



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 14, 2023 – 11:33 AM EDT

PDB ID : 4RPH  
Title : Crystal structure of Micobacterium tuberculosis UDP-Galactopyranose mutase in complex with substrate UDP-Galp (reduced)  
Authors : Van Straaten, K.E.; Sanders, D.A.R.  
Deposited on : 2014-10-30  
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

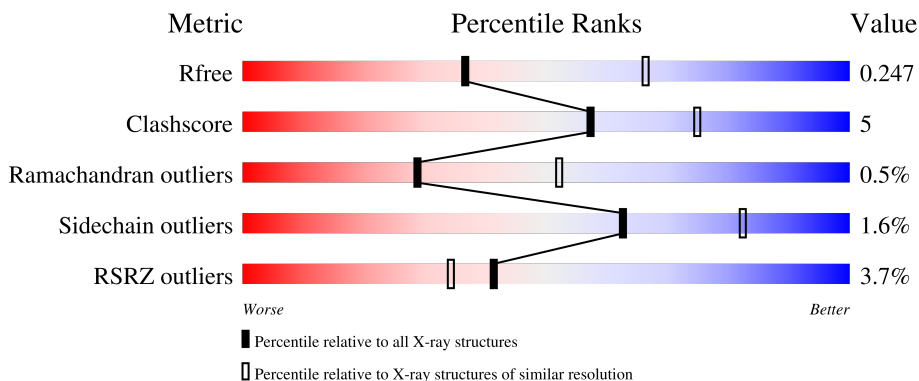
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


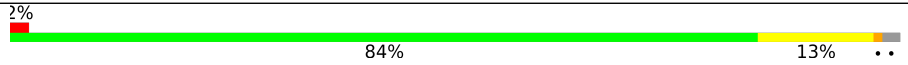
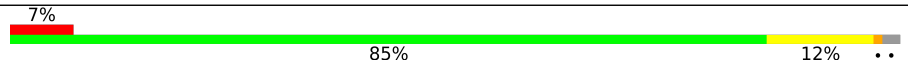
The reported resolution of this entry is 2.60 Å.

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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	399	 2% 88% 9% ..
1	B	399	 2% 84% 13% ..
1	C	399	 7% 85% 12% ..

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 9992 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.

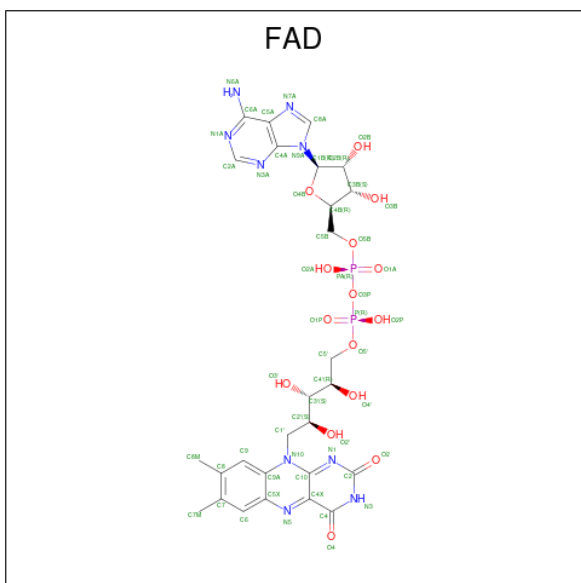
- Molecule 1 is a protein called UDP-galactopyranose mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	391	Total 3188	C 2034	N 556	O 590	S 8	0	0	0
1	A	392	Total 3208	C 2045	N 562	O 593	S 8	0	1	0
1	C	391	Total 3199	C 2040	N 560	O 591	S 8	0	1	0

There are 3 discrepancies between the modelled and reference sequences:

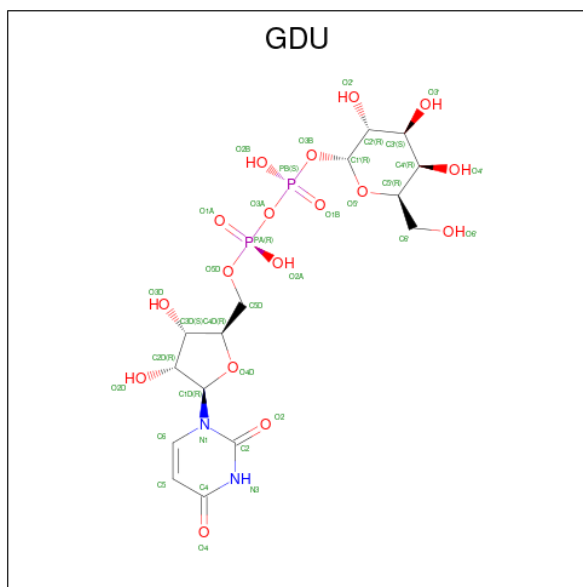
Chain	Residue	Modelled	Actual	Comment	Reference
B	306	ARG	PRO	engineered mutation	UNP P9WIK1
A	306	ARG	PRO	engineered mutation	UNP P9WIK1
C	306	ARG	PRO	engineered mutation	UNP P9WIK1

- 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	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	A	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		

- Molecule 3 is GALACTOSE-URIDINE-5'-DIPHOSPHATE (three-letter code: GDU) (formula:  $C_{15}H_{24}N_2O_{17}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	B	1	Total	C	N	O	P	0	0
			36	15	2	17	2		
3	A	1	Total	C	N	O	P	0	0
			36	15	2	17	2		
3	C	1	Total	C	N	O	P	0	0
			36	15	2	17	2		

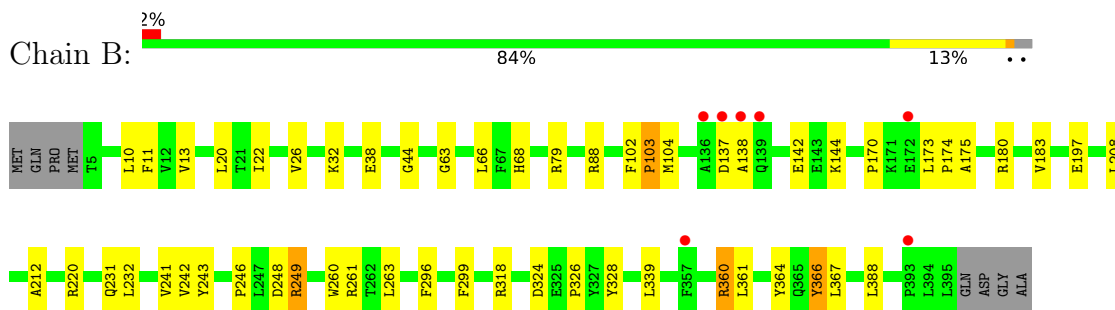
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	60	Total	O	0	0
			60	60		
4	A	38	Total	O	0	0
			38	38		
4	C	32	Total	O	0	0
			32	32		

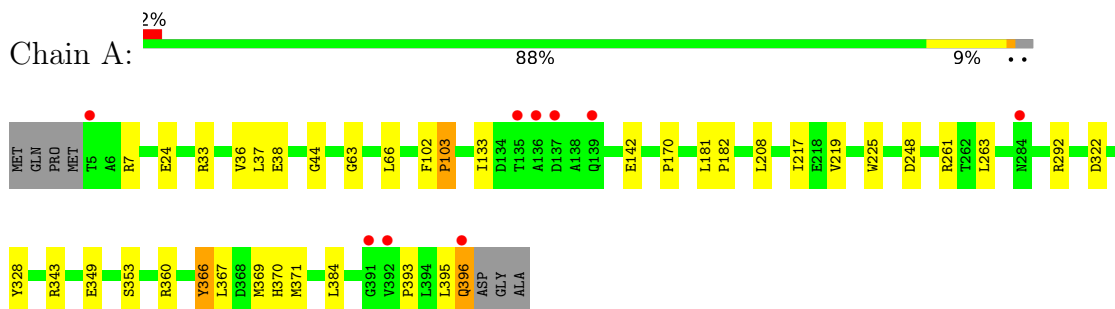
### 3 Residue-property plots [i](#)

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.

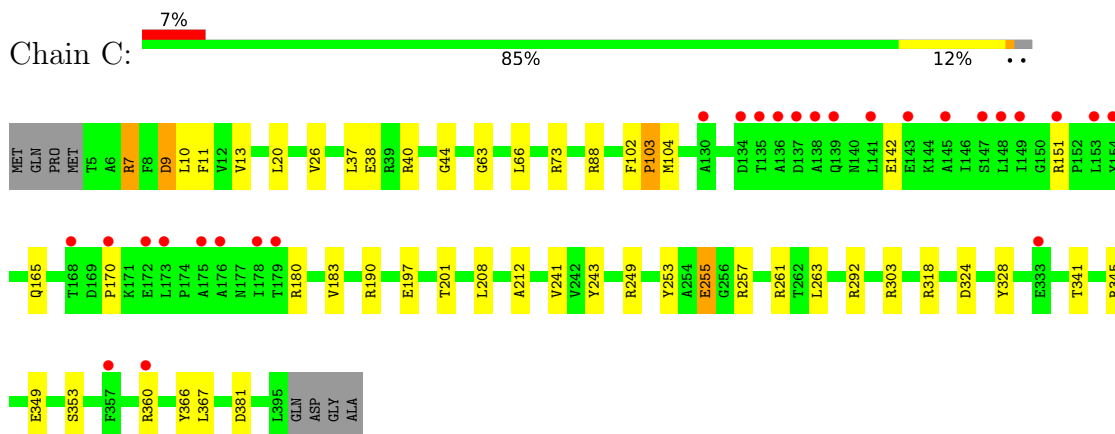
- Molecule 1: UDP-galactopyranose mutase



- Molecule 1: UDP-galactopyranose mutase



- Molecule 1: UDP-galactopyranose mutase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	171.73Å 99.47Å 100.24Å 90.00° 110.17° 90.00°	Depositor
Resolution (Å)	48.57 – 2.60 48.57 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (48.57-2.60) 98.8 (48.57-2.60)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 2.61Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.208 , 0.246 0.212 , 0.247	Depositor DCC
$R_{free}$ test set	2412 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	64.8	Xtrriage
Anisotropy	0.252	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 36.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9992	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.21	0/3297	0.37	0/4479
1	B	0.21	0/3277	0.38	0/4453
1	C	0.21	0/3288	0.37	0/4467
All	All	0.21	0/9862	0.37	0/13399

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3208	0	3067	29	0
1	B	3188	0	3047	31	0
1	C	3199	0	3059	29	0
2	A	53	0	30	8	0
2	B	53	0	30	6	0
2	C	53	0	30	5	0
3	A	36	0	22	4	0
3	B	36	0	22	3	0
3	C	36	0	22	4	0
4	A	38	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	60	0	0	2	0
4	C	32	0	0	0	0
All	All	9992	0	9329	94	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 5.

All (94) 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:63:GLY:HA3	2:A:401:FAD:HM72	1.66	0.76
1:C:63:GLY:HA3	2:C:401:FAD:HM72	1.70	0.73
1:A:142:GLU:HB2	1:A:170:PRO:HB2	1.75	0.69
1:B:63:GLY:HA3	2:B:401:FAD:HM72	1.79	0.65
1:C:104:MET:HG2	1:C:183:VAL:HG13	1.78	0.65
1:A:393:PRO:O	1:A:396:GLN:NE2	2.32	0.63
1:A:44:GLY:HA3	1:A:208:LEU:HD13	1.82	0.61
1:B:144:LYS:HD2	1:B:175:ALA:HB1	1.83	0.61
1:C:360[B]:ARG:NH2	2:C:401:FAD:O2A	2.34	0.59
1:A:360[A]:ARG:NH1	1:A:367:LEU:O	2.37	0.57
1:C:44:GLY:HA3	1:C:208:LEU:HD13	1.85	0.57
1:B:142:GLU:HB2	1:B:170:PRO:HB2	1.85	0.57
1:A:360[B]:ARG:NH2	2:A:401:FAD:O2A	2.37	0.57
1:C:73:ARG:NH2	1:C:381:ASP:OD2	2.38	0.56
1:B:44:GLY:HA3	1:B:208:LEU:HD13	1.87	0.56
1:C:263:LEU:HD11	1:C:328:TYR:HE2	1.72	0.55
1:C:201:THR:O	1:C:303:ARG:NH2	2.34	0.53
1:B:360:ARG:HD2	1:B:367:LEU:H	1.72	0.53
1:C:66:LEU:O	2:C:401:FAD:N3	2.38	0.53
1:C:360[A]:ARG:NH1	1:C:367:LEU:O	2.42	0.52
1:C:255:GLU:HG2	1:C:341:THR:HG21	1.92	0.52
1:B:38:GLU:OE2	2:B:401:FAD:O2B	2.27	0.52
1:B:366:TYR:OH	3:B:402:GDU:O2B	2.21	0.52
1:B:66:LEU:H	2:B:401:FAD:C4	2.23	0.52
1:A:248:ASP:OD1	1:A:248:ASP:N	2.43	0.52
1:C:102:PHE:CG	1:C:103:PRO:HA	2.45	0.52
1:B:102:PHE:CG	1:B:103:PRO:HA	2.45	0.51
1:C:318:ARG:NH1	1:C:324:ASP:OD2	2.42	0.51
1:A:102:PHE:CG	1:A:103:PRO:HA	2.46	0.51
2:A:401:FAD:N5	3:A:402:GDU:H2'	2.25	0.51
1:B:68:HIS:NE2	4:B:549:HOH:O	2.34	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LEU:O	2:B:401:FAD:N3	2.40	0.51
1:C:180:ARG:NH2	3:C:402:GDU:O2'	2.44	0.51
1:C:142:GLU:HB2	1:C:170:PRO:HB2	1.94	0.50
1:C:38:GLU:HG3	1:C:40:ARG:H	1.77	0.50
1:A:292:ARG:NH2	3:A:402:GDU:O1B	2.44	0.50
1:A:261:ARG:HB2	1:A:328:TYR:HB2	1.93	0.50
1:B:261:ARG:HB2	1:B:328:TYR:HB2	1.94	0.50
1:C:11:PHE:HB2	1:C:241:VAL:HG22	1.94	0.50
1:C:349:GLU:OE2	1:C:353:SER:OG	2.29	0.50
1:B:20:LEU:HD22	1:B:212:ALA:HB2	1.93	0.49
2:B:401:FAD:N5	3:B:402:GDU:H2'	2.27	0.49
1:A:366:TYR:OH	3:A:402:GDU:O2B	2.26	0.49
1:A:343:ARG:NH1	4:A:507:HOH:O	2.36	0.49
1:B:263:LEU:HD11	1:B:328:TYR:HE1	1.79	0.48
1:C:10:LEU:HD22	1:C:26:VAL:HG21	1.96	0.47
1:C:20:LEU:HD22	1:C:212:ALA:HB2	1.96	0.47
1:C:9:ASP:N	1:C:9:ASP:OD1	2.45	0.47
1:B:104:MET:HG2	1:B:183:VAL:HG13	1.96	0.47
1:B:326:PRO:HG2	1:B:361:LEU:HD22	1.97	0.46
1:B:318:ARG:NH1	1:B:324:ASP:OD2	2.37	0.46
1:A:263:LEU:HD11	1:A:328:TYR:HE1	1.81	0.46
1:B:88:ARG:NH1	1:B:197:GLU:O	2.48	0.46
1:B:248:ASP:N	1:B:248:ASP:OD1	2.47	0.46
1:A:322:ASP:OD1	1:A:322:ASP:N	2.49	0.46
1:C:292:ARG:NH2	3:C:402:GDU:O1B	2.48	0.46
1:B:13:VAL:HB	1:B:243:TYR:HA	1.98	0.45
1:A:66:LEU:H	2:A:401:FAD:C4	2.26	0.45
1:A:349:GLU:OE1	1:A:353:SER:OG	2.34	0.45
1:C:66:LEU:H	2:C:401:FAD:C4	2.28	0.45
1:A:44:GLY:HA2	1:A:208:LEU:HD22	1.98	0.45
1:B:180:ARG:NH2	3:B:402:GDU:O2'	2.49	0.45
1:A:36:VAL:HB	1:A:219:VAL:HG22	1.98	0.45
1:C:66:LEU:HD23	3:C:402:GDU:H6'2	1.99	0.45
1:C:88:ARG:NH1	1:C:197:GLU:O	2.50	0.45
1:B:220:ARG:HH12	1:B:232:LEU:HD21	1.82	0.45
1:A:38:GLU:OE2	2:A:401:FAD:O2B	2.30	0.44
1:A:66:LEU:O	2:A:401:FAD:N3	2.44	0.43
2:C:401:FAD:N5	3:C:402:GDU:H2'	2.34	0.43
1:B:10:LEU:HD22	1:B:26:VAL:HG21	2.01	0.42
1:B:246:PRO:HB2	1:B:249:ARG:HB3	2.01	0.42
1:A:181:LEU:HA	1:A:182:PRO:HD3	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:LYS:HE3	1:B:388:LEU:HB3	2.02	0.42
1:A:66:LEU:HD23	3:A:402:GDU:HG2	2.01	0.42
1:B:173:LEU:HA	1:B:174:PRO:HD3	1.94	0.42
2:A:401:FAD:H1'1	2:A:401:FAD:H9	1.79	0.42
2:B:401:FAD:O2	4:B:501:HOH:O	2.22	0.42
1:B:11:PHE:HB2	1:B:241:VAL:HG22	2.02	0.42
1:C:253:TYR:CD1	1:C:257:ARG:HG2	2.54	0.42
1:A:384:LEU:HD23	1:A:384:LEU:HA	1.90	0.41
1:A:7:ARG:NH1	1:A:33:ARG:HH11	2.19	0.41
1:B:339:LEU:HD22	1:B:364:TYR:HD2	1.85	0.41
1:B:296:PHE:HA	1:B:299:PHE:CD1	2.54	0.41
1:C:7:ARG:HA	1:C:7:ARG:HD2	1.84	0.41
1:C:88:ARG:HD2	1:C:190:ARG:CZ	2.51	0.41
1:A:37:LEU:HD13	1:A:225:TRP:HE3	1.86	0.40
1:A:369:MET:H	2:A:401:FAD:C2	2.34	0.40
1:C:261:ARG:HB2	1:C:328:TYR:HB2	2.02	0.40
1:B:260:TRP:CE3	1:B:326:PRO:HB2	2.56	0.40
1:A:24:GLU:HA	1:A:217:ILE:HD11	2.03	0.40
1:A:370:HIS:CE1	1:A:371:MET:HG3	2.57	0.40
1:A:133:ILE:HD12	1:A:133:ILE:HA	1.99	0.40
1:C:13:VAL:HB	1:C:243:TYR:HA	2.03	0.40
1:B:22:ILE:HG21	1:B:242:VAL:HG21	2.03	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	391/399 (98%)	381 (97%)	9 (2%)	1 (0%)	41 64
1	B	389/399 (98%)	373 (96%)	13 (3%)	3 (1%)	19 39
1	C	390/399 (98%)	374 (96%)	14 (4%)	2 (0%)	29 52

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1170/1197 (98%)	1128 (96%)	36 (3%)	6 (0%)	29	52

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	138	ALA
1	B	137	ASP
1	C	103	PRO
1	C	151	ARG
1	B	103	PRO
1	A	103	PRO

### 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	330/334 (99%)	327 (99%)	3 (1%)	78	91
1	B	328/334 (98%)	323 (98%)	5 (2%)	65	83
1	C	329/334 (98%)	321 (98%)	8 (2%)	49	74
All	All	987/1002 (98%)	971 (98%)	16 (2%)	62	82

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

Mol	Chain	Res	Type
1	B	79	ARG
1	B	231	GLN
1	B	249	ARG
1	B	360	ARG
1	B	366	TYR
1	A	366	TYR
1	A	395	LEU
1	A	396	GLN
1	C	7	ARG
1	C	9	ASP

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Mol	Chain	Res	Type
1	C	37	LEU
1	C	165	GLN
1	C	249	ARG
1	C	255	GLU
1	C	345	ARG
1	C	366	TYR

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

Mol	Chain	Res	Type
1	A	396	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	GDU	B	402	-	35,38,38	1.94	11 (31%)	53,58,58	1.74	11 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	FAD	C	401	-	53,58,58	2.02	14 (26%)	68,89,89	1.64	15 (22%)
2	FAD	A	401	-	53,58,58	2.02	14 (26%)	68,89,89	1.65	14 (20%)
2	FAD	B	401	-	53,58,58	2.02	14 (26%)	68,89,89	1.65	16 (23%)
3	GDU	C	402	-	35,38,38	1.93	11 (31%)	53,58,58	1.78	12 (22%)
3	GDU	A	402	-	35,38,38	1.93	11 (31%)	53,58,58	1.78	11 (20%)

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	GDU	B	402	-	-	4/23/59/59	0/3/3/3
2	FAD	C	401	-	-	6/30/50/50	0/6/6/6
2	FAD	A	401	-	-	7/30/50/50	0/6/6/6
2	FAD	B	401	-	-	8/30/50/50	0/6/6/6
3	GDU	C	402	-	-	4/23/59/59	0/3/3/3
3	GDU	A	402	-	-	4/23/59/59	0/3/3/3

All (75) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FAD	C10-N1	5.35	1.44	1.33
2	C	401	FAD	C10-N1	5.27	1.44	1.33
2	B	401	FAD	C10-N1	5.25	1.44	1.33
2	C	401	FAD	C1'-C2'	5.14	1.59	1.52
2	B	401	FAD	C2B-C3B	-5.14	1.39	1.53
2	A	401	FAD	C2B-C3B	-5.10	1.39	1.53
2	C	401	FAD	C2B-C3B	-5.09	1.39	1.53
2	A	401	FAD	C1'-C2'	5.01	1.59	1.52
2	B	401	FAD	C1'-C2'	4.81	1.59	1.52
3	A	402	GDU	O4-C4	4.70	1.33	1.24
3	B	402	GDU	O4-C4	4.70	1.33	1.24
3	C	402	GDU	O4-C4	4.69	1.33	1.24
3	A	402	GDU	PB-O3B	-4.06	1.49	1.60
3	C	402	GDU	PB-O3B	-4.05	1.49	1.60
3	B	402	GDU	PB-O3B	-4.04	1.49	1.60
2	B	401	FAD	C7M-C7	3.91	1.58	1.51
2	A	401	FAD	C7M-C7	3.84	1.58	1.51
2	C	401	FAD	C7M-C7	3.78	1.58	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	GDU	C3D-C2D	-3.50	1.43	1.53
3	A	402	GDU	C3D-C2D	-3.46	1.43	1.53
2	B	401	FAD	C2-N1	3.44	1.44	1.36
2	C	401	FAD	C2-N1	3.42	1.44	1.36
3	C	402	GDU	C3D-C2D	-3.40	1.44	1.53
2	B	401	FAD	C9A-C5X	-3.36	1.35	1.41
2	A	401	FAD	C2-N1	3.32	1.44	1.36
3	C	402	GDU	C4'-C3'	-3.22	1.44	1.52
3	B	402	GDU	C4'-C3'	-3.19	1.44	1.52
2	A	401	FAD	C9A-C5X	-3.17	1.35	1.41
2	C	401	FAD	C9A-C5X	-3.14	1.36	1.41
3	A	402	GDU	C4'-C3'	-3.13	1.44	1.52
2	B	401	FAD	O3'-C3'	-3.11	1.35	1.43
2	C	401	FAD	O3'-C3'	-3.07	1.35	1.43
2	A	401	FAD	O3'-C3'	-3.05	1.35	1.43
2	C	401	FAD	C9A-N10	-2.96	1.35	1.41
3	B	402	GDU	C3D-C4D	-2.88	1.45	1.53
2	A	401	FAD	C9A-N10	-2.85	1.36	1.41
3	C	402	GDU	C3D-C4D	-2.85	1.45	1.53
3	A	402	GDU	C3D-C4D	-2.83	1.45	1.53
2	B	401	FAD	C9A-N10	-2.81	1.36	1.41
2	C	401	FAD	O4'-C4'	2.80	1.49	1.43
2	A	401	FAD	O4'-C4'	2.78	1.49	1.43
2	A	401	FAD	C4X-N5	2.77	1.36	1.30
2	B	401	FAD	O4'-C4'	2.74	1.49	1.43
3	B	402	GDU	C4-N3	-2.72	1.33	1.38
3	C	402	GDU	C4-N3	-2.72	1.33	1.38
3	A	402	GDU	C4-N3	-2.70	1.33	1.38
3	B	402	GDU	C2-N1	-2.68	1.34	1.38
3	C	402	GDU	C2-N1	-2.59	1.34	1.38
2	C	401	FAD	P-O1P	2.59	1.60	1.50
2	B	401	FAD	P-O1P	2.58	1.60	1.50
2	C	401	FAD	C4X-N5	2.58	1.35	1.30
2	B	401	FAD	C4'-C3'	-2.57	1.48	1.53
2	A	401	FAD	P-O1P	2.57	1.60	1.50
2	A	401	FAD	C1'-N10	-2.53	1.41	1.48
3	A	402	GDU	C2-N1	-2.51	1.34	1.38
2	B	401	FAD	C1'-N10	-2.49	1.41	1.48
2	B	401	FAD	C4X-N5	2.48	1.35	1.30
2	C	401	FAD	C4'-C3'	-2.45	1.48	1.53
2	A	401	FAD	C4'-C3'	-2.43	1.48	1.53
2	C	401	FAD	C1'-N10	-2.43	1.42	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	FAD	C5X-N5	-2.42	1.34	1.39
3	C	402	GDU	C3'-C2'	-2.40	1.46	1.52
3	A	402	GDU	C3'-C2'	-2.36	1.46	1.52
3	B	402	GDU	C3'-C2'	-2.31	1.46	1.52
2	C	401	FAD	C5X-N5	-2.24	1.35	1.39
2	A	401	FAD	C5X-N5	-2.23	1.35	1.39
3	B	402	GDU	C5-C4	-2.17	1.38	1.43
3	A	402	GDU	C6-C5	2.17	1.40	1.35
3	B	402	GDU	C6-C5	2.16	1.40	1.35
3	C	402	GDU	O5'-C5'	2.16	1.49	1.44
3	C	402	GDU	C6-C5	2.15	1.40	1.35
3	A	402	GDU	C5-C4	-2.14	1.38	1.43
3	C	402	GDU	C5-C4	-2.14	1.38	1.43
3	B	402	GDU	O5'-C5'	2.10	1.49	1.44
3	A	402	GDU	O5'-C5'	2.10	1.49	1.44

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	402	GDU	N3-C2-N1	5.92	122.75	114.89
3	A	402	GDU	N3-C2-N1	5.90	122.72	114.89
3	B	402	GDU	N3-C2-N1	5.81	122.60	114.89
3	C	402	GDU	C4-N3-C2	-5.17	119.77	126.58
3	A	402	GDU	C4-N3-C2	-5.07	119.89	126.58
3	B	402	GDU	C4-N3-C2	-4.98	120.01	126.58
2	B	401	FAD	O5B-PA-O1A	-4.83	90.19	109.07
2	A	401	FAD	O5B-PA-O1A	-4.77	90.45	109.07
2	C	401	FAD	O5B-PA-O1A	-4.70	90.69	109.07
2	A	401	FAD	N3A-C2A-N1A	-4.48	121.67	128.68
2	B	401	FAD	N3A-C2A-N1A	-4.41	121.78	128.68
2	C	401	FAD	N3A-C2A-N1A	-4.40	121.81	128.68
3	A	402	GDU	PB-O3A-PA	-4.09	118.79	132.83
3	C	402	GDU	PB-O3A-PA	-4.07	118.87	132.83
3	B	402	GDU	PB-O3A-PA	-4.06	118.91	132.83
2	C	401	FAD	C1'-C2'-C3'	3.71	120.17	109.79
2	A	401	FAD	C1'-C2'-C3'	3.69	120.09	109.79
2	B	401	FAD	O5'-P-O1P	-3.36	95.93	109.07
2	B	401	FAD	C4-N3-C2	-3.35	119.45	125.64
2	C	401	FAD	O5'-P-O1P	-3.34	96.02	109.07
2	B	401	FAD	O5'-C5'-C4'	3.34	118.27	109.36
2	A	401	FAD	O5'-P-O1P	-3.31	96.12	109.07
2	A	401	FAD	O5'-C5'-C4'	3.28	118.11	109.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	FAD	C1'-C2'-C3'	3.26	118.91	109.79
2	C	401	FAD	O5'-C5'-C4'	3.23	117.97	109.36
2	A	401	FAD	C4-N3-C2	-3.17	119.78	125.64
3	C	402	GDU	C5-C4-N3	3.17	119.58	114.84
3	A	402	GDU	C5-C4-N3	3.13	119.52	114.84
3	B	402	GDU	C5-C4-N3	3.10	119.47	114.84
2	C	401	FAD	C4-N3-C2	-3.09	119.94	125.64
3	C	402	GDU	O2-C2-N1	-3.05	118.73	122.79
3	B	402	GDU	O2-C2-N1	-2.94	118.88	122.79
3	A	402	GDU	C3'-C4'-C5'	2.90	115.42	110.24
3	A	402	GDU	O2-C2-N1	-2.90	118.93	122.79
3	A	402	GDU	O5'-C5'-C4'	2.86	114.89	109.69
3	C	402	GDU	O5'-C5'-C4'	2.81	114.80	109.69
2	A	401	FAD	C4X-C4-N3	2.76	120.20	113.19
3	C	402	GDU	C3'-C4'-C5'	2.75	115.14	110.24
3	C	402	GDU	O4-C4-C5	-2.75	120.33	125.16
2	C	401	FAD	O5B-C5B-C4B	2.74	118.44	108.99
3	A	402	GDU	O4-C4-C5	-2.73	120.35	125.16
2	B	401	FAD	O5B-C5B-C4B	2.72	118.36	108.99
3	B	402	GDU	O4-C4-C5	-2.70	120.42	125.16
3	B	402	GDU	O5'-C5'-C4'	2.67	114.53	109.69
2	B	401	FAD	C4X-C4-N3	2.64	119.90	113.19
2	A	401	FAD	O5B-C5B-C4B	2.64	118.08	108.99
3	B	402	GDU	C3'-C4'-C5'	2.63	114.93	110.24
2	C	401	FAD	C4X-C4-N3	2.63	119.87	113.19
2	B	401	FAD	O4-C4-C4X	-2.58	119.75	126.60
2	C	401	FAD	O4-C4-C4X	-2.55	119.83	126.60
2	A	401	FAD	O4-C4-C4X	-2.52	119.92	126.60
3	B	402	GDU	O3A-PB-O3B	2.43	107.38	102.48
2	A	401	FAD	C3B-C2B-C1B	2.37	104.55	100.98
2	B	401	FAD	C4-C4X-C10	2.31	120.67	116.79
2	A	401	FAD	C10-C4X-N5	-2.27	120.03	124.86
3	A	402	GDU	C3D-C2D-C1D	2.26	105.72	101.43
2	C	401	FAD	C10-C4X-N5	-2.25	120.08	124.86
2	B	401	FAD	O2P-P-O5'	2.25	118.19	107.75
2	C	401	FAD	O2P-P-O5'	2.25	118.19	107.75
2	A	401	FAD	O2P-P-O5'	2.25	118.18	107.75
2	B	401	FAD	O2A-PA-O5B	2.23	118.12	107.75
3	A	402	GDU	C1D-N1-C2	2.21	121.58	117.57
2	A	401	FAD	O2A-PA-O5B	2.21	118.00	107.75
2	C	401	FAD	C3B-C2B-C1B	2.20	104.29	100.98
2	A	401	FAD	C2B-C3B-C4B	2.17	106.86	102.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	FAD	C4X-C10-N1	-2.16	119.72	124.73
2	B	401	FAD	C10-C4X-N5	-2.16	120.28	124.86
3	C	402	GDU	C3D-C2D-C1D	2.15	105.52	101.43
3	C	402	GDU	O3A-PB-O3B	2.15	106.83	102.48
2	C	401	FAD	C2B-C3B-C4B	2.15	106.82	102.64
3	C	402	GDU	C1D-N1-C2	2.14	121.45	117.57
3	B	402	GDU	O5D-C5D-C4D	2.12	116.28	108.99
2	B	401	FAD	C2B-C3B-C4B	2.10	106.72	102.64
3	C	402	GDU	O5D-C5D-C4D	2.06	116.08	108.99
3	B	402	GDU	C1D-N1-C2	2.06	121.30	117.57
2	B	401	FAD	C5X-C9A-N10	2.03	120.05	117.95
3	A	402	GDU	O3A-PB-O3B	2.02	106.56	102.48
2	C	401	FAD	O2A-PA-O5B	2.02	117.11	107.75
2	C	401	FAD	C4X-C10-N1	-2.00	120.08	124.73

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	FAD	O4'-C4'-C5'-O5'
2	A	401	FAD	O4B-C4B-C5B-O5B
2	A	401	FAD	C3'-C4'-C5'-O5'
2	A	401	FAD	O4'-C4'-C5'-O5'
2	C	401	FAD	O4B-C4B-C5B-O5B
2	C	401	FAD	C3'-C4'-C5'-O5'
2	C	401	FAD	O4'-C4'-C5'-O5'
3	B	402	GDU	PA-O3A-PB-O3B
3	B	402	GDU	O5'-C1'-O3B-PB
3	A	402	GDU	PA-O3A-PB-O3B
3	C	402	GDU	PA-O3A-PB-O3B
2	B	401	FAD	O4B-C4B-C5B-O5B
3	A	402	GDU	O5'-C5'-C6'-O6'
3	C	402	GDU	O5'-C5'-C6'-O6'
3	B	402	GDU	O5'-C5'-C6'-O6'
3	A	402	GDU	C4'-C5'-C6'-O6'
3	B	402	GDU	C4'-C5'-C6'-O6'
3	C	402	GDU	C4'-C5'-C6'-O6'
2	B	401	FAD	C3B-C4B-C5B-O5B
2	A	401	FAD	C3B-C4B-C5B-O5B
2	C	401	FAD	C3B-C4B-C5B-O5B
2	B	401	FAD	C3'-C4'-C5'-O5'
2	B	401	FAD	C5B-O5B-PA-O3P

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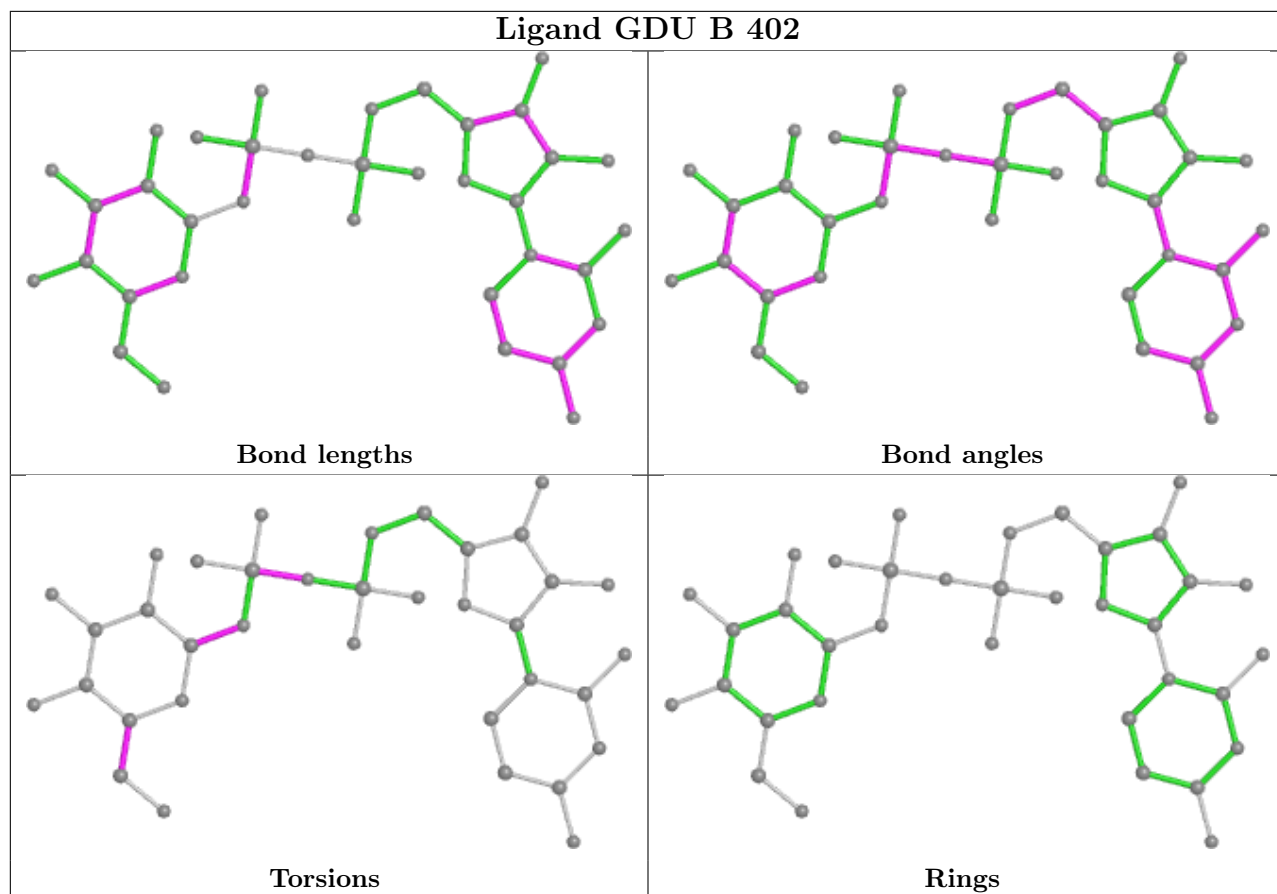
Mol	Chain	Res	Type	Atoms
2	A	401	FAD	C5B-O5B-PA-O3P
2	C	401	FAD	C5B-O5B-PA-O3P
2	C	401	FAD	C5'-O5'-P-O3P
2	B	401	FAD	C2'-C3'-C4'-C5'
2	B	401	FAD	O3'-C3'-C4'-C5'
3	A	402	GDU	O5'-C1'-O3B-PB
3	C	402	GDU	O5'-C1'-O3B-PB
2	B	401	FAD	C4'-C5'-O5'-P
2	A	401	FAD	O3'-C3'-C4'-C5'
2	A	401	FAD	C5'-O5'-P-O3P

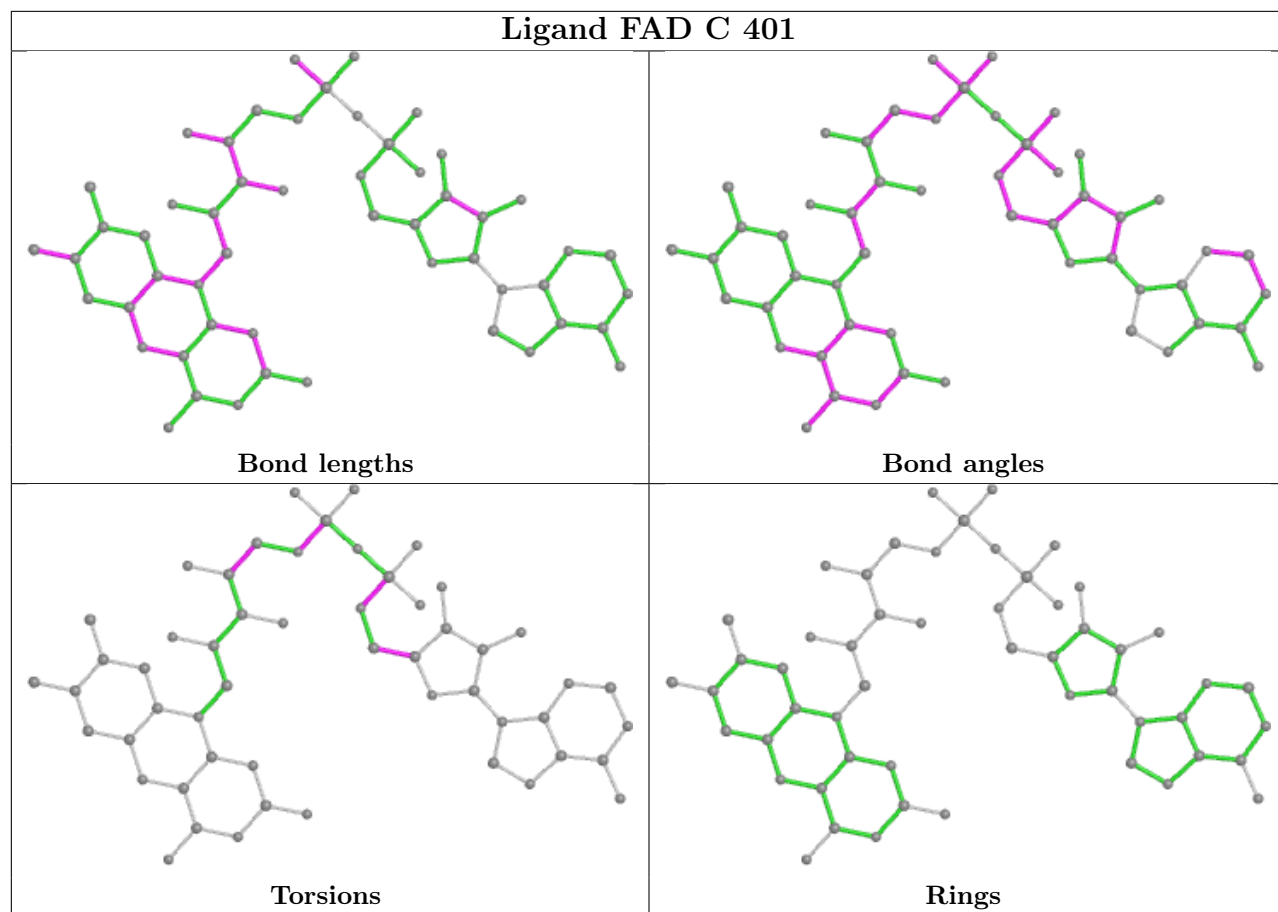
There are no ring outliers.

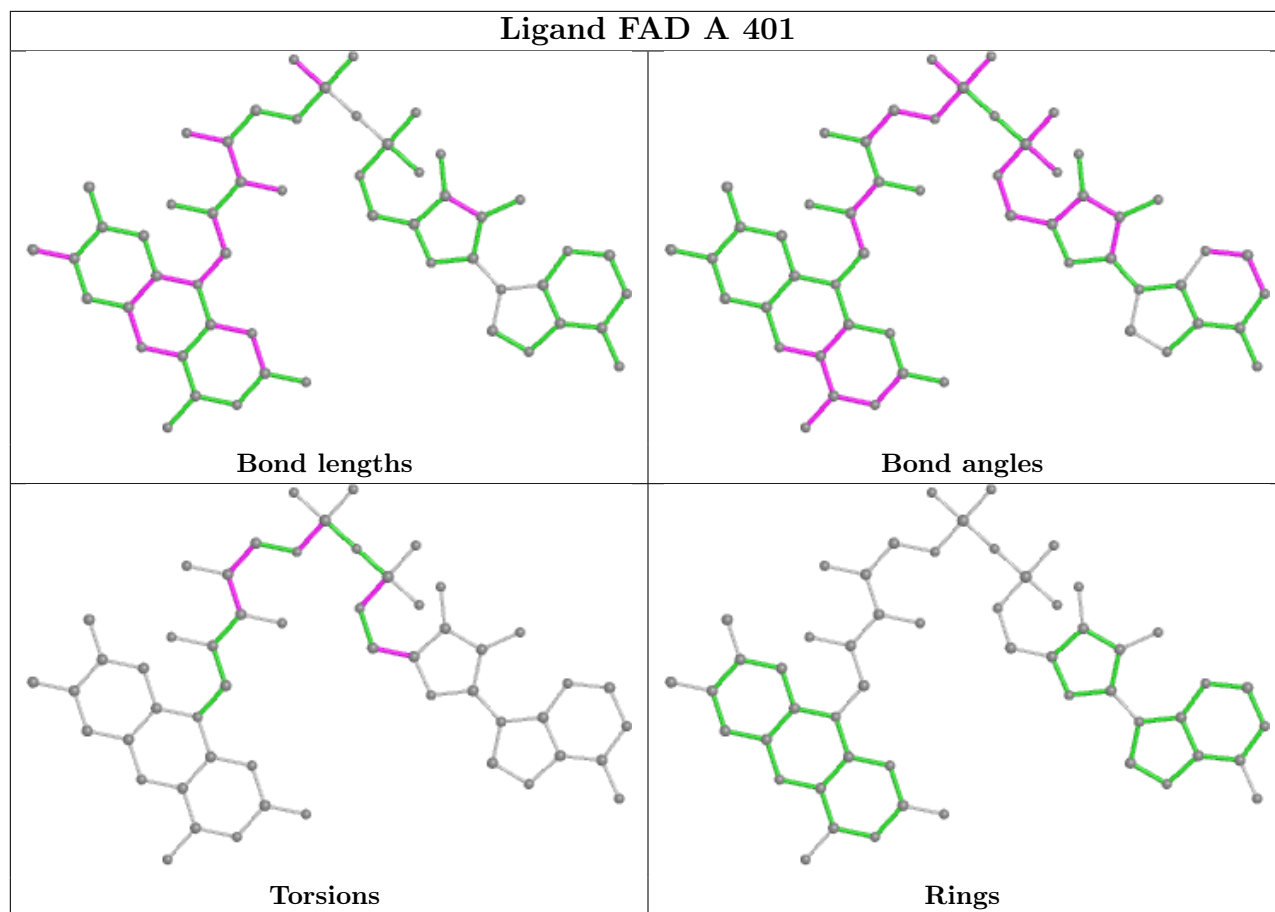
6 monomers are involved in 27 short contacts:

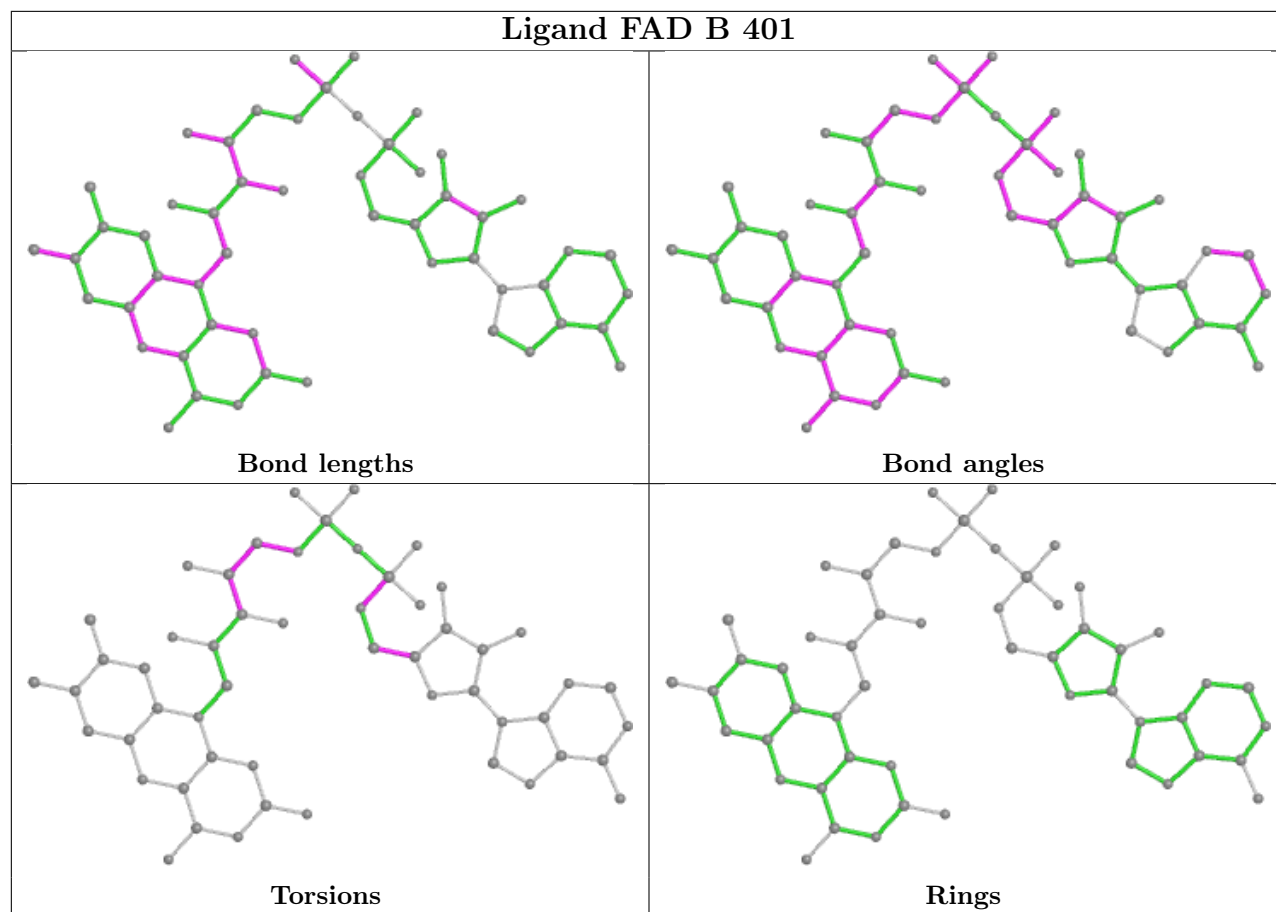
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	GDU	3	0
2	C	401	FAD	5	0
2	A	401	FAD	8	0
2	B	401	FAD	6	0
3	C	402	GDU	4	0
3	A	402	GDU	4	0

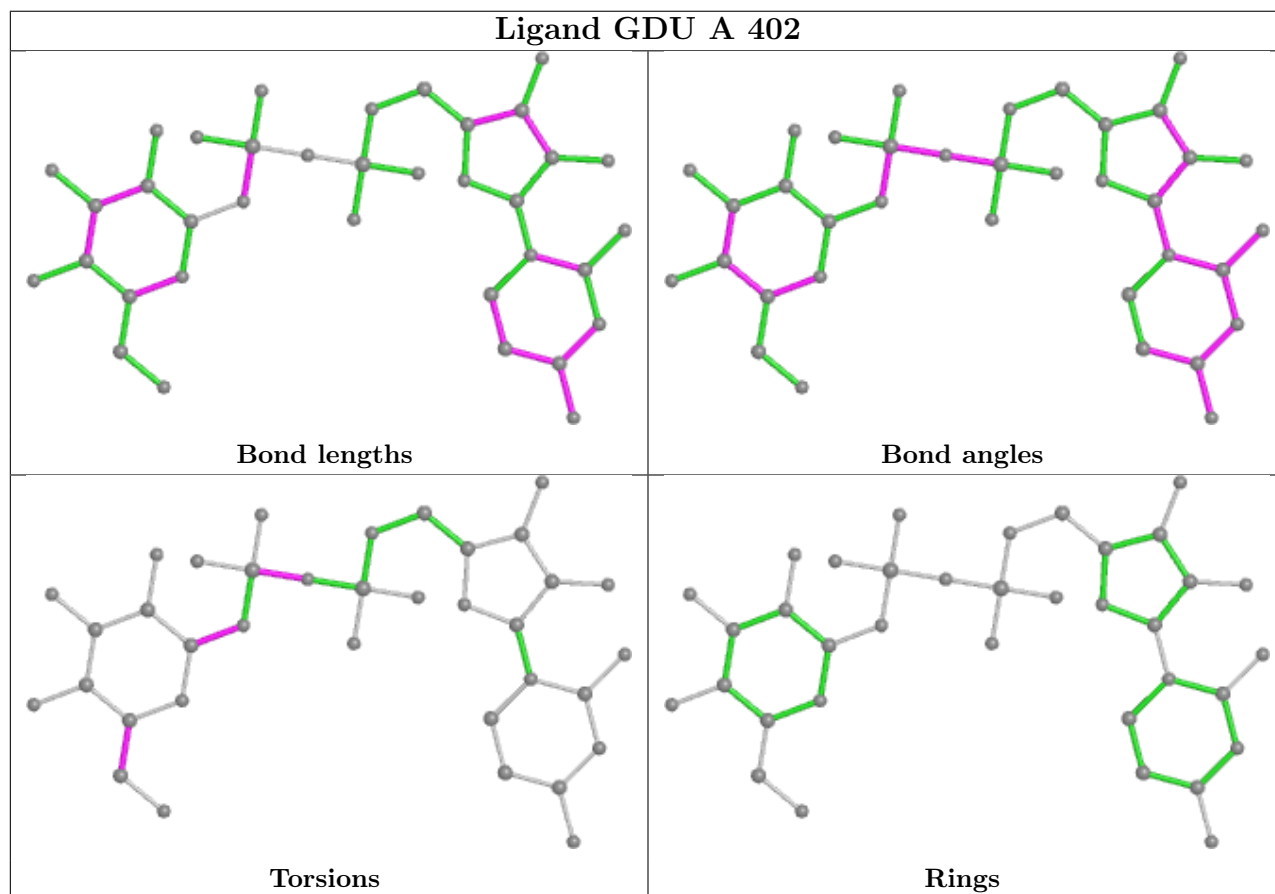
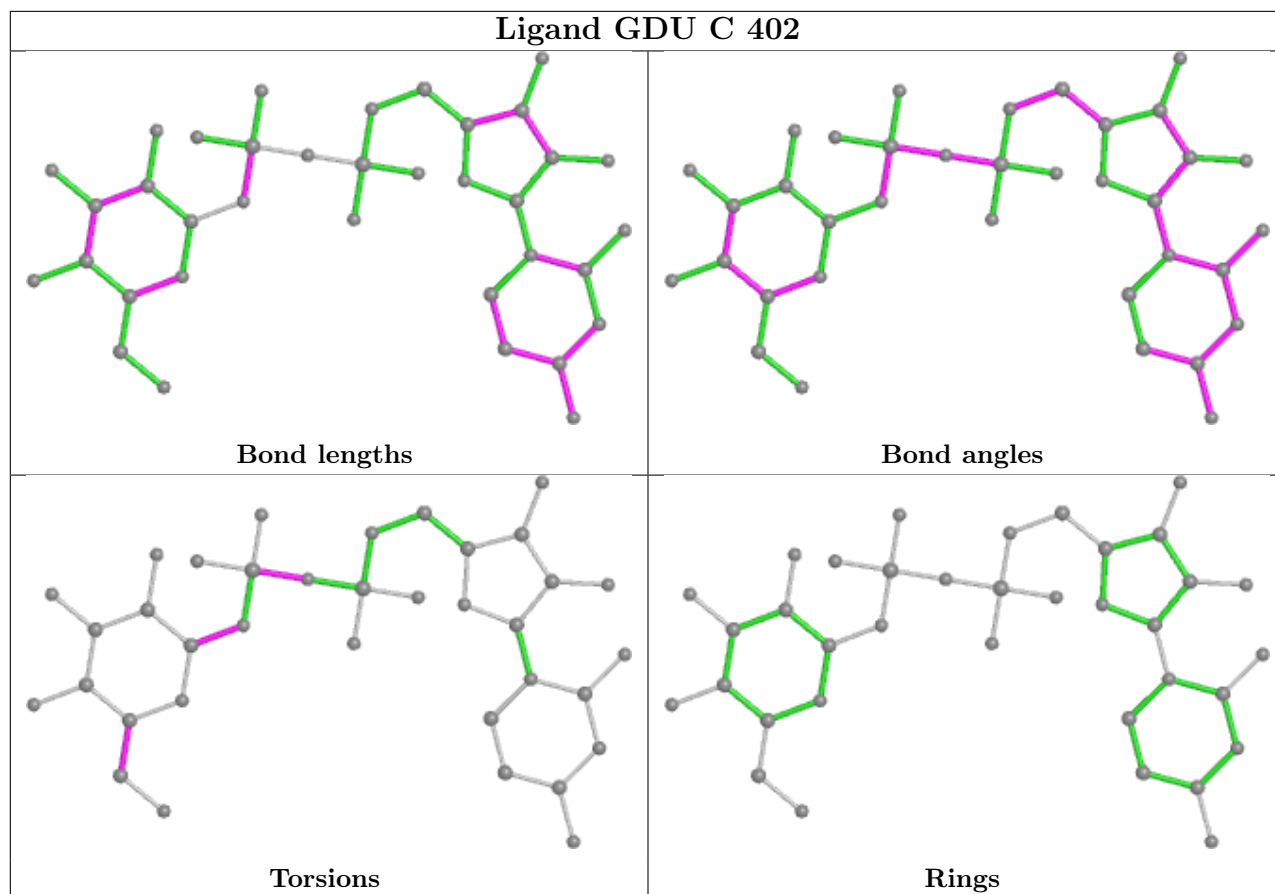
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	392/399 (98%)	0.19	9 (2%) 60 54	42, 61, 85, 108	0
1	B	391/399 (97%)	0.12	7 (1%) 68 64	41, 59, 83, 103	0
1	C	391/399 (97%)	0.49	27 (6%) 16 12	45, 69, 107, 130	0
All	All	1174/1197 (98%)	0.27	43 (3%) 41 34	41, 62, 94, 130	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	135	THR	7.0
1	C	136	ALA	5.4
1	C	149	ILE	5.3
1	C	143	GLU	5.0
1	C	137	ASP	4.8
1	C	134	ASP	4.7
1	C	154	TYR	4.6
1	C	139	GLN	4.4
1	A	137	ASP	3.9
1	C	153	LEU	3.8
1	C	145	ALA	3.7
1	C	170	PRO	3.3
1	C	172	GLU	3.2
1	C	179	THR	3.1
1	A	135	THR	3.1
1	C	147	SER	3.1
1	C	333	GLU	3.1
1	C	138	ALA	3.1
1	C	141	LEU	3.0
1	C	178	ILE	2.8
1	C	357	PHE	2.8
1	C	151	ARG	2.8
1	B	136	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	130	ALA	2.7
1	B	138	ALA	2.7
1	C	175	ALA	2.6
1	C	176	ALA	2.6
1	B	137	ASP	2.5
1	A	396	GLN	2.5
1	B	172	GLU	2.5
1	A	284	ASN	2.5
1	C	168	THR	2.4
1	C	148	LEU	2.4
1	A	136	ALA	2.3
1	A	5	THR	2.2
1	C	360[A]	ARG	2.2
1	A	391	GLY	2.2
1	A	139	GLN	2.1
1	B	357	PHE	2.1
1	B	139	GLN	2.1
1	B	393	PRO	2.0
1	A	392	VAL	2.0
1	C	173	LEU	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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

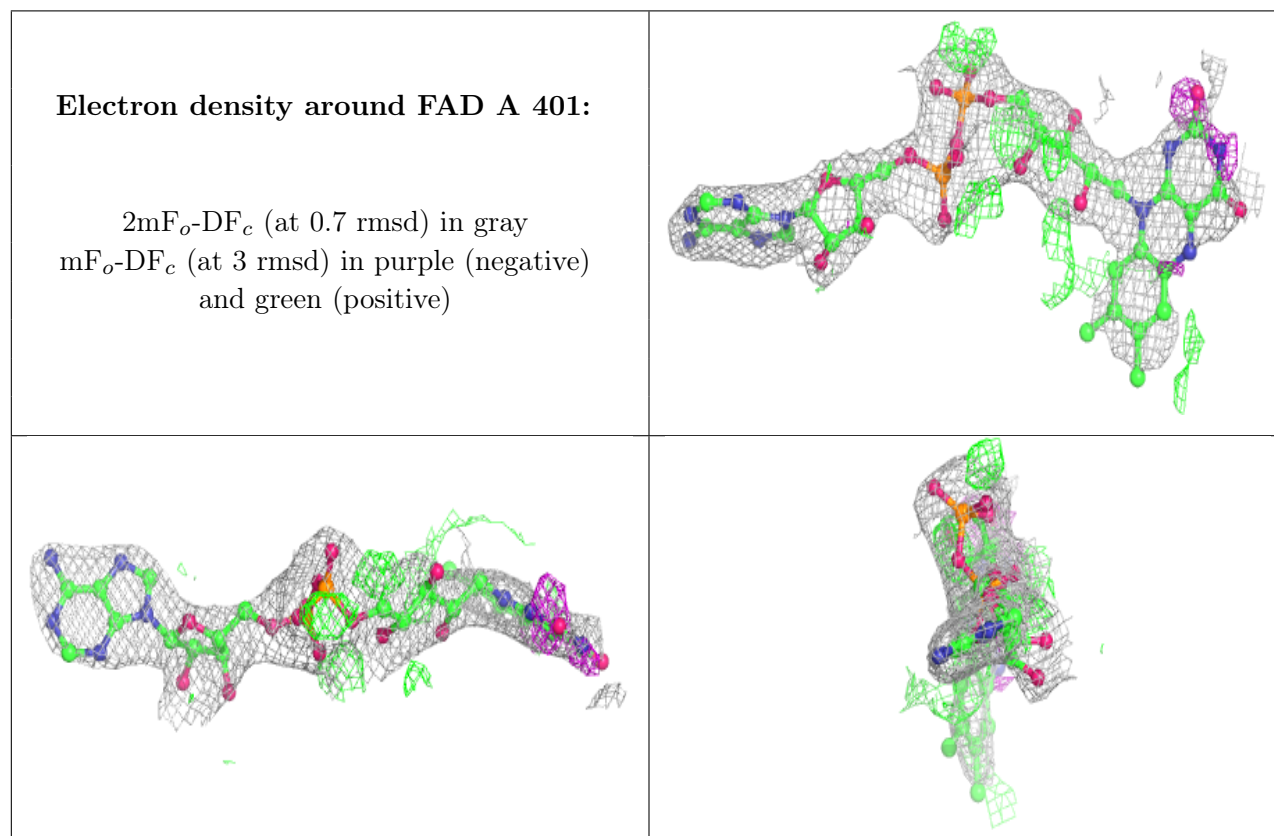
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAD	A	401	53/53	0.86	0.28	53,65,71,73	53
2	FAD	C	401	53/53	0.87	0.30	53,77,83,86	53
2	FAD	B	401	53/53	0.88	0.32	49,65,71,74	53

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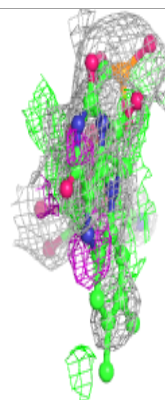
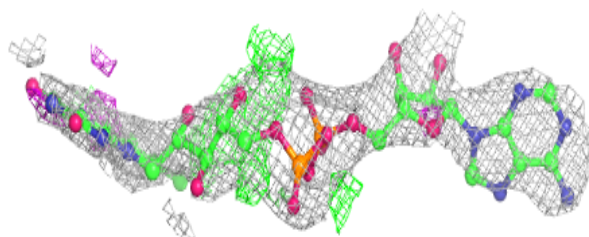
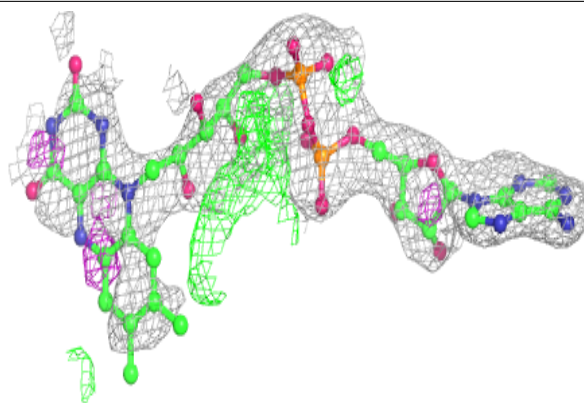
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GDU	C	402	36/36	0.92	0.20	68,80,95,97	9
3	GDU	A	402	36/36	0.97	0.18	51,62,69,72	9
3	GDU	B	402	36/36	0.97	0.18	44,55,61,62	9

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.

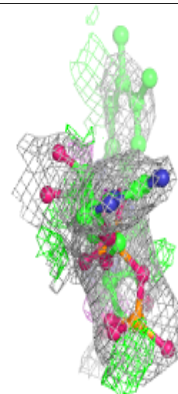
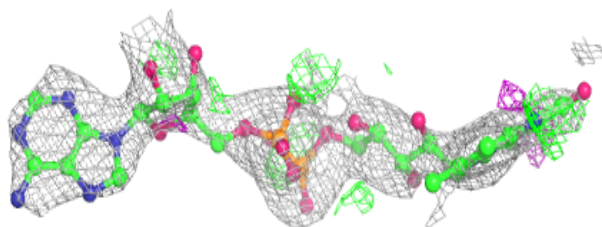
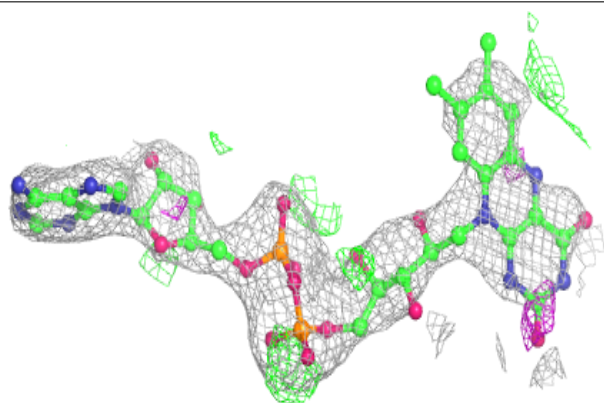


**Electron density around FAD C 401:**

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

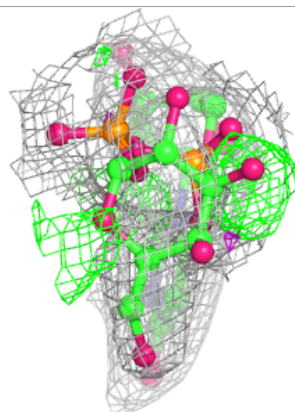
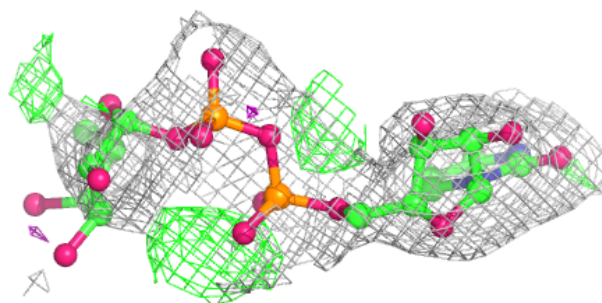
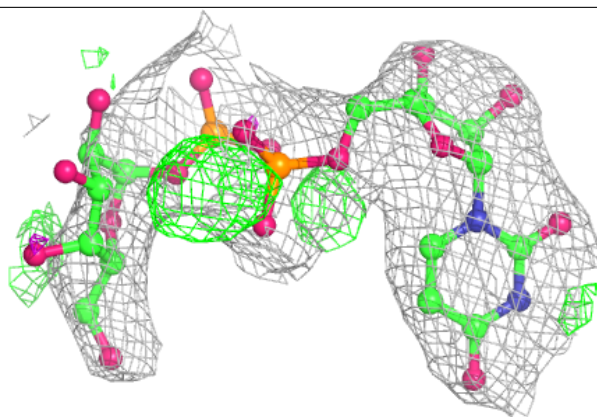
**Electron density around FAD B 401:**

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

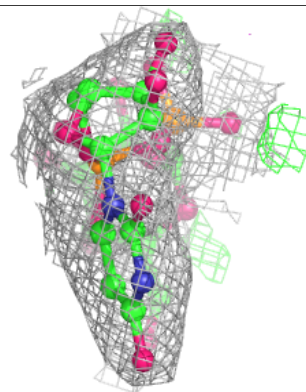
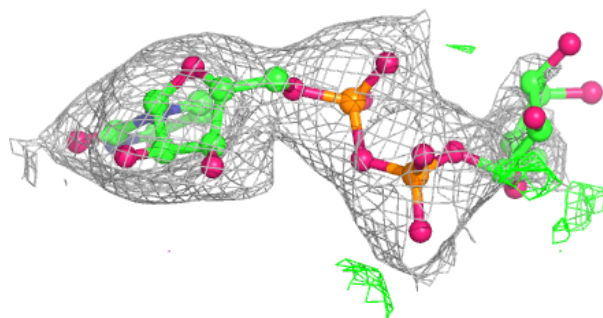
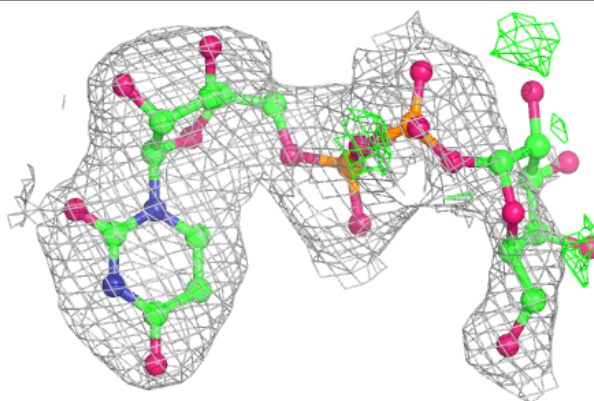


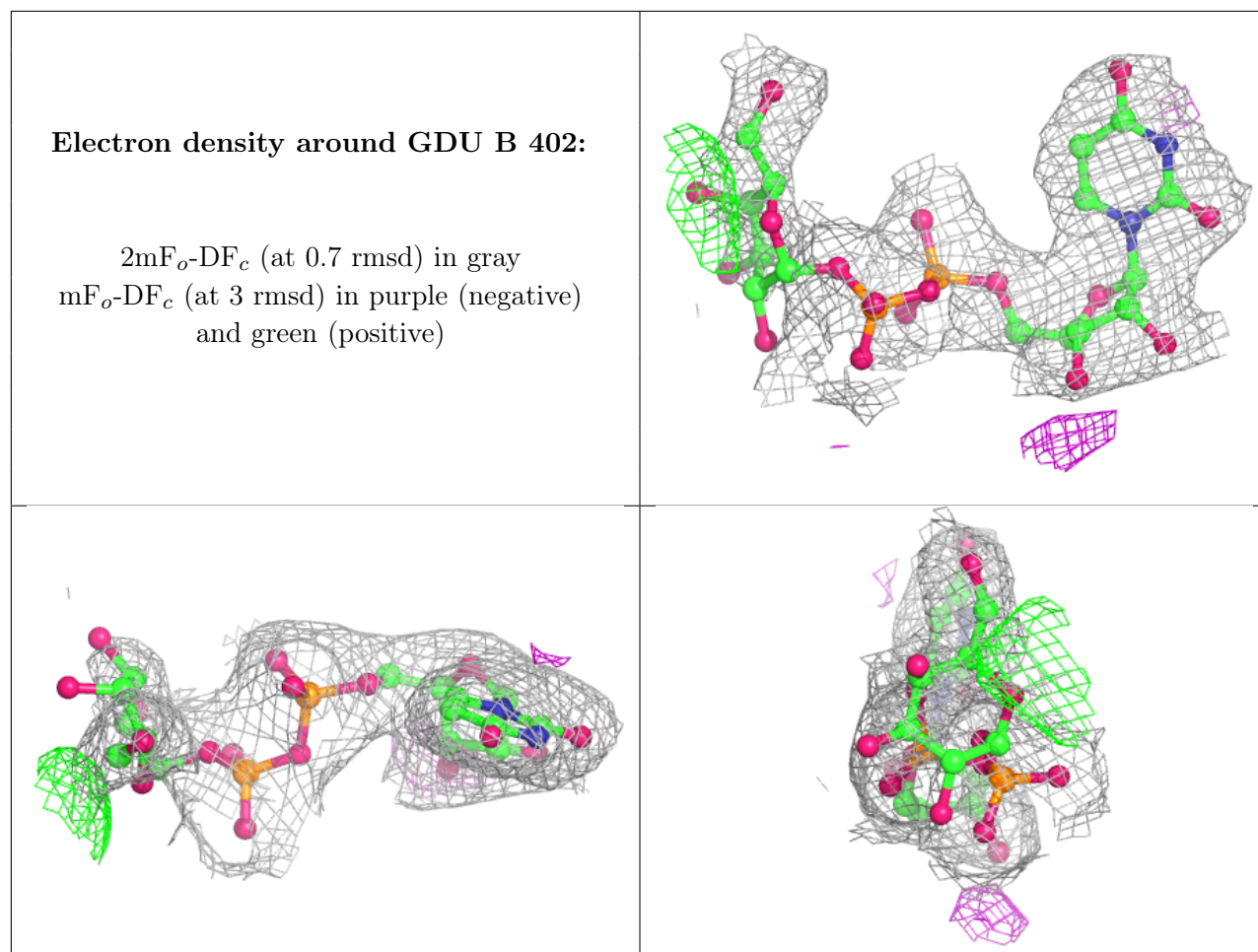
**Electron density around GDU C 402:**

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

**Electron density around GDU A 402:**

$2mF_o-DF_c$  (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.