	E	401	53/53	0.96	0.13	29,44,53,57	0
2	FAD	F	401	53/53	0.96	0.14	41,50,60,63	0
2	FAD	A	401	53/53	0.98	0.15	21,25,34,42	0
2	FAD	B	401	53/53	0.98	0.13	27,36,44,59	0
2	FAD	D	401	53/53	0.98	0.11	22,26,31,43	0
2	FAD	C	401	53/53	0.98	0.14	20,24,28,30	0

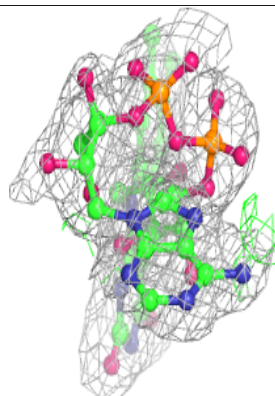
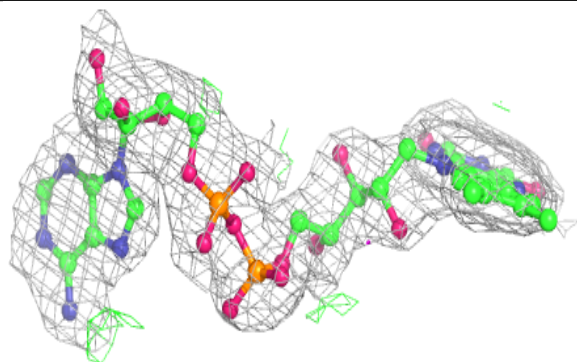
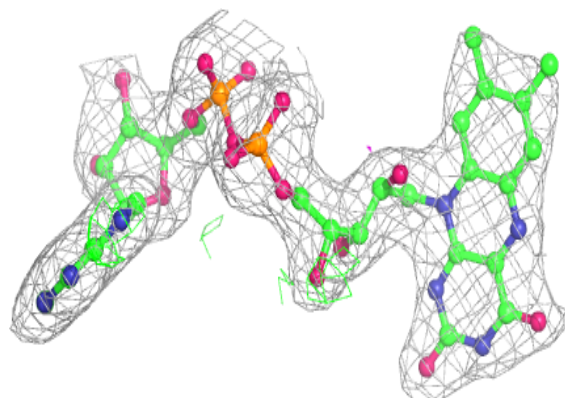
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD E 401:

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

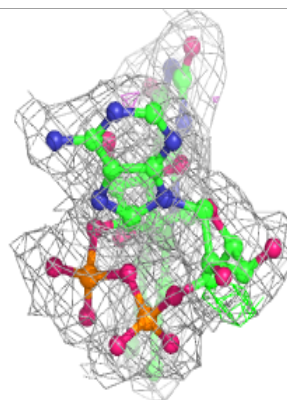
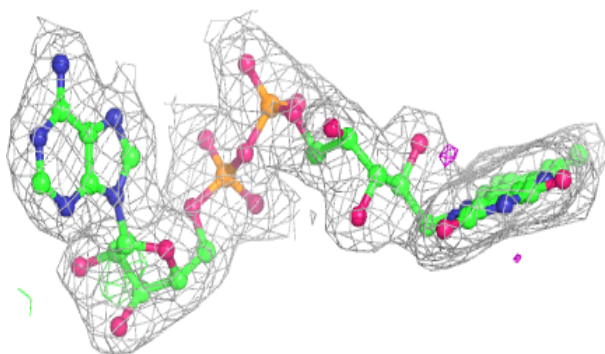
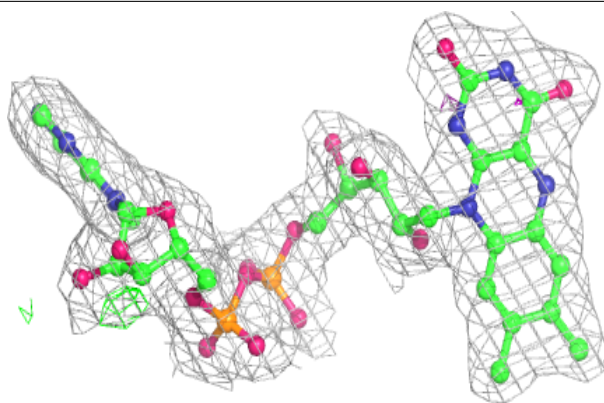
**Electron density around FAD F 401:**

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

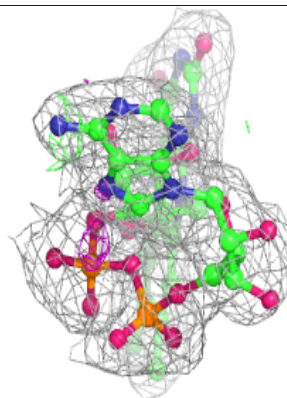
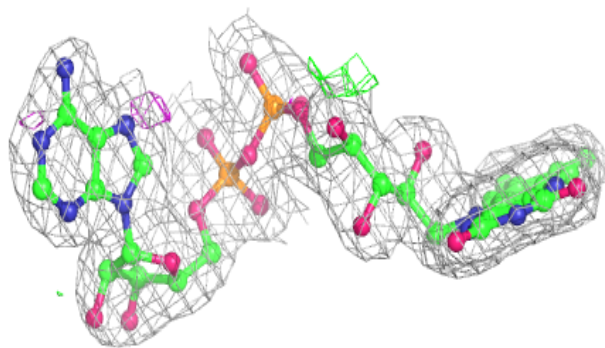
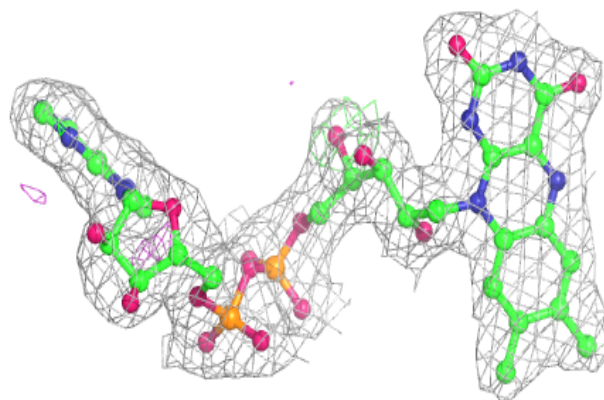


Electron density around FAD A 401:

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

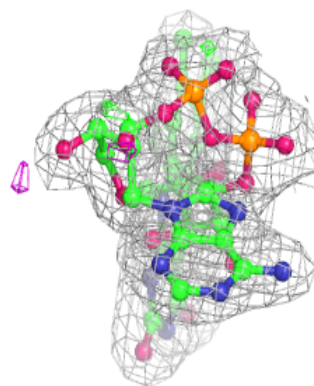
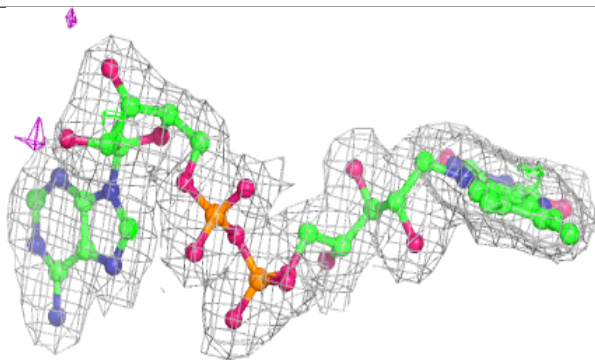
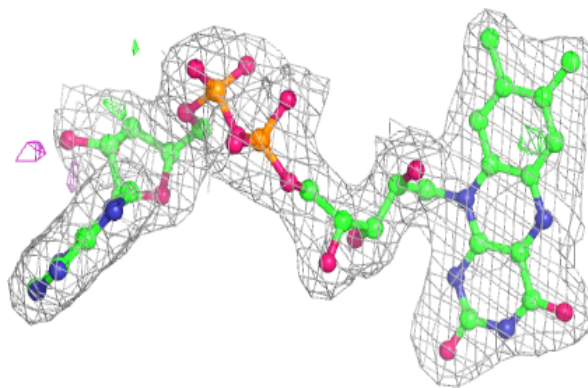
**Electron density around FAD B 401:**

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

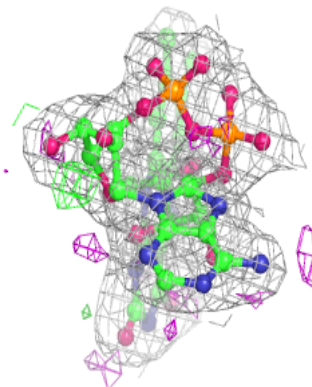
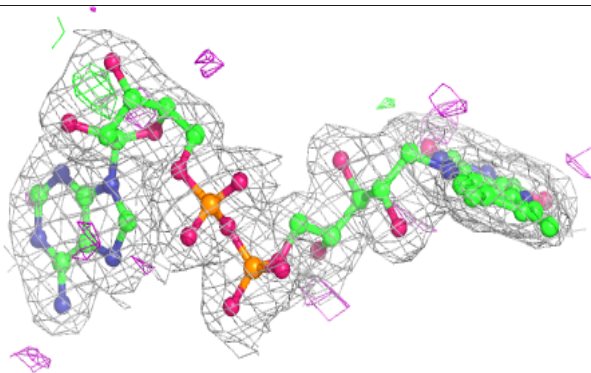
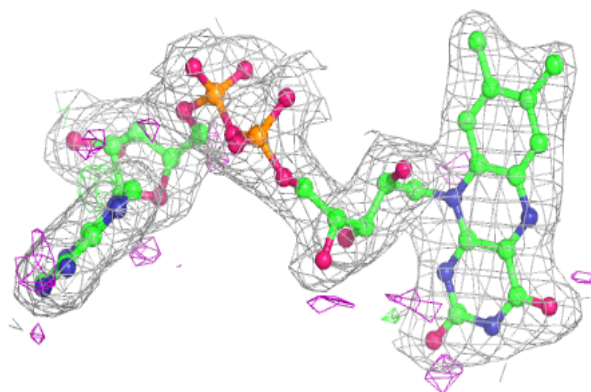


Electron density around FAD D 401:

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

**Electron density around FAD C 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.