



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:38 PM GMT

PDB ID : 1KVT  
Title : UDP-GALACTOSE 4-EPIMERASE COMPLEXED WITH UDP-PHENOL  
Authors : Thoden, J.B.; Gulick, A.; Holden, H.M.  
Deposited on : 1997-03-07  
Resolution : 2.15 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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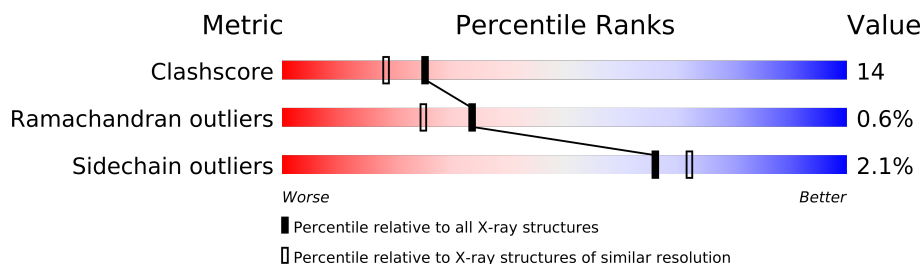
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.15 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))
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	338	

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 3095 atoms, of which 0 are hydrogen and 0 are deuterium.

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-GALACTOSE 4-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	338	2626	1657	463	494	12	0	0	0

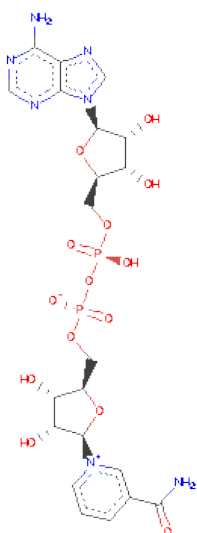
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	124	VAL	SER	ENGINEERED	UNP P09147
A	131	ASN	GLN	CONFLICT	UNP P09147

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

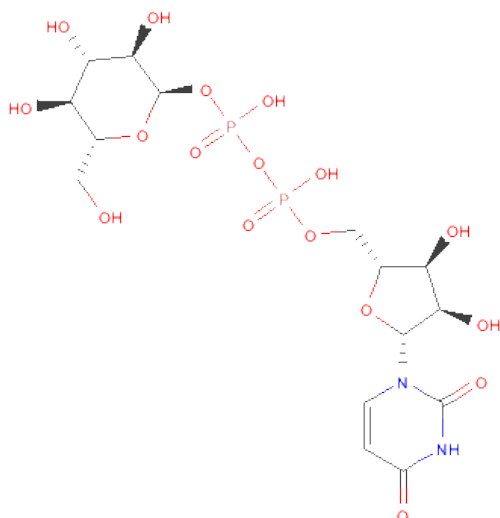
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

- 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>).



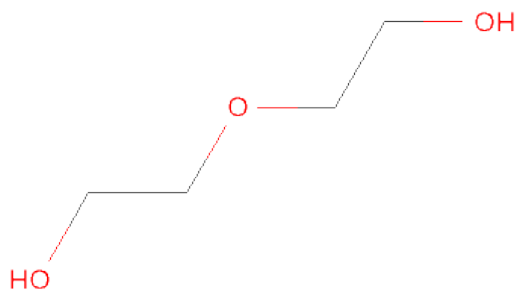
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is URIDINE-5'-DIPHOSPHATE-GLUCOSE (three-letter code: UPG) (formula:  $C_{15}H_{24}N_2O_{17}P_2$ ).



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	381	Total	O	0	0
			381	381		

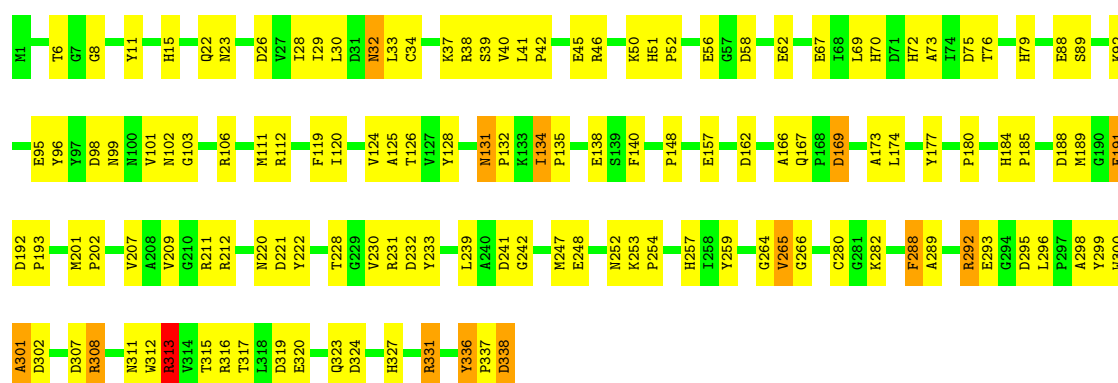
### 3 Residue-property plots

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.

Note EDS was not executed.

#### • Molecule 1: UDP-GALACTOSE 4-EPIMERASE

Chain A: 



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.50Å 83.50Å 108.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.15	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.15)	Depositor
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	TNT	Depositor
R, $R_{free}$	0.178 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3095	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: UPG, PEG, NAD, NA

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	1.02	12/2692 (0.4%)	1.65	60/3664 (1.6%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	191	GLU	CD-OE1	8.22	1.34	1.25
1	A	67	GLU	CD-OE1	7.24	1.33	1.25
1	A	62	GLU	CD-OE2	-7.24	1.17	1.25
1	A	95	GLU	CD-OE1	7.01	1.33	1.25
1	A	293	GLU	CD-OE1	5.73	1.31	1.25
1	A	138	GLU	CD-OE1	5.61	1.31	1.25
1	A	88	GLU	CD-OE1	-5.53	1.19	1.25
1	A	191	GLU	CD-OE2	-5.49	1.19	1.25
1	A	56	GLU	CD-OE2	5.47	1.31	1.25
1	A	248	GLU	CD-OE1	-5.46	1.19	1.25
1	A	140	PHE	CD1-CE1	5.23	1.49	1.39
1	A	188	ASP	CG-OD1	5.01	1.36	1.25

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	319	ASP	CB-CG-OD1	11.97	129.07	118.30
1	A	106	ARG	NE-CZ-NH2	-11.87	114.37	120.30
1	A	319	ASP	CB-CG-OD2	-11.82	107.66	118.30
1	A	58	ASP	CB-CG-OD2	-10.88	108.50	118.30
1	A	292	ARG	NE-CZ-NH1	9.06	124.83	120.30
1	A	188	ASP	CB-CG-OD1	-8.73	110.44	118.30
1	A	38	ARG	NE-CZ-NH1	8.70	124.65	120.30
1	A	241	ASP	CB-CG-OD1	8.06	125.56	118.30
1	A	241	ASP	CB-CG-OD2	-7.96	111.14	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ARG	NE-CZ-NH1	7.85	124.23	120.30
1	A	233	TYR	CB-CG-CD1	-7.82	116.31	121.00
1	A	169	ASP	CB-CG-OD1	-7.70	111.37	118.30
1	A	315	THR	CA-CB-CG2	-7.37	102.09	112.40
1	A	26	ASP	CB-CG-OD2	6.99	124.59	118.30
1	A	112	ARG	NE-CZ-NH2	-6.90	116.85	120.30
1	A	228	THR	CA-CB-CG2	-6.76	102.94	112.40
1	A	221	ASP	CB-CG-OD2	6.71	124.33	118.30
1	A	221	ASP	CB-CG-OD1	-6.41	112.53	118.30
1	A	211	ARG	NE-CZ-NH1	6.32	123.46	120.30
1	A	192	ASP	CB-CG-OD2	-6.29	112.64	118.30
1	A	338	ASP	CB-CG-OD2	6.24	123.91	118.30
1	A	128	TYR	CB-CG-CD1	-6.20	117.28	121.00
1	A	212	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	A	131	ASN	N-CA-CB	6.10	121.58	110.60
1	A	101	VAL	CA-CB-CG1	-6.05	101.83	110.90
1	A	26	ASP	CB-CG-OD1	-6.02	112.88	118.30
1	A	298	ALA	CB-CA-C	-5.95	101.17	110.10
1	A	75	ASP	CB-CA-C	-5.93	98.53	110.40
1	A	259	TYR	CB-CG-CD2	5.93	124.56	121.00
1	A	192	ASP	CB-CG-OD1	5.92	123.62	118.30
1	A	302	ASP	N-CA-C	-5.84	95.23	111.00
1	A	324	ASP	CB-CG-OD2	-5.82	113.06	118.30
1	A	177	TYR	CB-CG-CD2	-5.81	117.52	121.00
1	A	75	ASP	CB-CG-OD1	-5.78	113.09	118.30
1	A	233	TYR	CB-CG-CD2	5.74	124.45	121.00
1	A	336	TYR	CB-CG-CD2	-5.71	117.57	121.00
1	A	301	ALA	N-CA-CB	-5.71	102.10	110.10
1	A	301	ALA	CB-CA-C	-5.69	101.56	110.10
1	A	289	ALA	CB-CA-C	5.68	118.62	110.10
1	A	96	TYR	CA-CB-CG	-5.67	102.62	113.40
1	A	313	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	A	58	ASP	CB-CG-OD1	5.56	123.30	118.30
1	A	298	ALA	N-CA-CB	-5.55	102.32	110.10
1	A	308	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	338	ASP	CB-CG-OD1	-5.49	113.36	118.30
1	A	211	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	A	119	PHE	CB-CG-CD2	-5.46	116.98	120.80
1	A	331	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	40	VAL	CG1-CB-CG2	-5.40	102.27	110.90
1	A	265	VAL	CG1-CB-CG2	-5.31	102.40	110.90
1	A	259	TYR	CB-CG-CD1	-5.28	117.83	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	ASP	CB-CG-OD2	5.27	123.05	118.30
1	A	299	TYR	CB-CG-CD1	-5.26	117.84	121.00
1	A	288	PHE	CB-CG-CD1	5.21	124.44	120.80
1	A	162	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	162	ASP	CB-CG-OD1	-5.13	113.68	118.30
1	A	222	TYR	CB-CG-CD2	-5.12	117.93	121.00
1	A	76	THR	CA-CB-CG2	-5.08	105.29	112.40
1	A	46	ARG	NE-CZ-NH1	-5.07	117.77	120.30
1	A	207	VAL	CA-CB-CG1	-5.06	103.31	110.90

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2626	0	2559	73	1
2	A	1	0	0	0	0
3	A	44	0	26	0	0
4	A	36	0	22	2	0
5	A	7	0	10	8	0
6	A	381	0	0	8	9
All	All	3095	0	2617	73	9

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (73) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:253:LYS:NZ	5:A:410:PEG:H12	1.80	0.96
1:A:37:LYS:HE2	1:A:338:ASP:OD1	1.66	0.95
1:A:253:LYS:HZ3	5:A:410:PEG:H12	1.33	0.94
1:A:22:GLN:HG2	6:A:687:HOH:O	1.68	0.92
1:A:184:HIS:ND1	1:A:185:PRO:HD2	1.85	0.91

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:124:VAL:HG21	4:A:341:UPG:O4'	1.74	0.88
1:A:23:ASN:OD1	6:A:424:HOH:O	1.93	0.87
1:A:253:LYS:CE	5:A:410:PEG:H12	2.04	0.86
1:A:184:HIS:CG	1:A:185:PRO:HD2	2.15	0.81
1:A:131:ASN:OD1	1:A:132:PRO:HD2	1.89	0.71
1:A:124:VAL:HG22	1:A:126:THR:H	1.56	0.71
1:A:265:VAL:HG22	1:A:266:GLY:N	2.06	0.70
1:A:253:LYS:HE2	5:A:410:PEG:H12	1.74	0.68
1:A:184:HIS:CE1	1:A:185:PRO:HD2	2.31	0.66
1:A:265:VAL:CG2	1:A:266:GLY:N	2.59	0.66
1:A:30:LEU:HD13	1:A:30:LEU:C	2.17	0.66
1:A:191:GLU:HG2	1:A:193:PRO:HD3	1.78	0.65
1:A:124:VAL:CG2	4:A:341:UPG:O4'	2.45	0.64
1:A:89:SER:O	1:A:148:PRO:HG2	2.00	0.62
1:A:180:PRO:HB3	1:A:239:LEU:CD2	2.30	0.61
1:A:253:LYS:HZ3	5:A:410:PEG:C1	2.10	0.61
1:A:327:HIS:O	1:A:331:ARG:HG3	2.02	0.59
1:A:22:GLN:NE2	6:A:449:HOH:O	2.22	0.58
1:A:292:ARG:HB3	1:A:295:ASP:OD2	2.05	0.57
1:A:8:GLY:HA3	1:A:29:ILE:HG23	1.87	0.57
1:A:120:ILE:HD11	1:A:247:MET:HA	1.88	0.56
1:A:253:LYS:NZ	5:A:410:PEG:C1	2.64	0.56
1:A:242:GLY:HA2	1:A:312:TRP:CD1	2.42	0.55
1:A:232:ASP:HB2	1:A:300:TRP:HA	1.89	0.54
1:A:280:CYS:SG	1:A:282:LYS:HG2	2.47	0.54
1:A:32:ASN:HD22	1:A:33:LEU:N	2.06	0.54
1:A:265:VAL:HG23	6:A:623:HOH:O	2.08	0.54
1:A:32:ASN:C	1:A:32:ASN:HD22	2.11	0.53
1:A:252:ASN:O	1:A:254:PRO:HD3	2.09	0.52
1:A:220:ASN:HB3	1:A:288:PHE:CD1	2.45	0.52
1:A:232:ASP:HB2	1:A:300:TRP:CB	2.39	0.52
1:A:201:MET:HB3	1:A:202:PRO:HD3	1.90	0.52
1:A:99:ASN:O	1:A:103:GLY:HA3	2.11	0.51
1:A:184:HIS:CG	1:A:185:PRO:CD	2.91	0.50
1:A:98:ASP:O	1:A:102:ASN:HB3	2.12	0.50
1:A:15:HIS:HA	1:A:189:MET:HE1	1.95	0.49
1:A:313:ARG:HD2	6:A:653:HOH:O	2.13	0.48
1:A:39:SER:HB2	1:A:337:PRO:HD2	1.96	0.48
1:A:317:THR:OG1	1:A:320:GLU:HG3	2.13	0.48
1:A:157:GLU:HB3	1:A:174:LEU:HD13	1.96	0.47
1:A:41:LEU:N	1:A:42:PRO:HD2	2.29	0.47
1:A:28:ILE:N	1:A:28:ILE:HD12	2.29	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:69:LEU:HD13	1:A:111:MET:HA	1.96	0.47
1:A:32:ASN:ND2	1:A:34:CYS:H	2.12	0.47
1:A:135:PRO:HB3	1:A:300:TRP:CD1	2.51	0.45
1:A:134:ILE:HA	1:A:135:PRO:C	2.38	0.44
1:A:124:VAL:HG22	1:A:125:ALA:N	2.33	0.44
1:A:169:ASP:N	1:A:169:ASP:OD1	2.36	0.44
1:A:264:GLY:HA2	1:A:301:ALA:O	2.18	0.43
1:A:253:LYS:HZ3	5:A:410:PEG:C3	2.32	0.43
1:A:189:MET:HA	1:A:336:TYR:OH	2.19	0.43
1:A:232:ASP:HB2	1:A:300:TRP:HB2	2.01	0.43
1:A:166:ALA:C	1:A:167:GLN:HG2	2.38	0.43
1:A:45:GLU:HG3	1:A:51:HIS:HA	2.00	0.43
1:A:6:THR:OG1	1:A:79:HIS:HA	2.19	0.42
1:A:230:VAL:C	1:A:231:ARG:HG2	2.39	0.42
1:A:92:LYS:HE2	6:A:498:HOH:O	2.20	0.42
1:A:307:ASP:O	1:A:311:ASN:HA	2.20	0.42
1:A:296:LEU:HA	1:A:296:LEU:HD23	1.86	0.42
1:A:41:LEU:N	1:A:42:PRO:CD	2.82	0.42
1:A:307:ASP:O	1:A:311:ASN:N	2.53	0.41
1:A:173:ALA:HA	1:A:257:HIS:O	2.19	0.41
1:A:45:GLU:HG3	1:A:50:LYS:O	2.20	0.41
1:A:135:PRO:HD2	6:A:518:HOH:O	2.20	0.41
1:A:253:LYS:HE2	5:A:410:PEG:C1	2.44	0.41
1:A:72:HIS:O	1:A:73:ALA:C	2.59	0.41
1:A:323:GLN:HG2	6:A:749:HOH:O	2.20	0.40
1:A:316:ARG:HA	1:A:320:GLU:OE1	2.22	0.40

All (9) 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	Distance(Å)	Clash(Å)
6:A:645:HOH:O	6:A:645:HOH:O[6_765]	0.35	1.85
6:A:520:HOH:O	6:A:520:HOH:O[4_646]	0.46	1.74
6:A:692:HOH:O	6:A:692:HOH:O[5_556]	0.49	1.71
6:A:519:HOH:O	6:A:519:HOH:O[4_646]	0.62	1.58
6:A:724:HOH:O	6:A:724:HOH:O[5_556]	0.65	1.55
6:A:652:HOH:O	6:A:652:HOH:O[4_646]	0.74	1.46
6:A:528:HOH:O	6:A:528:HOH:O[5_556]	1.95	0.25
6:A:532:HOH:O	6:A:724:HOH:O[5_556]	2.12	0.08
1:A:167:GLN:NE2	6:A:500:HOH:O[5_556]	2.18	0.02

## 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	336/338 (99%)	322 (96%)	12 (4%)	2 (1%)	33	25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	TYR
1	A	52	PRO

### 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	282/282 (100%)	276 (98%)	6 (2%)	66	72

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	70	HIS
1	A	134	ILE
1	A	209	VAL
1	A	308	ARG
1	A	313	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN

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Mol	Chain	Res	Type
1	A	158	GLN
1	A	194	GLN
1	A	267	ASN
1	A	274	ASN
1	A	323	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	$\# Z  > 2$
3	NAD	A	340	-	48,48,48	2.15	11 (22%)	73,73,73	1.69	12 (16%)
4	UPG	A	341	-	38,38,38	1.47	6 (15%)	54,58,58	1.62	14 (25%)
5	PEG	A	410	-	6,6,6	0.86	0	5,5,5	1.50	1 (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	NAD	A	340	-	-	0/30/62/62	0/3/5/5
4	UPG	A	341	-	-	0/21/59/59	0/3/3/3
5	PEG	A	410	-	-	0/4/4/4	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	340	NAD	C4N-C3N	7.42	1.52	1.39
3	A	340	NAD	C2N-C3N	6.61	1.48	1.38
3	A	340	NAD	C5N-C4N	5.89	1.52	1.39
4	A	341	UPG	PB-O3B	4.91	1.75	1.60
3	A	340	NAD	C2N-N1N	3.91	1.40	1.35
3	A	340	NAD	PA-O3	-3.31	1.53	1.59
4	A	341	UPG	PB-O3A	-3.16	1.54	1.59
4	A	341	UPG	C6-C5	-2.96	1.31	1.36
3	A	340	NAD	O4D-C1D	-2.76	1.37	1.41
3	A	340	NAD	C2A-N1A	2.75	1.39	1.33
3	A	340	NAD	C3N-C7N	-2.69	1.45	1.50
4	A	341	UPG	C6-N1	2.50	1.39	1.35
4	A	341	UPG	O3B-C1'	2.40	1.46	1.42
3	A	340	NAD	C6N-N1N	2.39	1.42	1.35
3	A	340	NAD	C2D-C1D	-2.34	1.50	1.53
3	A	340	NAD	O4B-C1B	2.14	1.44	1.41
4	A	341	UPG	PB-O2B	-2.06	1.45	1.55

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	340	NAD	C5N-C4N-C3N	-7.47	110.62	120.32
4	A	341	UPG	O3B-C1'-C2'	5.13	117.77	108.38
3	A	340	NAD	O7N-C7N-C3N	-4.67	114.31	119.58
3	A	340	NAD	C8A-N9A-C4A	-3.90	103.92	106.90
3	A	340	NAD	C5A-C6A-N6A	3.84	129.40	120.72
3	A	340	NAD	C6N-N1N-C2N	-3.83	117.71	122.04
3	A	340	NAD	N6A-C6A-N1A	-3.58	112.34	119.36
4	A	341	UPG	C1'-C2'-C3'	3.36	116.54	110.00
4	A	341	UPG	C2-N1-C1C	3.26	120.25	118.21
4	A	341	UPG	O5'-C5'-C4'	2.89	115.11	109.76
5	A	410	PEG	C3-O2-C2	2.89	126.07	113.38
4	A	341	UPG	O2C-C2C-C1C	-2.69	103.08	111.23
3	A	340	NAD	C3N-C7N-N7N	2.66	120.80	117.77
3	A	340	NAD	C2N-C3N-C4N	2.65	121.31	118.31
4	A	341	UPG	C4C-O4C-C1C	2.62	112.59	109.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	340	NAD	C6N-C5N-C4N	2.52	123.44	119.44
4	A	341	UPG	O5'-C1'-C2'	-2.44	105.31	110.31
4	A	341	UPG	PB-O3B-C1'	2.44	129.89	119.58
4	A	341	UPG	C3C-C2C-C1C	2.34	104.57	100.91
4	A	341	UPG	O3A-PB-O3B	2.27	109.22	103.48
4	A	341	UPG	O2C-C2C-C3C	-2.27	104.46	111.83
4	A	341	UPG	O4C-C1C-C2C	-2.22	103.36	106.77
3	A	340	NAD	O4B-C1B-N9A	2.17	110.46	108.44
3	A	340	NAD	N3A-C2A-N1A	2.14	130.49	128.71
3	A	340	NAD	O4D-C1D-N1N	2.12	110.12	107.95
4	A	341	UPG	O2'-C2'-C1'	2.08	114.58	110.04
4	A	341	UPG	O3A-PA-O5C	-2.02	94.38	103.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.