



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2016 – 01:33 PM EST

PDB ID : 5GY7  
Title : X-Ray structure of H243I mutant of UDP-Galactose 4-epimerase from E.coli:evidence for existence of open and closed active site during catalysis.  
Authors : Singh, N.; Tiwari, P.; Phulera, S.; Dixit, A.; Choudhury, D.  
Deposited on : 2016-09-21  
Resolution : 1.43 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028320  
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) : rb-20028320

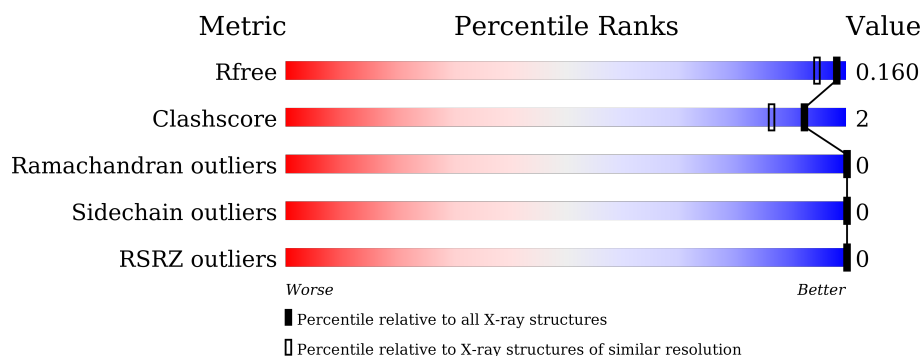
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.43 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	1164 (1.46-1.42)
Clashscore	102246	1219 (1.46-1.42)
Ramachandran outliers	100387	1200 (1.46-1.42)
Sidechain outliers	100360	1200 (1.46-1.42)
RSRZ outliers	91569	1166 (1.46-1.42)

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	344	
1	B	344	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	403	-	-	-	X
5	NO3	B	403	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5934 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.

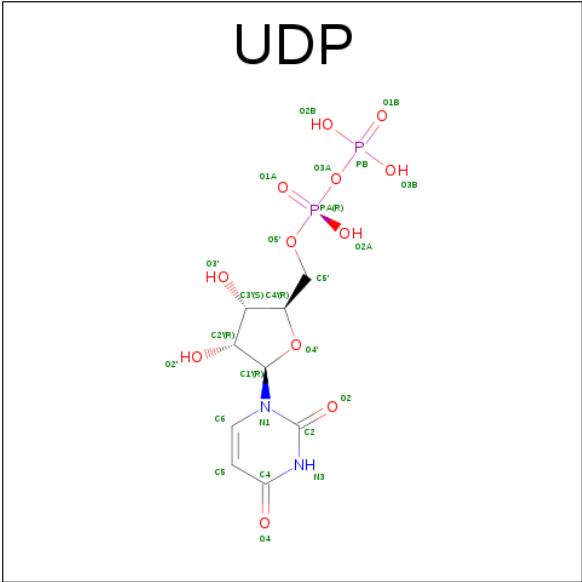
- Molecule 1 is a protein called UDP-glucose 4-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2581	1632	452	486	11			
1	B	338	Total	C	N	O	S	0	0	0
			2561	1618	447	485	11			

There are 14 discrepancies between the modelled and reference sequences:

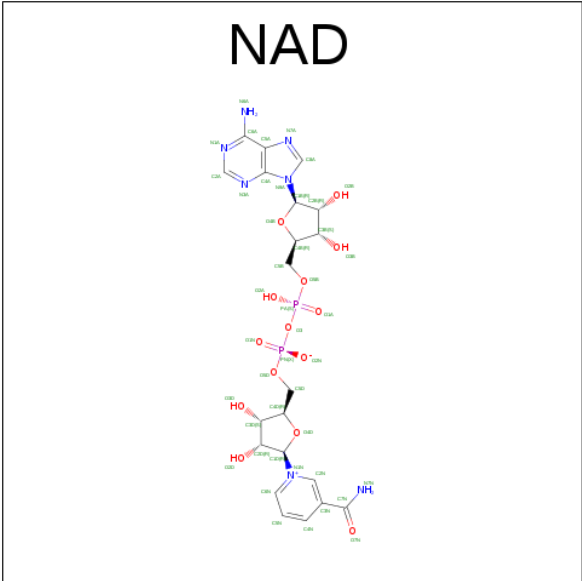
Chain	Residue	Modelled	Actual	Comment	Reference
A	243	ILE	HIS	engineered mutation	UNP P09147
A	339	HIS	-	expression tag	UNP P09147
A	340	HIS	-	expression tag	UNP P09147
A	341	HIS	-	expression tag	UNP P09147
A	342	HIS	-	expression tag	UNP P09147
A	343	HIS	-	expression tag	UNP P09147
A	344	HIS	-	expression tag	UNP P09147
B	243	ILE	HIS	engineered mutation	UNP P09147
B	339	HIS	-	expression tag	UNP P09147
B	340	HIS	-	expression tag	UNP P09147
B	341	HIS	-	expression tag	UNP P09147
B	342	HIS	-	expression tag	UNP P09147
B	343	HIS	-	expression tag	UNP P09147
B	344	HIS	-	expression tag	UNP P09147

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>2</sub>O<sub>12</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



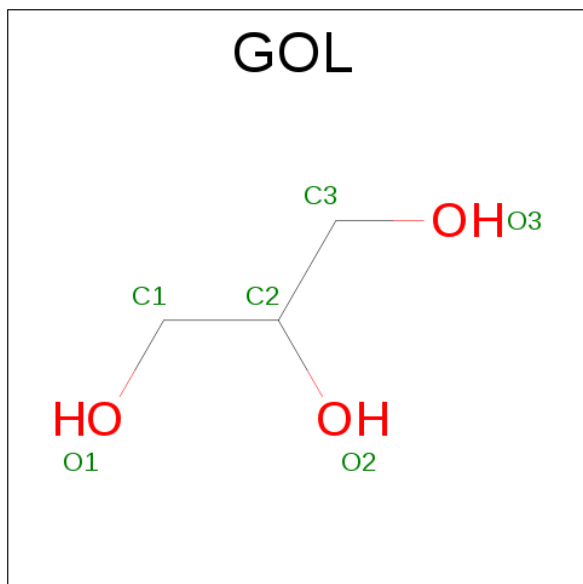
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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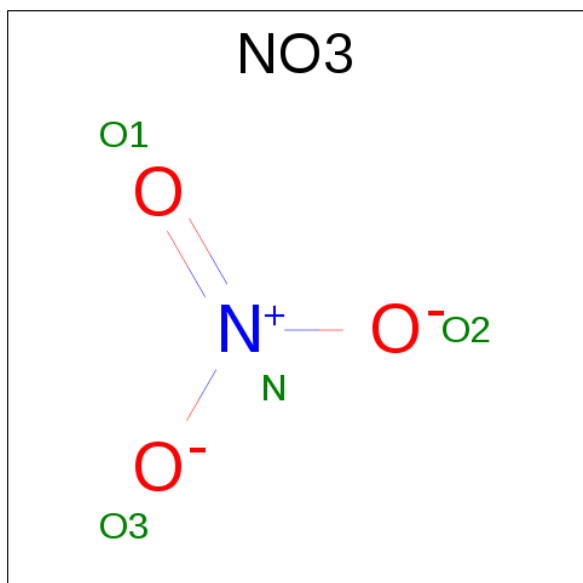
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	
			44	21	7	14	2	
								0
								0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O		
			6	3	3	0	0

- Molecule 5 is NITRATE ION (three-letter code: NO3) (formula:  $NO_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	N	O	0	0
			4	1	3		

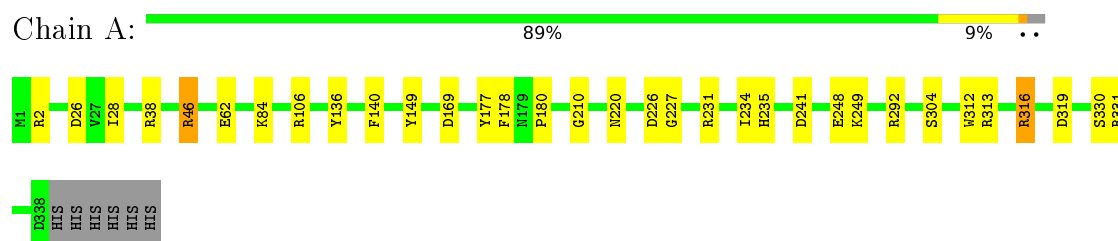
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	332	Total	O	0	0
			332	332		
6	B	312	Total	O	0	0
			312	312		

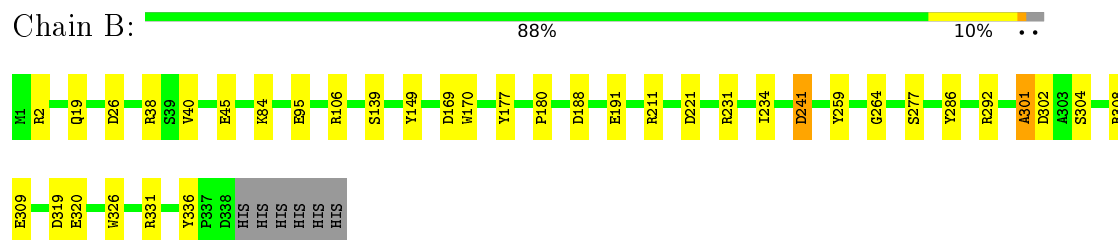
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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 ( $RSRZ > 2$ ). 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-glucose 4-epimerase



- Molecule 1: UDP-glucose 4-epimerase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	47.11Å 91.16Å 80.26Å 90.00° 99.87° 90.00°	Depositor
Resolution (Å)	79.07 – 1.43 41.36 – 1.43	Depositor EDS
% Data completeness (in resolution range)	98.7 (79.07-1.43) 98.7 (41.36-1.43)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	7.73 (at 1.43Å)	Xtriage
Refinement program	REFMAC 5.8.0151	Depositor
R, $R_{free}$	0.134 , 0.160 0.134 , 0.160	Depositor DCC
$R_{free}$ test set	6013 reflections (5.21%)	DCC
Wilson B-factor (Å <sup>2</sup> )	9.3	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 46.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5934	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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: GOL, UDP, NAD, NO3

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	1.34	9/2646 (0.3%)	1.34	22/3611 (0.6%)
1	B	1.36	12/2624 (0.5%)	1.31	25/3584 (0.7%)
All	All	1.35	21/5270 (0.4%)	1.32	47/7195 (0.7%)

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 (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	139	SER	CB-OG	-9.45	1.29	1.42
1	A	210	GLY	CA-C	-7.46	1.40	1.51
1	A	227	GLY	N-CA	-6.32	1.36	1.46
1	A	304	SER	CB-OG	-6.19	1.34	1.42
1	B	45	GLU	CD-OE2	6.08	1.32	1.25
1	B	286	TYR	CZ-OH	5.94	1.48	1.37
1	A	248	GLU	CD-OE1	5.75	1.31	1.25
1	B	95	GLU	CG-CD	5.70	1.60	1.51
1	B	304	SER	CB-OG	-5.54	1.35	1.42
1	A	136	TYR	CE1-CZ	5.53	1.45	1.38
1	B	170	TRP	CZ3-CH2	-5.46	1.31	1.40
1	B	320	GLU	CG-CD	-5.35	1.44	1.51
1	A	292	ARG	CZ-NH1	-5.33	1.26	1.33
1	B	95	GLU	CD-OE1	5.31	1.31	1.25
1	B	169	ASP	CB-CG	-5.29	1.40	1.51
1	A	62	GLU	CD-OE2	5.26	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	326	TRP	CE3-CZ3	5.21	1.47	1.38
1	B	286	TYR	CE2-CZ	-5.19	1.31	1.38
1	B	277	SER	CB-OG	5.17	1.49	1.42
1	A	312	TRP	CZ3-CH2	5.12	1.48	1.40
1	A	330	SER	CB-OG	-5.08	1.35	1.42

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	38	ARG	NE-CZ-NH2	-11.24	114.68	120.30
1	B	308	ARG	NE-CZ-NH2	-10.87	114.86	120.30
1	A	292	ARG	NE-CZ-NH2	-10.38	115.11	120.30
1	B	26	ASP	CB-CG-OD2	-10.29	109.04	118.30
1	A	46	ARG	NE-CZ-NH1	10.26	125.43	120.30
1	B	211	ARG	NE-CZ-NH2	-9.95	115.32	120.30
1	B	292	ARG	NE-CZ-NH1	-8.55	116.03	120.30
1	A	241	ASP	CB-CG-OD1	8.35	125.81	118.30
1	A	319	ASP	CB-CG-OD1	8.05	125.55	118.30
1	B	177	TYR	CB-CG-CD1	7.93	125.76	121.00
1	A	169	ASP	CB-CG-OD1	7.91	125.42	118.30
1	B	319	ASP	CB-CG-OD1	7.55	125.10	118.30
1	A	106	ARG	NE-CZ-NH1	7.49	124.05	120.30
1	B	211	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	A	231	ARG	NE-CZ-NH2	-7.37	116.61	120.30
1	B	188	ASP	CB-CG-OD1	7.32	124.89	118.30
1	A	46	ARG	NE-CZ-NH2	-7.27	116.67	120.30
1	B	169	ASP	CB-CG-OD2	-7.01	111.99	118.30
1	B	169	ASP	CB-CG-OD1	6.80	124.42	118.30
1	B	309	GLU	OE1-CD-OE2	-6.77	115.18	123.30
1	A	331	ARG	NE-CZ-NH2	-6.72	116.94	120.30
1	B	231	ARG	NE-CZ-NH2	-6.66	116.97	120.30
1	A	292	ARG	NH1-CZ-NH2	6.61	126.67	119.40
1	A	177	TYR	CB-CG-CD1	6.53	124.92	121.00
1	B	38	ARG	NE-CZ-NH2	6.47	123.53	120.30
1	A	2	ARG	NE-CZ-NH2	-6.37	117.12	120.30
1	B	221	ASP	CB-CG-OD1	6.31	123.98	118.30
1	B	191	GLU	OE1-CD-OE2	6.00	130.50	123.30
1	B	302	ASP	CB-CG-OD1	5.92	123.63	118.30
1	A	313	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	B	292	ARG	NH1-CZ-NH2	5.90	125.89	119.40
1	B	286	TYR	CB-CG-CD2	-5.90	117.46	121.00
1	A	38	ARG	NE-CZ-NH1	5.88	123.24	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	331	ARG	NE-CZ-NH1	-5.61	117.50	120.30
1	A	136	TYR	CB-CG-CD1	5.42	124.25	121.00
1	A	226	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	A	316	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	B	286	TYR	CB-CG-CD1	5.36	124.22	121.00
1	A	140	PHE	CB-CG-CD2	-5.23	117.14	120.80
1	B	259	TYR	CZ-CE2-CD2	-5.21	115.11	119.80
1	A	231	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	B	2	ARG	CG-CD-NE	-5.17	100.94	111.80
1	A	26	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	B	106	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	B	188	ASP	CB-CG-OD2	-5.03	113.77	118.30
1	A	178	PHE	CB-CA-C	5.01	120.43	110.40
1	B	241	ASP	CB-CG-OD1	5.01	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	301	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	2581	0	2481	13	0
1	B	2561	0	2451	8	0
2	A	25	0	11	0	0
2	B	25	0	11	0	0
3	A	44	0	26	0	0
3	B	44	0	26	0	0
4	A	6	0	8	0	0
5	B	4	0	0	0	0
6	A	332	0	0	7	2
6	B	312	0	0	1	2
All	All	5934	0	5014	21	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (21) 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:19:GLN:NE2	1:B:241:ASP:OD1	1.68	1.25
1:A:46:ARG:HD3	6:A:782:HOH:O	1.38	1.22
1:A:249:LYS:HE2	6:A:501:HOH:O	1.05	1.18
1:B:19:GLN:HG3	6:B:566:HOH:O	1.52	1.07
1:A:249:LYS:CE	6:A:501:HOH:O	1.66	0.94
1:A:46:ARG:CD	6:A:782:HOH:O	2.05	0.92
1:A:249:LYS:NZ	6:A:501:HOH:O	1.90	0.77
1:A:46:ARG:CG	6:A:782:HOH:O	2.35	0.62
1:B:19:GLN:HE22	1:B:241:ASP:CG	1.98	0.58
1:A:46:ARG:HG3	6:A:782:HOH:O	2.07	0.52
1:A:84:LYS:HA	1:A:149:TYR:CE1	2.49	0.47
1:B:84:LYS:HA	1:B:149:TYR:CE1	2.49	0.47
1:A:235:HIS:CD2	1:A:316:ARG:HD3	2.50	0.46
1:B:180:PRO:HA	1:B:234:ILE:O	2.15	0.46
1:A:180:PRO:HA	1:A:234:ILE:O	2.15	0.45
1:A:220:ASN:OD1	1:A:220:ASN:C	2.52	0.45
1:B:40:VAL:HG23	1:B:336:TYR:CZ	2.52	0.44
1:A:235:HIS:CG	1:A:316:ARG:HD3	2.54	0.42
1:A:28:ILE:HD12	1:A:28:ILE:N	2.35	0.41
1:B:19:GLN:HE22	1:B:241:ASP:HA	1.86	0.41
1:B:264:GLY:HA2	1:B:301:ALA:O	2.20	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:504:HOH:O	6:B:591:HOH:O[1_454]	1.92	0.28
6:A:782:HOH:O	6:B:642:HOH:O[1_554]	2.10	0.10

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	336/344 (98%)	333 (99%)	3 (1%)	0	100	100
1	B	336/344 (98%)	332 (99%)	4 (1%)	0	100	100
All	All	672/688 (98%)	665 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	271/288 (94%)	271 (100%)	0	100	100
1	B	267/288 (93%)	267 (100%)	0	100	100
All	All	538/576 (93%)	538 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	B	19	GLN

### 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 [i](#)

There are no carbohydrates 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
2	UDP	A	401	-	20,26,26	1.57	4 (20%)	24,40,40	2.17	3 (12%)
3	NAD	A	402	-	42,48,48	1.48	7 (16%)	46,73,73	1.41	5 (10%)
4	GOL	A	403	-	5,5,5	1.85	1 (20%)	5,5,5	1.92	1 (20%)
2	UDP	B	401	-	20,26,26	1.24	2 (10%)	24,40,40	2.21	2 (8%)
3	NAD	B	402	-	42,48,48	1.74	9 (21%)	46,73,73	1.76	9 (19%)
5	NO3	B	403	-	1,3,3	2.01	1 (100%)	0,3,3	0.00	-

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	UDP	A	401	-	-	0/12/32/32	0/2/2/2
3	NAD	A	402	-	-	0/22/62/62	0/5/5/5
4	GOL	A	403	-	-	0/4/4/4	0/0/0/0
2	UDP	B	401	-	-	0/12/32/32	0/2/2/2
3	NAD	B	402	-	-	0/22/62/62	0/5/5/5
5	NO3	B	403	-	-	0/0/0/0	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	NAD	C3N-C7N	-3.85	1.44	1.50
3	B	402	NAD	C2B-C1B	-3.13	1.48	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	UDP	C2'-C1'	-2.91	1.49	1.53
3	B	402	NAD	C2N-C3N	-2.82	1.34	1.39
2	A	401	UDP	C2-N3	-2.80	1.32	1.38
3	A	402	NAD	C2N-C3N	-2.66	1.34	1.39
3	B	402	NAD	C6N-C5N	-2.40	1.33	1.38
3	A	402	NAD	C2B-C3B	-2.28	1.47	1.53
2	A	401	UDP	PA-O2A	-2.12	1.46	1.55
3	A	402	NAD	C6N-C5N	-2.11	1.34	1.38
5	B	403	NO3	O1-N	2.01	1.29	1.23
3	A	402	NAD	C6N-N1N	2.01	1.41	1.35
2	A	401	UDP	C6-N1	2.05	1.38	1.35
2	B	401	UDP	C6-N1	2.38	1.38	1.35
3	B	402	NAD	C2A-N3A	2.73	1.37	1.32
3	B	402	NAD	O4D-C1D	3.19	1.45	1.41
3	B	402	NAD	O4B-C1B	3.21	1.45	1.41
3	A	402	NAD	C4N-C3N	3.30	1.44	1.39
3	B	402	NAD	C5N-C4N	3.51	1.45	1.38
4	A	403	GOL	O2-C2	3.52	1.53	1.43
3	A	402	NAD	O4B-C1B	3.83	1.46	1.41
3	A	402	NAD	C5N-C4N	3.98	1.46	1.38
2	A	401	UDP	C4-N3	4.00	1.40	1.33
3	B	402	NAD	C4N-C3N	4.80	1.47	1.39

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	NAD	N3A-C2A-N1A	-5.36	124.66	128.87
3	A	402	NAD	C1B-N9A-C4A	-4.89	121.35	126.81
3	B	402	NAD	C5N-C4N-C3N	-3.97	115.61	120.35
3	B	402	NAD	C4N-C3N-C7N	-2.91	113.39	121.11
2	A	401	UDP	C5-C4-N3	-2.83	116.32	123.28
3	A	402	NAD	C5N-C4N-C3N	-2.80	117.01	120.35
3	A	402	NAD	N3A-C2A-N1A	-2.69	126.75	128.87
3	B	402	NAD	O7N-C7N-N7N	-2.13	119.55	122.58
2	A	401	UDP	O2A-PA-O3A	2.07	114.13	105.27
3	A	402	NAD	O4D-C4D-C3D	2.08	109.39	105.16
2	B	401	UDP	O2B-PB-O1B	2.15	117.64	110.63
3	B	402	NAD	C2A-N1A-C6A	2.18	122.65	118.77
3	A	402	NAD	C3N-C2N-N1N	2.20	122.86	120.34
3	B	402	NAD	C2D-C1D-N1N	2.66	118.74	113.53
3	B	402	NAD	O7N-C7N-C3N	2.72	122.61	119.60
3	B	402	NAD	C2N-C3N-C7N	3.56	129.46	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	403	GOL	O2-C2-C1	3.74	126.44	108.47
3	B	402	NAD	C3N-C2N-N1N	4.32	125.28	120.34
2	A	401	UDP	C4-N3-C2	8.77	123.44	114.21
2	B	401	UDP	C4-N3-C2	9.66	124.39	114.21

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	338/344 (98%)	-0.56	0 100 100	6, 10, 17, 28	0
1	B	338/344 (98%)	-0.52	0 100 100	5, 9, 16, 30	0
All	All	676/688 (98%)	-0.54	0 100 100	5, 9, 17, 30	0

There are no RSRZ outliers to report.

### 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(Å <sup>2</sup> )	Q<0.9
5	NO3	B	403	4/4	0.93	0.15	13.02	21,21,22,26	0
4	GOL	A	403	6/6	0.94	0.10	5.65	12,18,22,27	0
3	NAD	A	402	44/44	0.99	0.06	-0.04	5,6,9,11	0
3	NAD	B	402	44/44	0.99	0.05	-0.49	5,6,9,11	0
2	UDP	A	401	25/25	0.99	0.05	-1.22	7,7,10,15	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	UDP	B	401	25/25	0.99	0.05	-1.52	5,6,9,13	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.