



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:03 PM GMT

PDB ID : 1NAI
Title : UDP-GALACTOSE 4-EPIMERASE FROM ESCHERICHIA COLI, OXIDIZED
Authors : Thoden, J.B.; Frey, P.A.; Holden, H.M.
Deposited on : 1995-11-22
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

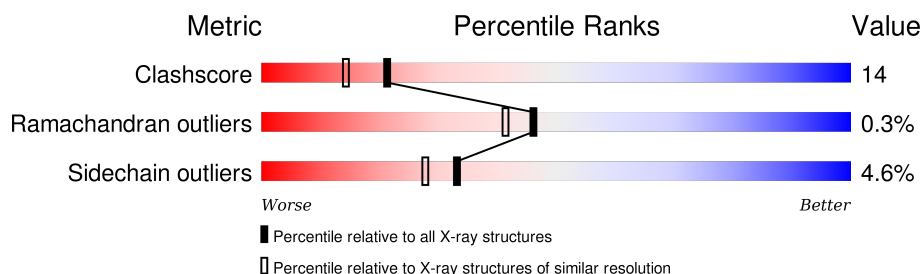
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	338	

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
5	PDO	A	351	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3084 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-GALACTOSE 4-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	3	0
			2643	1671	463	497	12			

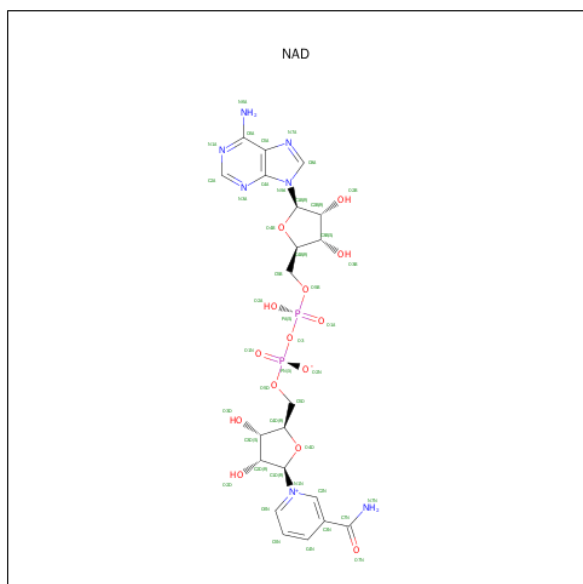
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	131	ASN	GLN	CONFLICT	UNP P09147

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

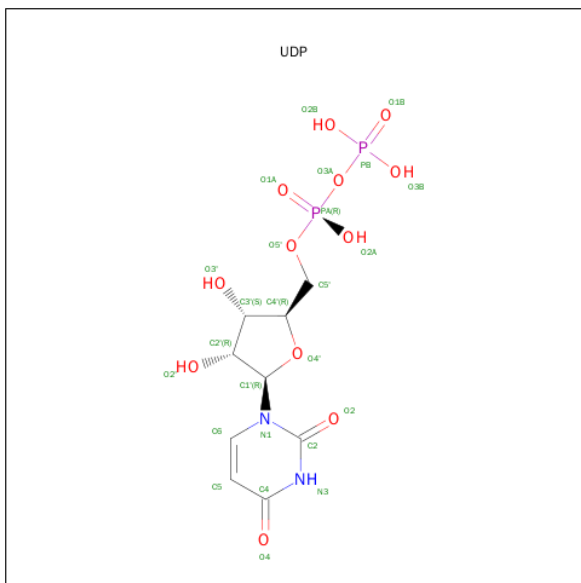
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Na	0	0
			2	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



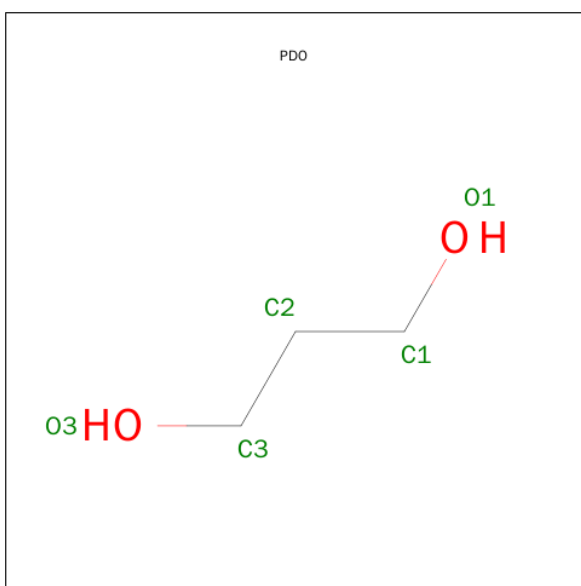
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $\text{C}_9\text{H}_{14}\text{N}_2\text{O}_{12}\text{P}_2$).



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

- Molecule 5 is 1,3-PROPANDIOL (three-letter code: PDO) (formula: $\text{C}_3\text{H}_8\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			5	3	2		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 7 is water.

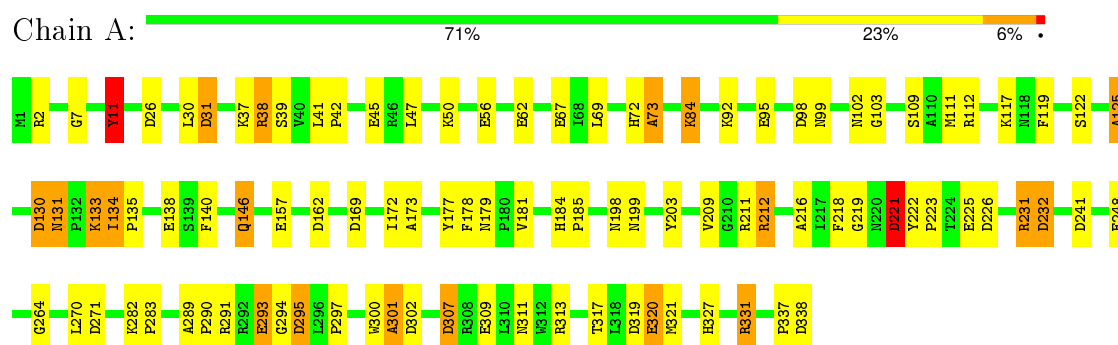
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	358	Total	O	0	0
			358	358		

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



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, α , β , γ	83.50Å 83.50Å 108.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.193 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3084	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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: PDO, NA, UDP, PEG, NAD

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	0.97	11/2723 (0.4%)	1.40	40/3705 (1.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	157	GLU	CD-OE1	7.52	1.33	1.25
1	A	138	GLU	CD-OE1	7.09	1.33	1.25
1	A	225	GLU	CD-OE2	6.85	1.33	1.25
1	A	95	GLU	CD-OE1	6.44	1.32	1.25
1	A	67	GLU	CD-OE1	6.10	1.32	1.25
1	A	293	GLU	CD-OE1	5.89	1.32	1.25
1	A	62	GLU	CD-OE1	5.71	1.31	1.25
1	A	320	GLU	CD-OE1	-5.38	1.19	1.25
1	A	56	GLU	CD-OE2	5.28	1.31	1.25
1	A	248	GLU	CD-OE2	5.18	1.31	1.25
1	A	309	GLU	CD-OE1	5.11	1.31	1.25

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	231	ARG	NE-CZ-NH1	11.93	126.27	120.30
1	A	232[A]	ASP	CB-CG-OD1	10.79	128.01	118.30
1	A	232[B]	ASP	CB-CG-OD1	10.79	128.01	118.30
1	A	221	ASP	CB-CG-OD2	-10.01	109.29	118.30
1	A	295	ASP	CB-CG-OD2	-7.88	111.21	118.30
1	A	122	SER	CB-CA-C	-7.52	95.82	110.10
1	A	271	ASP	CB-CG-OD2	-7.15	111.86	118.30
1	A	221	ASP	N-CA-CB	-6.91	98.17	110.60
1	A	26	ASP	CB-CG-OD1	-6.84	112.14	118.30
1	A	241	ASP	CB-CG-OD1	6.78	124.40	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	271	ASP	CB-CG-OD1	6.76	124.38	118.30
1	A	232[A]	ASP	CB-CG-OD2	-6.75	112.23	118.30
1	A	232[B]	ASP	CB-CG-OD2	-6.75	112.23	118.30
1	A	221	ASP	CB-CG-OD1	6.61	124.25	118.30
1	A	319	ASP	CB-CG-OD2	-6.55	112.41	118.30
1	A	38	ARG	NE-CZ-NH1	6.44	123.52	120.30
1	A	307	ASP	CB-CG-OD1	-6.19	112.73	118.30
1	A	295	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	211	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	A	313	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	A	319	ASP	CB-CG-OD1	5.86	123.57	118.30
1	A	169	ASP	CB-CG-OD1	-5.84	113.04	118.30
1	A	231	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	A	307	ASP	CB-CG-OD2	5.83	123.55	118.30
1	A	11	TYR	CB-CG-CD1	5.79	124.48	121.00
1	A	26	ASP	CB-CG-OD2	5.75	123.48	118.30
1	A	302	ASP	CB-CG-OD2	5.72	123.45	118.30
1	A	131	ASN	N-CA-CB	5.71	120.88	110.60
1	A	173	ALA	N-CA-CB	5.69	118.06	110.10
1	A	31	ASP	CB-CG-OD2	5.64	123.38	118.30
1	A	125	ALA	CB-CA-C	5.58	118.47	110.10
1	A	226	ASP	CB-CG-OD2	-5.55	113.31	118.30
1	A	338	ASP	CB-CG-OD1	-5.54	113.31	118.30
1	A	226	ASP	CB-CG-OD1	5.47	123.22	118.30
1	A	146	GLN	N-CA-CB	-5.47	100.75	110.60
1	A	301	ALA	CB-CA-C	-5.30	102.15	110.10
1	A	289	ALA	CB-CA-C	5.28	118.03	110.10
1	A	241	ASP	CB-CG-OD2	-5.23	113.59	118.30
1	A	162	ASP	CB-CG-OD1	-5.17	113.65	118.30
1	A	73	ALA	N-CA-CB	-5.15	102.89	110.10

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	2643	0	2565	72	0
2	A	2	0	0	0	0
3	A	44	0	26	4	0
4	A	25	0	11	2	0
5	A	5	0	8	5	0
6	A	7	0	10	0	0
7	A	358	0	0	13	5
All	All	3084	0	2620	76	5

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 14.

All (76) 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:178[B]:PHE:CZ	1:A:301:ALA:HB2	2.00	0.95
1:A:178[B]:PHE:HB2	7:A:593:HOH:O	1.65	0.95
1:A:84:LYS:HZ1	5:A:351:PDO:H22	1.37	0.90
1:A:37:LYS:HE3	7:A:580:HOH:O	1.83	0.78
1:A:317:THR:OG1	1:A:320:GLU:HG3	1.88	0.73
1:A:218[A]:PHE:HD1	1:A:290:PRO:O	1.72	0.71
1:A:125:ALA:HB3	1:A:178[A]:PHE:CE1	2.26	0.70
1:A:84:LYS:NZ	5:A:351:PDO:H22	2.05	0.70
1:A:178[B]:PHE:CE1	1:A:301:ALA:HB2	2.27	0.69
1:A:218[A]:PHE:CD1	1:A:290:PRO:O	2.45	0.69
1:A:184:HIS:CG	1:A:185:PRO:HD2	2.29	0.67
3:A:349:NAD:N7N	3:A:349:NAD:O1N	2.27	0.67
1:A:198:ASN:ND2	7:A:681:HOH:O	2.26	0.66
1:A:146:GLN:NE2	1:A:294:GLY:O	2.28	0.66
3:A:349:NAD:H5N	7:A:672:HOH:O	1.96	0.65
1:A:222:TYR:HB3	1:A:223:PRO:HD2	1.78	0.65
1:A:125:ALA:HB1	1:A:178[A]:PHE:CZ	2.31	0.65
1:A:221:ASP:HB3	1:A:291:ARG:HH21	1.62	0.65
1:A:177:TYR:HA	1:A:178[B]:PHE:CD1	2.32	0.64
1:A:177:TYR:HA	1:A:178[B]:PHE:HD1	1.62	0.64
1:A:39:SER:O	1:A:42:PRO:HD2	2.00	0.61
1:A:125:ALA:HB3	1:A:178[A]:PHE:CD1	2.35	0.61
1:A:125:ALA:CB	1:A:178[A]:PHE:CZ	2.84	0.61
1:A:11:TYR:HB2	5:A:351:PDO:H11	1.83	0.60
1:A:133:LYS:O	1:A:140:PHE:HE2	1.83	0.60
1:A:231:ARG:HD2	4:A:350:UDP:O2B	2.03	0.58
1:A:30:LEU:HD13	1:A:30:LEU:C	2.23	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:LYS:HE3	7:A:552:HOH:O	2.04	0.56
1:A:135:PRO:HB3	1:A:300:TRP:CD1	2.41	0.56
1:A:125:ALA:CB	1:A:178[A]:PHE:CE1	2.88	0.56
1:A:39:SER:OG	1:A:337:PRO:HD2	2.05	0.55
1:A:203:TYR:CE2	1:A:212:ARG:HD3	2.42	0.54
1:A:327:HIS:O	1:A:331:ARG:HG3	2.09	0.53
1:A:84:LYS:HD3	3:A:349:NAD:N7N	2.23	0.53
1:A:178[B]:PHE:CZ	1:A:301:ALA:CB	2.85	0.52
1:A:181:VAL:HG23	1:A:321:MET:HB3	1.92	0.52
1:A:181:VAL:HG12	7:A:371:HOH:O	2.09	0.51
1:A:131:ASN:HB2	7:A:708:HOH:O	2.14	0.47
3:A:349:NAD:O2A	5:A:351:PDO:H12	2.15	0.46
1:A:216:ALA:HB3	7:A:682:HOH:O	2.14	0.46
1:A:222:TYR:HB3	1:A:223:PRO:CD	2.45	0.46
1:A:178[B]:PHE:CE2	1:A:301:ALA:HB2	2.48	0.45
1:A:133:LYS:HB3	1:A:140:PHE:CE2	2.52	0.45
1:A:99:ASN:O	1:A:103:GLY:HA3	2.16	0.44
1:A:133:LYS:HB3	1:A:140:PHE:CZ	2.52	0.44
1:A:219:GLY:HA2	7:A:601:HOH:O	2.17	0.44
1:A:270:LEU:HD23	1:A:270:LEU:HA	1.86	0.44
1:A:199:ASN:HA	4:A:350:UDP:O2A	2.18	0.43
1:A:264:GLY:HA2	1:A:301:ALA:O	2.18	0.43
1:A:177:TYR:C	1:A:178[B]:PHE:CD1	2.92	0.43
1:A:282:LYS:HB2	1:A:283:PRO:HD2	2.01	0.42
1:A:41:LEU:N	1:A:42:PRO:CD	2.82	0.42
1:A:72:HIS:O	1:A:73:ALA:HB3	2.18	0.42
1:A:7:GLY:HA2	1:A:31:ASP:OD2	2.20	0.42
1:A:295:ASP:HA	7:A:412:HOH:O	2.19	0.42
1:A:47:LEU:HA	1:A:47:LEU:HD23	1.84	0.42
1:A:135:PRO:HB3	1:A:300:TRP:NE1	2.35	0.42
1:A:307:ASP:O	1:A:311:ASN:HA	2.20	0.42
1:A:69:LEU:HD13	1:A:111:MET:HA	2.00	0.42
1:A:223:PRO:HD2	1:A:297:PRO:HB3	2.02	0.41
1:A:327:HIS:HB3	7:A:432:HOH:O	2.20	0.41
5:A:351:PDO:H32	7:A:535:HOH:O	2.20	0.41
1:A:98:ASP:O	1:A:102:ASN:HB3	2.20	0.41
1:A:109:SER:O	1:A:112:ARG:HB2	2.19	0.41
1:A:45:GLU:HG3	1:A:50:LYS:O	2.21	0.41
1:A:130:ASP:O	7:A:643:HOH:O	2.21	0.41
1:A:222:TYR:CB	1:A:223:PRO:HD2	2.47	0.41
1:A:232[A]:ASP:HB2	1:A:300:TRP:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:PHE:O	1:A:172:ILE:HA	2.21	0.41
1:A:134:ILE:HG13	1:A:134:ILE:H	1.62	0.41

All (5) 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)
7:A:489:HOH:O	7:A:489:HOH:O[5_555]	1.07	1.13
7:A:562:HOH:O	7:A:562:HOH:O[6_555]	1.08	1.12
7:A:695:HOH:O	7:A:695:HOH:O[6_555]	1.59	0.61
7:A:527:HOH:O	7:A:527:HOH:O[4_556]	1.75	0.45
7:A:567:HOH:O	7:A:567:HOH:O[5_555]	1.95	0.25

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	339/338 (100%)	324 (96%)	14 (4%)	1 (0%)	46 41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	TYR

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	285/282 (101%)	272 (95%)	13 (5%)	33	28

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	38	ARG
1	A	84	LYS
1	A	117	LYS
1	A	130	ASP
1	A	133	LYS
1	A	134	ILE
1	A	179	ASN
1	A	209	VAL
1	A	212	ARG
1	A	221	ASP
1	A	293	GLU
1	A	331	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	GLN
1	A	167	GLN
1	A	274	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

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	349	-	38,48,48	1.20	4 (10%)	47,73,73	1.77	10 (21%)
4	UDP	A	350	-	18,26,26	1.49	3 (16%)	26,40,40	3.67	8 (30%)
5	PDO	A	351	-	4,4,4	0.27	0	3,3,3	0.60	0
6	PEG	A	352	-	6,6,6	0.60	0	5,5,5	1.75	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	349	-	-	0/22/62/62	0/5/5/5
4	UDP	A	350	-	-	0/12/32/32	0/2/2/2
5	PDO	A	351	-	-	0/2/2/2	0/0/0/0
6	PEG	A	352	-	-	0/4/4/4	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	350	UDP	C6-C5	-2.84	1.31	1.38
3	A	349	NAD	C3N-C7N	2.15	1.53	1.50
3	A	349	NAD	C6N-N1N	2.17	1.41	1.35
4	A	350	UDP	C6-N1	2.66	1.39	1.35
3	A	349	NAD	C4N-C3N	3.14	1.44	1.39
3	A	349	NAD	C2A-N1A	3.23	1.40	1.33
4	A	350	UDP	C4-N3	3.43	1.39	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	350	UDP	O3B-PB-O3A	-6.11	77.38	105.09
3	A	349	NAD	C5N-C4N-C3N	-5.05	113.98	120.33
3	A	349	NAD	C5N-C6N-N1N	-4.34	112.97	120.47
4	A	350	UDP	O3A-PA-O5'	-4.33	91.46	102.94
4	A	350	UDP	O2'-C2'-C3'	-3.66	99.91	111.83
4	A	350	UDP	C5-C4-N3	-3.21	114.88	123.12
3	A	349	NAD	C3N-C2N-N1N	-2.29	117.72	120.36
3	A	349	NAD	O3D-C3D-C4D	-2.15	104.62	111.05
3	A	349	NAD	C2D-C3D-C4D	-2.04	98.41	102.61
4	A	350	UDP	O2B-PB-O3A	2.00	114.18	105.09
4	A	350	UDP	O3'-C3'-C4'	2.26	117.82	111.05
3	A	349	NAD	C3N-C7N-N7N	2.60	120.67	117.82
3	A	349	NAD	O3-PN-O5D	2.89	110.59	102.94
4	A	350	UDP	O3B-PB-O2B	2.97	118.67	107.38
3	A	349	NAD	O4D-C1D-N1N	3.19	111.63	108.13
3	A	349	NAD	C2N-C3N-C4N	3.70	122.41	118.29
6	A	352	PEG	C3-O2-C2	3.82	129.75	113.31
3	A	349	NAD	C6N-C5N-C4N	4.81	126.72	119.44
4	A	350	UDP	C4-N3-C2	15.10	129.09	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	349	NAD	4	0
4	A	350	UDP	2	0
5	A	351	PDO	5	0

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.