



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:29 AM GMT

PDB ID : 3INR  
Title : Structure of UDP-galactopyranose mutase bound to UDP-galactose (oxidized)  
Authors : Gruber, T.D.; Kiessling, L.L.; Forest, K.T.  
Deposited on : 2009-08-12  
Resolution : 2.30 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

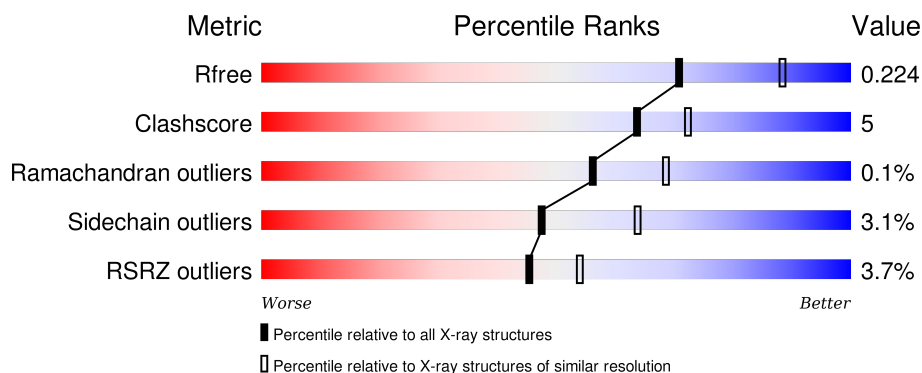
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	390	<div> <div>3%</div> <div>85%</div> <div>12%</div> <div>..</div> </div>
1	B	390	<div> <div>4%</div> <div>85%</div> <div>10%</div> <div>..</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6666 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
1	A	383	Total	C	N	O	S	0	0	0
			3119	1994	525	583	17			
1	B	378	Total	C	N	O	S	0	0	0
			3082	1971	519	575	17			

There are 22 discrepancies between the modelled and reference sequences:

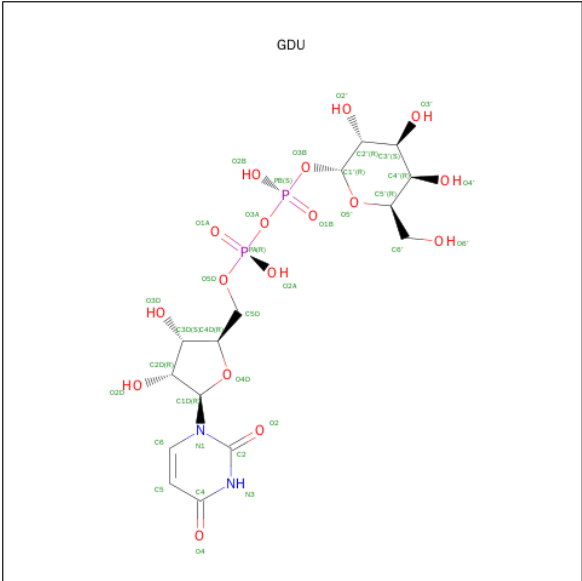
Chain	Residue	Modelled	Actual	Comment	Reference
A	73	ILE	VAL	SEE REMARK 999	UNP Q48485
A	222	ASP	GLU	SEE REMARK 999	UNP Q48485
A	258	ILE	THR	SEE REMARK 999	UNP Q48485
A	372	ASP	GLU	SEE REMARK 999	UNP Q48485
A	384	GLY	ARG	ENGINEERED	UNP Q48485
A	385	HIS	-	EXPRESSION TAG	UNP Q48485
A	386	HIS	-	EXPRESSION TAG	UNP Q48485
A	387	HIS	-	EXPRESSION TAG	UNP Q48485
A	388	HIS	-	EXPRESSION TAG	UNP Q48485
A	389	HIS	-	EXPRESSION TAG	UNP Q48485
A	390	HIS	-	EXPRESSION TAG	UNP Q48485
B	73	ILE	VAL	SEE REMARK 999	UNP Q48485
B	222	ASP	GLU	SEE REMARK 999	UNP Q48485
B	258	ILE	THR	SEE REMARK 999	UNP Q48485
B	372	ASP	GLU	SEE REMARK 999	UNP Q48485
B	384	GLY	ARG	ENGINEERED	UNP Q48485
B	385	HIS	-	EXPRESSION TAG	UNP Q48485
B	386	HIS	-	EXPRESSION TAG	UNP Q48485
B	387	HIS	-	EXPRESSION TAG	UNP Q48485
B	388	HIS	-	EXPRESSION TAG	UNP Q48485
B	389	HIS	-	EXPRESSION TAG	UNP Q48485
B	390	HIS	-	EXPRESSION TAG	UNP Q48485

- 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		

- Molecule 3 is SUGAR (GALACTOSE-URIDINE-5'-DIPHOSPHATE) (three-letter code: GDU) (formula: C<sub>15</sub>H<sub>24</sub>N<sub>2</sub>O<sub>17</sub>P<sub>2</sub>).



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

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

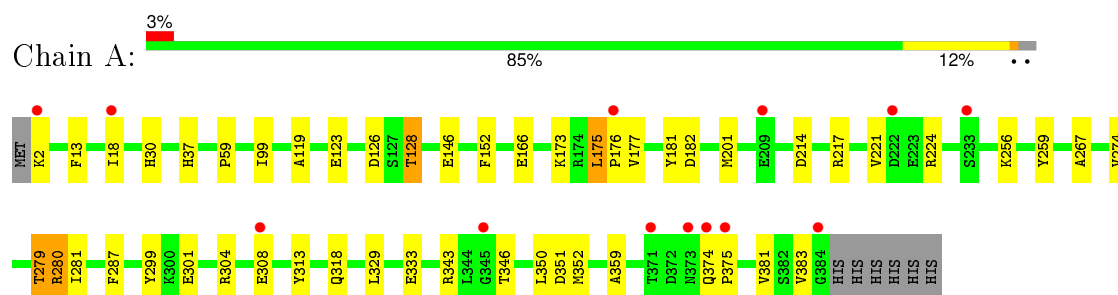
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	143	Total	O	0	0
			143	143		
4	B	144	Total	O	0	0
			144	144		

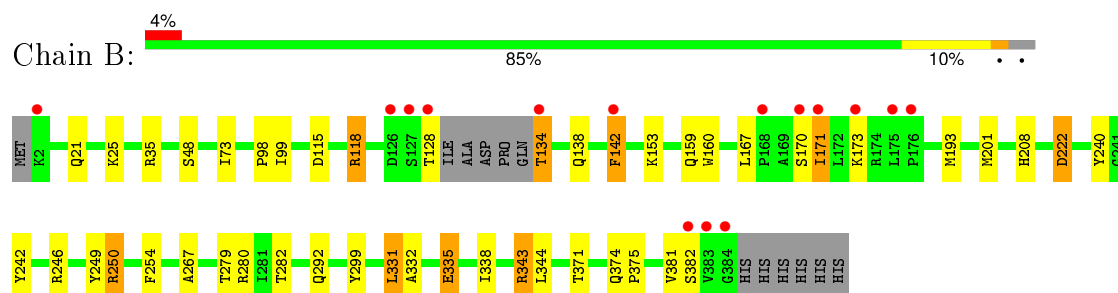
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.80 Å 93.80 Å 128.58 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 29.66 – 2.29	Depositor EDS
% Data completeness (in resolution range)	98.3 (30.00-2.30) 97.7 (29.66-2.29)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.81 (at 2.29 Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.187 , 0.233 0.182 , 0.224	Depositor DCC
$R_{free}$ test set	2421 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.9	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 31.6	EDS
Estimated twinning fraction	0.036 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48602 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6666	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GDU, 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	1.03	7/3202 (0.2%)	0.89	6/4336 (0.1%)
1	B	1.07	2/3163 (0.1%)	0.91	6/4280 (0.1%)
All	All	1.05	9/6365 (0.1%)	0.90	12/8616 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	146	GLU	CD-OE1	6.06	1.32	1.25
1	A	301	GLU	CB-CG	5.76	1.63	1.52
1	B	335	GLU	CD-OE2	-5.75	1.19	1.25
1	A	119	ALA	CA-CB	5.50	1.64	1.52
1	A	274	VAL	CB-CG1	5.27	1.64	1.52
1	A	177	VAL	CA-CB	-5.20	1.43	1.54
1	A	181	TYR	CD1-CE1	5.18	1.47	1.39
1	A	166	GLU	CG-CD	5.13	1.59	1.51
1	B	249	TYR	CD2-CE2	5.06	1.47	1.39

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	331	LEU	CB-CG-CD1	-10.29	93.50	111.00
1	A	343	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	B	343	ARG	NE-CZ-NH2	-6.55	117.02	120.30
1	A	280	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	A	182	ASP	CB-CG-OD2	-6.43	112.52	118.30
1	B	280	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	B	35	ARG	NE-CZ-NH1	5.93	123.26	120.30
1	A	217	ARG	NE-CZ-NH2	-5.77	117.41	120.30
1	B	118	ARG	NE-CZ-NH2	-5.35	117.63	120.30
1	B	222	ASP	CB-CG-OD2	5.15	122.94	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	182	ASP	CB-CG-OD1	5.09	122.88	118.30
1	A	304	ARG	CB-CA-C	-5.01	100.38	110.40

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3119	0	2998	29	0
1	B	3082	0	2962	36	0
2	A	53	0	31	3	0
2	B	53	0	31	0	0
3	A	36	0	22	1	0
3	B	36	0	22	0	0
4	A	143	0	0	4	0
4	B	144	0	0	5	0
All	All	6666	0	6066	65	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 (65) 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:B:73:ILE:HG22	1:B:201:MET:HE1	1.46	0.95
1:B:73:ILE:CG2	1:B:201:MET:HE1	2.02	0.89
1:B:73:ILE:HG22	1:B:201:MET:CE	2.09	0.81
1:B:160:TRP:HA	1:B:250:ARG:HD2	1.64	0.79
1:B:335:GLU:HG2	1:B:338:ILE:HG21	1.68	0.75
1:A:13:PHE:HE2	1:A:201:MET:HE2	1.53	0.73
1:A:99:ILE:HD11	1:A:152:PHE:HE1	1.53	0.73
1:B:153:LYS:HE3	4:B:648:HOH:O	1.91	0.70
1:A:99:ILE:HD11	1:A:152:PHE:CE1	2.27	0.69
1:A:329:LEU:O	1:A:333:GLU:HG2	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:ASP:OD1	1:A:128:THR:HB	1.94	0.68
1:B:208:HIS:O	4:B:513:HOH:O	2.10	0.68
1:A:351:ASP:OD1	4:A:427:HOH:O	2.15	0.65
1:B:335:GLU:HG2	1:B:338:ILE:CG2	2.27	0.64
1:A:99:ILE:CD1	1:A:152:PHE:HE1	2.13	0.62
1:B:170:SER:O	1:B:171:ILE:HB	2.01	0.60
1:A:99:ILE:CD1	1:A:152:PHE:CE1	2.83	0.60
1:B:159:GLN:O	1:B:250:ARG:HD3	2.01	0.60
1:B:134:THR:HA	1:B:167:LEU:O	2.02	0.59
1:B:240:TYR:CZ	1:B:331:LEU:CD1	2.85	0.59
1:B:331:LEU:HD12	1:B:331:LEU:C	2.23	0.58
1:B:118:ARG:HD3	4:B:640:HOH:O	2.02	0.57
1:B:73:ILE:CG2	1:B:201:MET:CE	2.76	0.56
1:A:175:LEU:HD21	3:A:392:GDU:H5'2	1.90	0.54
1:A:13:PHE:HE2	1:A:201:MET:CE	2.19	0.53
1:B:240:TYR:OH	1:B:332:ALA:HA	2.09	0.52
1:B:240:TYR:CD1	1:B:331:LEU:HD11	2.45	0.52
1:B:240:TYR:CE1	1:B:331:LEU:HD11	2.46	0.51
1:B:138:GLN:NE2	1:B:142:PHE:HD2	2.09	0.50
1:B:21:GLN:OE1	4:B:621:HOH:O	2.20	0.50
1:A:30:HIS:NE2	1:A:214:ASP:OD2	2.25	0.50
1:A:374:GLN:HB3	1:A:375:PRO:HD2	1.92	0.49
1:B:240:TYR:CE2	1:B:331:LEU:CD1	2.95	0.49
1:B:242:TYR:CD1	1:B:246:ARG:HG2	2.48	0.48
1:A:13:PHE:CE2	1:A:201:MET:CE	2.96	0.47
1:B:343:ARG:HG3	1:B:344:LEU:HG	1.95	0.47
1:B:240:TYR:CE2	1:B:331:LEU:HD13	2.50	0.47
1:A:18:ILE:HG13	1:A:359:ALA:HB1	1.96	0.47
1:A:99:ILE:HD12	1:A:152:PHE:CZ	2.50	0.46
1:A:346:THR:HB	1:A:383:VAL:HG11	1.96	0.46
1:B:267:ALA:O	1:B:282:THR:HA	2.15	0.46
1:A:175:LEU:HA	1:A:176:PRO:HD3	1.85	0.46
1:B:193:MET:HB3	1:B:193:MET:HE3	1.45	0.45
1:B:159:GLN:O	1:B:250:ARG:CD	2.64	0.45
1:A:279:THR:HG22	1:A:280:ARG:HG2	1.99	0.45
1:B:331:LEU:HD12	1:B:332:ALA:N	2.31	0.45
1:A:37:HIS:CE1	4:A:585:HOH:O	2.70	0.45
1:A:308:GLU:HB2	4:A:635:HOH:O	2.16	0.44
1:A:259:TYR:OH	1:B:115:ASP:OD2	2.24	0.44
1:A:13:PHE:CE2	1:A:201:MET:HE2	2.42	0.43
1:B:374:GLN:HB3	1:B:375:PRO:HD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:391:FAD:H1'1	2:A:391:FAD:H9	1.82	0.43
1:A:267:ALA:HB1	1:A:287:PHE:CE2	2.54	0.42
1:A:256:LYS:HG3	1:A:299:TYR:CZ	2.54	0.42
1:A:350:LEU:O	2:A:391:FAD:H1'2	2.19	0.42
1:A:37:HIS:HE1	4:A:585:HOH:O	2.03	0.42
1:A:313:TYR:HD2	2:A:391:FAD:HM82	1.84	0.42
1:A:352:MET:HB2	1:A:352:MET:HE3	1.91	0.42
1:B:254:PHE:HB3	1:B:299:TYR:CD2	2.55	0.42
1:B:240:TYR:CZ	1:B:331:LEU:HD11	2.55	0.41
1:B:99:ILE:HD12	4:B:578:HOH:O	2.18	0.41
1:B:381:VAL:HG22	1:B:382:SER:N	2.36	0.41
1:B:25:LYS:HD3	1:B:25:LYS:HA	1.87	0.41
1:B:98:PRO:O	1:B:99:ILE:C	2.59	0.41
1:A:281:ILE:HA	1:A:299:TYR:O	2.21	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	381/390 (98%)	373 (98%)	8 (2%)	0	100	100
1	B	374/390 (96%)	361 (96%)	12 (3%)	1 (0%)	46	57
All	All	755/780 (97%)	734 (97%)	20 (3%)	1 (0%)	56	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	ILE

### 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	336/343 (98%)	325 (97%)	11 (3%)	45	61
1	B	332/343 (97%)	322 (97%)	10 (3%)	48	65
All	All	668/686 (97%)	647 (97%)	21 (3%)	47	64

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	59	PRO
1	A	123	GLU
1	A	128	THR
1	A	173	LYS
1	A	175	LEU
1	A	221	VAL
1	A	224	ARG
1	A	279	THR
1	A	318	GLN
1	A	381	VAL
1	B	48	SER
1	B	128	THR
1	B	134	THR
1	B	142	PHE
1	B	173	LYS
1	B	222	ASP
1	B	250	ARG
1	B	279	THR
1	B	292	GLN
1	B	371	THR

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

Mol	Chain	Res	Type
1	A	51	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

4 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$
2	FAD	A	391	-	48,58,58	1.38	6 (12%)	54,89,89	2.29	9 (16%)
3	GDU	A	392	-	29,38,38	0.84	0	43,58,58	1.60	6 (13%)
2	FAD	B	391	-	48,58,58	1.33	8 (16%)	54,89,89	2.31	9 (16%)
3	GDU	B	392	-	29,38,38	0.81	0	43,58,58	1.86	7 (16%)

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	A	391	-	-	0/30/50/50	0/6/6/6
3	GDU	A	392	-	-	0/19/59/59	0/3/3/3
2	FAD	B	391	-	-	0/30/50/50	0/6/6/6
3	GDU	B	392	-	-	0/19/59/59	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	391	FAD	C6-C5X	-2.22	1.38	1.41
2	B	391	FAD	P-O2P	-2.18	1.45	1.54
2	B	391	FAD	C4X-N5	2.10	1.36	1.33
2	B	391	FAD	C2A-N1A	2.10	1.37	1.33
2	B	391	FAD	C1'-N10	2.13	1.50	1.48
2	A	391	FAD	C5X-N5	2.41	1.39	1.35
2	A	391	FAD	C5'-C4'	2.59	1.55	1.51
2	B	391	FAD	C4-N3	2.62	1.38	1.33
2	A	391	FAD	C2A-N1A	2.65	1.38	1.33
2	B	391	FAD	C5X-N5	2.80	1.39	1.35
2	A	391	FAD	C4-N3	2.93	1.38	1.33
2	A	391	FAD	C4X-N5	3.40	1.38	1.33
2	A	391	FAD	C2A-N3A	3.87	1.39	1.32
2	B	391	FAD	C2A-N3A	4.26	1.39	1.32

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	391	FAD	N3A-C2A-N1A	-13.43	118.61	128.89
2	A	391	FAD	N3A-C2A-N1A	-12.30	119.48	128.89
3	B	392	GDU	O3A-PB-O3B	-3.55	93.42	103.63
2	B	391	FAD	C4X-C4-N3	-3.44	118.88	123.59
2	A	391	FAD	C4X-C4-N3	-3.09	119.36	123.59
2	B	391	FAD	P-O3P-PA	-3.03	124.21	132.73
3	B	392	GDU	C6-N1-C2	-2.59	117.08	121.28
2	A	391	FAD	P-O3P-PA	-2.45	125.84	132.73
3	A	392	GDU	O5'-C1'-O3B	-2.25	108.40	111.36
3	B	392	GDU	PB-O3A-PA	-2.10	126.83	132.73
3	B	392	GDU	O6'-C6'-C5'	-2.07	104.48	111.33
2	B	391	FAD	C4X-N5-C5X	2.02	119.08	116.76
2	B	391	FAD	C4X-C10-N10	2.08	121.74	120.52
2	B	391	FAD	O4B-C1B-N9A	2.12	112.53	108.10
2	B	391	FAD	O3P-PA-O5B	2.13	108.58	102.94
2	A	391	FAD	O3P-P-O5'	2.22	108.83	102.94
3	A	392	GDU	O3'-C3'-C2'	2.24	115.38	110.34
2	A	391	FAD	O2A-PA-O3P	2.29	115.49	105.09
3	A	392	GDU	O2'-C2'-C3'	2.33	115.57	110.34
3	B	392	GDU	O3B-C1'-C2'	2.41	112.90	108.39
2	A	391	FAD	C4-C4X-N5	2.63	121.91	118.72
3	A	392	GDU	C3'-C4'-C5'	2.91	115.27	110.20
2	A	391	FAD	C4X-N5-C5X	2.95	120.16	116.76
3	B	392	GDU	C4D-O4D-C1D	2.97	112.98	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	392	GDU	O3B-C1'-C2'	3.12	114.23	108.39
2	A	391	FAD	C5X-C9A-N10	3.35	120.16	117.62
2	B	391	FAD	C4B-O4B-C1B	3.36	113.41	109.72
2	B	391	FAD	C4-N3-C2	4.93	119.51	115.25
2	A	391	FAD	C4-N3-C2	6.23	120.63	115.25
3	A	392	GDU	C4-N3-C2	6.38	120.46	114.14
3	B	392	GDU	C4-N3-C2	8.32	122.38	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	391	FAD	3	0
3	A	392	GDU	1	0

## 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	383/390 (98%)	-0.08	13 (3%) 49 58	20, 31, 56, 69	0
1	B	378/390 (96%)	-0.04	15 (3%) 42 51	18, 30, 59, 77	0
All	All	761/780 (97%)	-0.06	28 (3%) 45 54	18, 31, 57, 77	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	384	GLY	5.4
1	A	373	ASN	4.6
1	B	383	VAL	4.5
1	B	171	ILE	4.2
1	A	374	GLN	3.9
1	A	176	PRO	3.8
1	B	168	PRO	3.6
1	A	308	GLU	3.2
1	A	384	GLY	3.2
1	A	371	THR	3.0
1	A	2	LYS	2.9
1	B	142	PHE	2.9
1	B	128	THR	2.8
1	A	375	PRO	2.8
1	B	127	SER	2.7
1	B	382	SER	2.7
1	B	2	LYS	2.6
1	B	134	THR	2.4
1	A	209	GLU	2.4
1	B	126	ASP	2.3
1	B	176	PRO	2.3
1	A	233	SER	2.2
1	B	175	LEU	2.2
1	A	222	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	345	GLY	2.1
1	A	18	ILE	2.1
1	B	170	SER	2.1
1	B	173	LYS	2.1

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

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

## 6.3 Carbohydrates [i](#)

There are no 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GDU	B	392	36/36	0.87	0.16	0.12	45,61,64,64	0
2	FAD	B	391	53/53	0.98	0.15	0.07	16,23,28,33	0
3	GDU	A	392	36/36	0.94	0.12	-0.07	22,34,51,54	0
2	FAD	A	391	53/53	0.97	0.14	-0.11	19,26,31,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.