



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 12:46 pm BST

PDB ID : 1LRJ
Title : Crystal Structure of E. coli UDP-Galactose 4-Epimerase Complexed with UDP-N-Acetylglucosamine
Authors : Thoden, J.B.; Henderson, J.M.; Fridovich-Keil, J.L.; Holden, H.M.
Deposited on : 2002-05-15
Resolution : 1.90 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

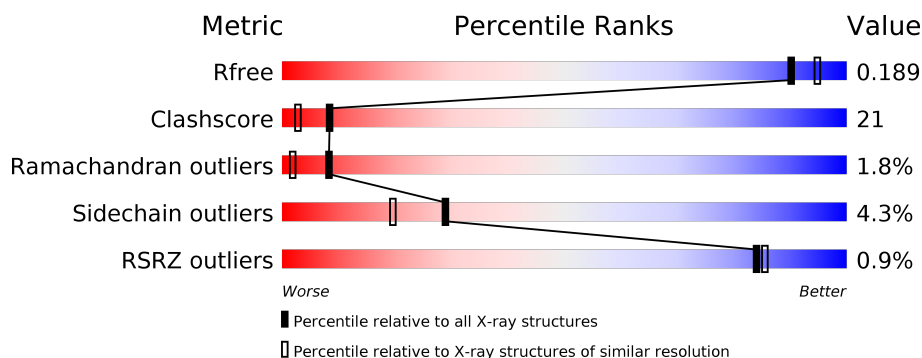
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 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	338	<div> <div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green);"></div> <div style="margin-left: 10px;"> % 62%33%•• </div> </div> </div>

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	PGE	A	342	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3149 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
			Total	C	N	O	S			
1	A	338	2635	1661	463	499	12	0	2	0

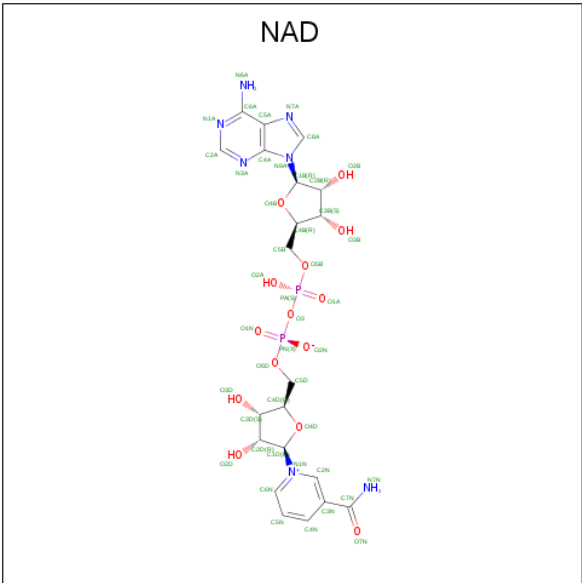
There is a discrepancy between the modelled and reference sequences:

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

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

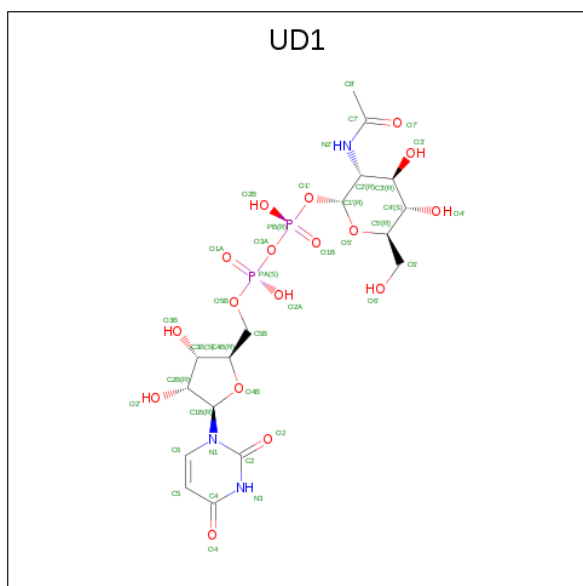
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Na	0	0
			2	2		

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



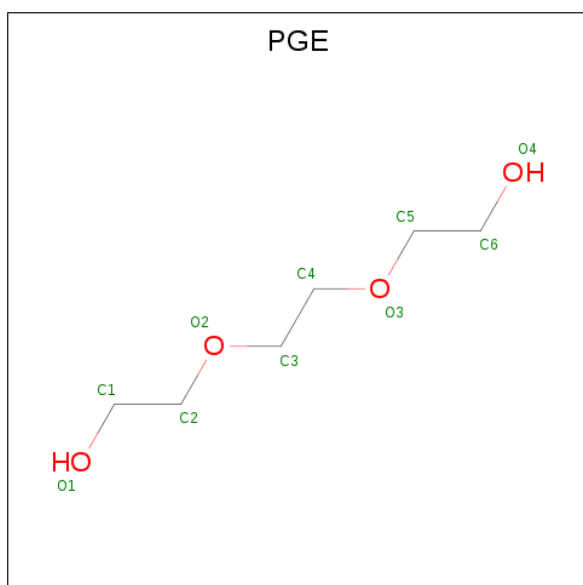
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (three-letter code: UD1) (formula: $C_{17}H_{27}N_3O_{17}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		

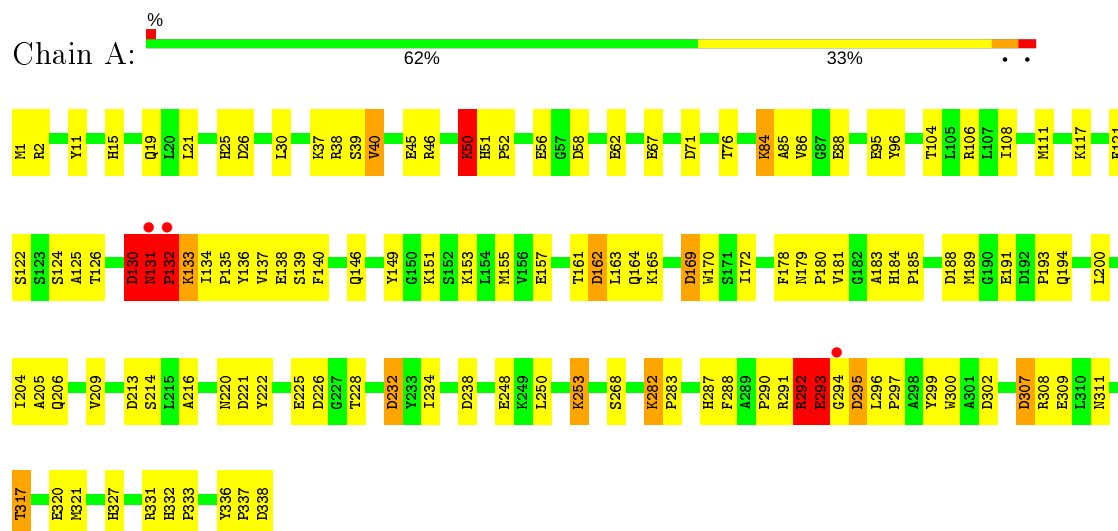
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	419	Total	O	0	0
			419	419		

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 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 ($\text{RSRZ} > 2$). 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-glucose 4-epimerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	83.20 Å 83.20 Å 108.30 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 1.90 29.99 – 1.75	Depositor EDS
% Data completeness (in resolution range)	93.4 (30.00-1.90) 94.7 (29.99-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.75 Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.192 , 0.268 0.193 , 0.189	Depositor DCC
R_{free} test set	4301 reflections (10.25%)	wwPDB-VP
Wilson B-factor (Å ²)	17.5	Xtriage
Anisotropy	0.417	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 126.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3149	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: NA, PGE, UD1, 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	1.08	13/2713 (0.5%)	1.39	37/3693 (1.0%)

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	A	1	2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	191	GLU	CD-OE2	8.36	1.34	1.25
1	A	293	GLU	CD-OE2	7.47	1.33	1.25
1	A	225	GLU	CD-OE2	7.33	1.33	1.25
1	A	157	GLU	CD-OE2	7.32	1.33	1.25
1	A	88	GLU	CD-OE2	6.56	1.32	1.25
1	A	62	GLU	CD-OE2	6.15	1.32	1.25
1	A	67	GLU	CD-OE2	6.15	1.32	1.25
1	A	309	GLU	CD-OE2	6.15	1.32	1.25
1	A	95	GLU	CD-OE2	6.09	1.32	1.25
1	A	56	GLU	CD-OE2	6.02	1.32	1.25
1	A	248	GLU	CD-OE2	5.21	1.31	1.25
1	A	138	GLU	CD-OE1	-5.17	1.20	1.25
1	A	56	GLU	CD-OE1	-5.08	1.20	1.25

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ARG	NE-CZ-NH1	10.68	125.64	120.30
1	A	132	PRO	C-N-CA	10.61	148.23	121.70
1	A	162	ASP	CB-CG-OD2	-9.23	109.99	118.30
1	A	221	ASP	CB-CG-OD2	-7.31	111.72	118.30
1	A	302	ASP	CB-CG-OD1	7.09	124.68	118.30
1	A	232[A]	ASP	CB-CG-OD2	-6.95	112.05	118.30
1	A	232[B]	ASP	CB-CG-OD2	-6.95	112.05	118.30
1	A	130	ASP	CB-CG-OD2	-6.72	112.25	118.30
1	A	307	ASP	CB-CG-OD2	-6.72	112.25	118.30
1	A	295	ASP	CB-CG-OD2	-6.63	112.33	118.30
1	A	295	ASP	CB-CG-OD1	6.55	124.19	118.30
1	A	46	ARG	NE-CZ-NH1	6.34	123.47	120.30
1	A	26	ASP	CB-CG-OD2	-6.31	112.62	118.30
1	A	307	ASP	CB-CG-OD1	6.27	123.94	118.30
1	A	232[A]	ASP	CB-CG-OD1	6.18	123.86	118.30
1	A	232[B]	ASP	CB-CG-OD1	6.18	123.86	118.30
1	A	26	ASP	CB-CG-OD1	6.14	123.83	118.30
1	A	213	ASP	CB-CG-OD2	-6.08	112.83	118.30
1	A	162	ASP	CB-CG-OD1	6.05	123.75	118.30
1	A	226	ASP	CB-CG-OD2	-6.05	112.85	118.30
1	A	71	ASP	CB-CG-OD2	-5.77	113.11	118.30
1	A	169	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	A	169	ASP	CB-CG-OD1	5.71	123.44	118.30
1	A	131	ASN	N-CA-CB	5.62	120.72	110.60
1	A	131	ASN	CB-CA-C	5.58	121.57	110.40
1	A	161	THR	CA-CB-CG2	-5.55	104.63	112.40
1	A	130	ASP	CB-CG-OD1	5.49	123.24	118.30
1	A	132	PRO	N-CA-CB	5.47	109.86	103.30
1	A	221	ASP	CB-CG-OD1	5.47	123.22	118.30
1	A	58	ASP	CB-CG-OD2	-5.41	113.44	118.30
1	A	290	PRO	N-CA-CB	5.34	109.71	103.30
1	A	317	THR	CA-CB-CG2	-5.33	104.94	112.40
1	A	50	LYS	CB-CA-C	5.23	120.85	110.40
1	A	132	PRO	N-CD-CG	5.21	111.02	103.20
1	A	213	ASP	CB-CG-OD1	5.14	122.93	118.30
1	A	238	ASP	CB-CG-OD1	5.08	122.87	118.30
1	A	38	ARG	NE-CZ-NH1	5.03	122.81	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	131	ASN	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	PRO	Mainchain,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	2635	0	2562	108	0
2	A	2	0	0	0	0
3	A	44	0	26	2	0
4	A	39	0	25	7	0
5	A	10	0	14	6	0
6	A	419	0	0	14	0
All	All	3149	0	2627	112	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 21.

All (112) 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:124:SER:OG	1:A:126[B]:THR:HG22	1.64	0.97
4:A:341:UD1:H8'3	6:A:710:HOH:O	1.64	0.95
1:A:296:LEU:HA	6:A:711:HOH:O	1.74	0.84
1:A:184:HIS:CG	1:A:185:PRO:HD2	2.13	0.83
1:A:133:LYS:O	1:A:140:PHE:HE2	1.61	0.83
1:A:194:GLN:HA	1:A:194:GLN:NE2	1.92	0.82
1:A:25:HIS:O	1:A:50:LYS:HE2	1.83	0.77
1:A:45:GLU:HG3	1:A:50:LYS:O	1.86	0.76
1:A:181:VAL:HG23	1:A:321:MET:HB3	1.69	0.75
1:A:130:ASP:C	1:A:132:PRO:HD2	2.08	0.73
1:A:132:PRO:HB3	1:A:136:TYR:HE1	1.54	0.73
1:A:40:VAL:HG23	1:A:336:TYR:CE1	2.24	0.72
1:A:194:GLN:HA	1:A:194:GLN:HE21	1.55	0.70
1:A:292:ARG:HG3	1:A:293:GLU:H	1.56	0.70
1:A:130:ASP:C	1:A:132:PRO:CD	2.60	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:LYS:O	1:A:140:PHE:CE2	2.44	0.69
1:A:296:LEU:HB3	1:A:297:PRO:HD2	1.74	0.69
1:A:131:ASN:N	1:A:132:PRO:HD3	2.09	0.68
1:A:124:SER:OG	1:A:126[C]:THR:HG23	1.95	0.67
1:A:253:LYS:HD2	5:A:342:PGE:H52	1.75	0.67
1:A:253:LYS:NZ	5:A:342:PGE:O1	2.27	0.67
1:A:37:LYS:NZ	1:A:338:ASP:OD1	2.17	0.66
1:A:131:ASN:N	1:A:132:PRO:CD	2.57	0.66
1:A:232[A]:ASP:OD1	6:A:637:HOH:O	2.14	0.65
1:A:50:LYS:HG2	1:A:51:HIS:N	2.10	0.65
1:A:292:ARG:HG2	1:A:295:ASP:OD2	1.97	0.64
1:A:292:ARG:NH2	4:A:341:UD1:H8'2	2.11	0.64
1:A:317:THR:OG1	1:A:320:GLU:HG3	1.99	0.62
1:A:134:ILE:HD13	1:A:299:TYR:HB3	1.81	0.62
1:A:253:LYS:HE3	5:A:342:PGE:C5	2.30	0.62
1:A:40:VAL:HG23	1:A:336:TYR:CZ	2.36	0.61
1:A:21:LEU:HD22	1:A:50:LYS:HD3	1.84	0.59
1:A:30:LEU:HD13	1:A:30:LEU:C	2.23	0.59
1:A:232[A]:ASP:HB2	1:A:300:TRP:CB	2.33	0.59
4:A:341:UD1:H4'	6:A:719:HOH:O	2.02	0.58
1:A:307:ASP:O	1:A:311:ASN:HA	2.03	0.58
1:A:253:LYS:HE3	5:A:342:PGE:H5	1.85	0.58
1:A:214:SER:OG	1:A:287:HIS:ND1	2.29	0.56
1:A:86:VAL:HG23	4:A:341:UD1:O3'	2.05	0.56
1:A:282:LYS:HB2	1:A:283:PRO:HD2	1.87	0.56
1:A:292:ARG:NH2	4:A:341:UD1:C8'	2.69	0.55
1:A:50:LYS:O	1:A:52:PRO:HD3	2.06	0.55
1:A:146:GLN:NE2	6:A:755:HOH:O	2.33	0.55
1:A:193:PRO:HB3	6:A:523:HOH:O	2.07	0.54
4:A:341:UD1:H6	4:A:341:UD1:O5B	2.08	0.54
1:A:125:ALA:HB3	1:A:178:PHE:CE2	2.42	0.54
1:A:39:SER:OG	1:A:337:PRO:HD2	2.07	0.53
1:A:132:PRO:HB3	1:A:136:TYR:CE1	2.41	0.53
1:A:134:ILE:HG12	1:A:136:TYR:CE1	2.44	0.53
1:A:292:ARG:HG3	1:A:293:GLU:N	2.22	0.52
1:A:181:VAL:CG2	1:A:321:MET:HB3	2.39	0.52
1:A:84:LYS:HE3	3:A:340:NAD:O3	2.09	0.52
1:A:220:ASN:HB3	1:A:288:PHE:CD2	2.44	0.52
1:A:200:LEU:O	1:A:204:ILE:HG13	2.10	0.52
1:A:178:PHE:O	1:A:180:PRO:HD3	2.10	0.51
1:A:234:ILE:O	1:A:234:ILE:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ALA:HB1	1:A:178:PHE:CZ	2.45	0.51
1:A:130:ASP:O	1:A:132:PRO:HD2	2.11	0.50
1:A:124:SER:CB	1:A:126[B]:THR:HG22	2.43	0.49
1:A:220:ASN:HB2	6:A:596:HOH:O	2.12	0.49
1:A:125:ALA:CB	1:A:178:PHE:CZ	2.96	0.49
1:A:188:ASP:OD2	1:A:331:ARG:NH2	2.46	0.48
1:A:21:LEU:CD2	1:A:50:LYS:HD3	2.42	0.48
1:A:86:VAL:N	6:A:750:HOH:O	2.46	0.48
1:A:222:TYR:CE1	1:A:291:ARG:HD3	2.49	0.48
1:A:104:THR:O	1:A:108:ILE:HG12	2.14	0.48
1:A:164:GLN:HB2	1:A:172:ILE:HD12	1.96	0.47
1:A:132:PRO:HB3	1:A:134:ILE:HG13	1.96	0.47
1:A:130:ASP:C	1:A:132:PRO:HD3	2.33	0.47
1:A:134:ILE:HD13	1:A:299:TYR:CB	2.44	0.47
1:A:206:GLN:HA	1:A:209:VAL:HG22	1.97	0.46
1:A:205:ALA:O	1:A:209:VAL:HG13	2.16	0.46
1:A:169:ASP:OD2	1:A:169:ASP:N	2.47	0.46
1:A:133:LYS:HB2	1:A:140:PHE:CZ	2.50	0.46
1:A:331:ARG:NH1	6:A:742:HOH:O	2.45	0.45
1:A:111:MET:HE2	1:A:170:TRP:CH2	2.51	0.45
1:A:136:TYR:HA	1:A:140:PHE:CE2	2.51	0.45
1:A:296:LEU:HB3	1:A:297:PRO:CD	2.45	0.45
1:A:121:PHE:CZ	1:A:153:LYS:HE3	2.52	0.45
1:A:228:THR:OG1	1:A:268:SER:HB2	2.17	0.45
1:A:183:ALA:HB1	1:A:189:MET:O	2.16	0.44
1:A:84:LYS:HA	1:A:149:TYR:CE2	2.52	0.44
1:A:178:PHE:N	6:A:648:HOH:O	2.24	0.44
1:A:184:HIS:CD2	1:A:185:PRO:HD2	2.52	0.44
1:A:232[A]:ASP:HB2	1:A:300:TRP:HB2	2.00	0.44
1:A:39:SER:CB	1:A:337:PRO:HD2	2.47	0.44
1:A:184:HIS:ND1	1:A:185:PRO:HD2	2.32	0.43
1:A:134:ILE:HG23	1:A:135:PRO:HA	2.00	0.43
1:A:134:ILE:HA	1:A:135:PRO:C	2.39	0.43
1:A:162:ASP:O	1:A:163:LEU:C	2.56	0.43
1:A:1:MET:HE2	1:A:76:THR:HG22	2.01	0.43
1:A:50:LYS:HG2	1:A:51:HIS:H	1.83	0.42
1:A:179:ASN:HA	1:A:180:PRO:HD2	1.82	0.42
1:A:151:LYS:O	1:A:155:MET:HG3	2.20	0.42
1:A:111:MET:HB3	1:A:111:MET:HE2	1.83	0.42
3:A:340:NAD:H4N	4:A:341:UD1:H6'2	2.02	0.42
1:A:327:HIS:O	1:A:331:ARG:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ALA:HB3	1:A:96:TYR:OH	2.20	0.41
1:A:282:LYS:CB	1:A:283:PRO:HD2	2.50	0.41
1:A:137:VAL:O	1:A:140:PHE:HB2	2.21	0.41
1:A:15:HIS:O	1:A:19:GLN:HG2	2.21	0.41
1:A:232[A]:ASP:HB2	1:A:300:TRP:HB3	2.02	0.41
1:A:250:LEU:O	1:A:253:LYS:HB2	2.21	0.41
1:A:51:HIS:HA	1:A:52:PRO:HD3	1.70	0.41
1:A:126[A]:THR:HG23	6:A:368:HOH:O	2.19	0.41
1:A:294:GLY:C	6:A:454:HOH:O	2.58	0.41
1:A:332:HIS:HA	1:A:333:PRO:HD2	1.89	0.41
1:A:216:ALA:HB3	6:A:704:HOH:O	2.21	0.41
1:A:117:LYS:NZ	6:A:696:HOH:O	2.39	0.40
1:A:137:VAL:HG12	1:A:139:SER:H	1.87	0.40
1:A:253:LYS:HZ3	5:A:342:PGE:H22	1.87	0.40
1:A:253:LYS:HE3	5:A:342:PGE:H52	2.00	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	339/338 (100%)	326 (96%)	7 (2%)	6 (2%)	8 2

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	133	LYS
1	A	293	GLU
1	A	131	ASN
1	A	11	TYR
1	A	132	PRO
1	A	292	ARG

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	285/282 (101%)	273 (96%)	12 (4%)	30	20

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	40	VAL
1	A	50	LYS
1	A	84	LYS
1	A	122	SER
1	A	130	ASP
1	A	165	LYS
1	A	253	LYS
1	A	282	LYS
1	A	292	ARG
1	A	293	GLU
1	A	308	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
1	A	158	GLN
1	A	179	ASN
1	A	194	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 [i](#)

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](#)

Of 5 ligands modelled in this entry, 2 are 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$
5	PGE	A	342	-	9,9,9	0.46	0	8,8,8	1.74	2 (25%)
3	NAD	A	340	-	42,48,48	1.92	9 (21%)	50,73,73	1.93	10 (20%)
4	UD1	A	341	-	34,41,41	2.01	8 (23%)	45,62,62	1.80	10 (22%)

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
5	PGE	A	342	-	-	4/7/7/7	-
3	NAD	A	340	-	-	4/26/62/62	0/5/5/5
4	UD1	A	341	-	-	7/24/63/63	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	340	NAD	C4N-C3N	5.99	1.49	1.39
4	A	341	UD1	C4-N3	5.40	1.42	1.33
3	A	340	NAD	C5N-C4N	5.16	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	340	NAD	C2N-C3N	4.93	1.46	1.39
4	A	341	UD1	C6-N1	4.46	1.41	1.35
4	A	341	UD1	C2B-C1B	4.40	1.60	1.53
3	A	340	NAD	C2N-N1N	4.37	1.40	1.35
4	A	341	UD1	O5'-C1'	3.79	1.51	1.41
3	A	340	NAD	C6N-C5N	-3.27	1.31	1.38
4	A	341	UD1	PB-O1'	3.05	1.68	1.60
3	A	340	NAD	C6N-N1N	2.92	1.42	1.35
4	A	341	UD1	C3'-C2'	2.44	1.57	1.53
4	A	341	UD1	O5'-C5'	2.39	1.50	1.44
4	A	341	UD1	O2'-C2B	2.28	1.48	1.43
3	A	340	NAD	C2D-C1D	-2.23	1.50	1.53
3	A	340	NAD	C2A-N1A	2.18	1.38	1.33
3	A	340	NAD	C7N-N7N	2.06	1.36	1.33

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	340	NAD	C5N-C4N-C3N	-6.99	112.07	120.34
3	A	340	NAD	C6N-N1N-C2N	-5.90	116.59	121.97
3	A	340	NAD	C3N-C7N-N7N	4.80	123.51	117.75
4	A	341	UD1	C1'-O5'-C5'	4.54	122.60	113.69
4	A	341	UD1	C2'-N2'-C7'	-4.05	113.34	123.18
4	A	341	UD1	C5-C4-N3	-3.84	114.87	123.31
4	A	341	UD1	O3A-PB-O1'	3.72	109.99	102.48
5	A	342	PGE	C5-O3-C4	3.59	128.82	113.29
4	A	341	UD1	O2'-C2B-C3B	-3.39	100.86	111.82
5	A	342	PGE	C3-O2-C2	3.19	127.12	113.29
3	A	340	NAD	C2N-C3N-C4N	-3.17	114.67	118.26
3	A	340	NAD	C5A-C6A-N6A	3.15	125.14	120.35
3	A	340	NAD	O7N-C7N-C3N	-2.77	116.31	119.63
3	A	340	NAD	C1B-N9A-C4A	-2.56	122.14	126.64
3	A	340	NAD	C6N-C5N-C4N	2.53	123.11	119.44
4	A	341	UD1	O4B-C1B-C2B	2.52	110.61	106.93
4	A	341	UD1	C8'-C7'-N2'	2.40	120.16	116.10
4	A	341	UD1	C3'-C2'-N2'	2.37	115.10	110.62
3	A	340	NAD	C2N-N1N-C1D	2.37	124.41	119.14
4	A	341	UD1	O5'-C1'-C2'	2.32	115.10	110.58
4	A	341	UD1	C1'-C2'-N2'	2.07	114.56	111.00
3	A	340	NAD	C5N-C6N-N1N	2.02	123.29	120.40

There are no chirality outliers.

All (15) torsion outliers are listed below:

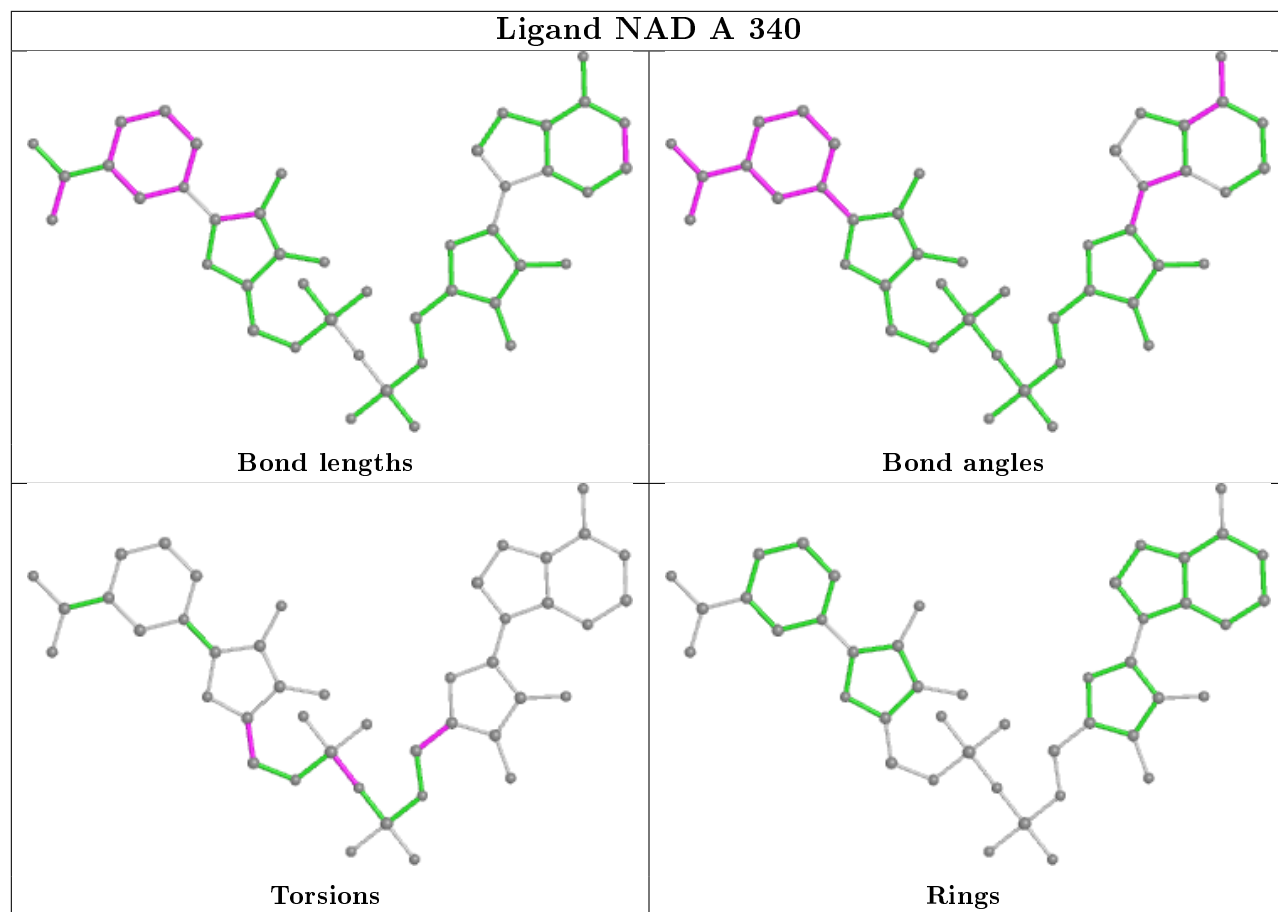
Mol	Chain	Res	Type	Atoms
3	A	340	NAD	O4D-C4D-C5D-O5D
3	A	340	NAD	C3D-C4D-C5D-O5D
4	A	341	UD1	O4B-C1B-N1-C6
4	A	341	UD1	O5'-C5'-C6'-O6'
4	A	341	UD1	C8'-C7'-N2'-C2'
5	A	342	PGE	O2-C3-C4-O3
4	A	341	UD1	C4'-C5'-C6'-O6'
5	A	342	PGE	O1-C1-C2-O2
4	A	341	UD1	O7'-C7'-N2'-C2'
5	A	342	PGE	C3-C4-O3-C5
3	A	340	NAD	PA-O3-PN-O2N
4	A	341	UD1	C1'-O1'-PB-O3A
4	A	341	UD1	PA-O3A-PB-O1B
5	A	342	PGE	C6-C5-O3-C4
3	A	340	NAD	O4B-C4B-C5B-O5B

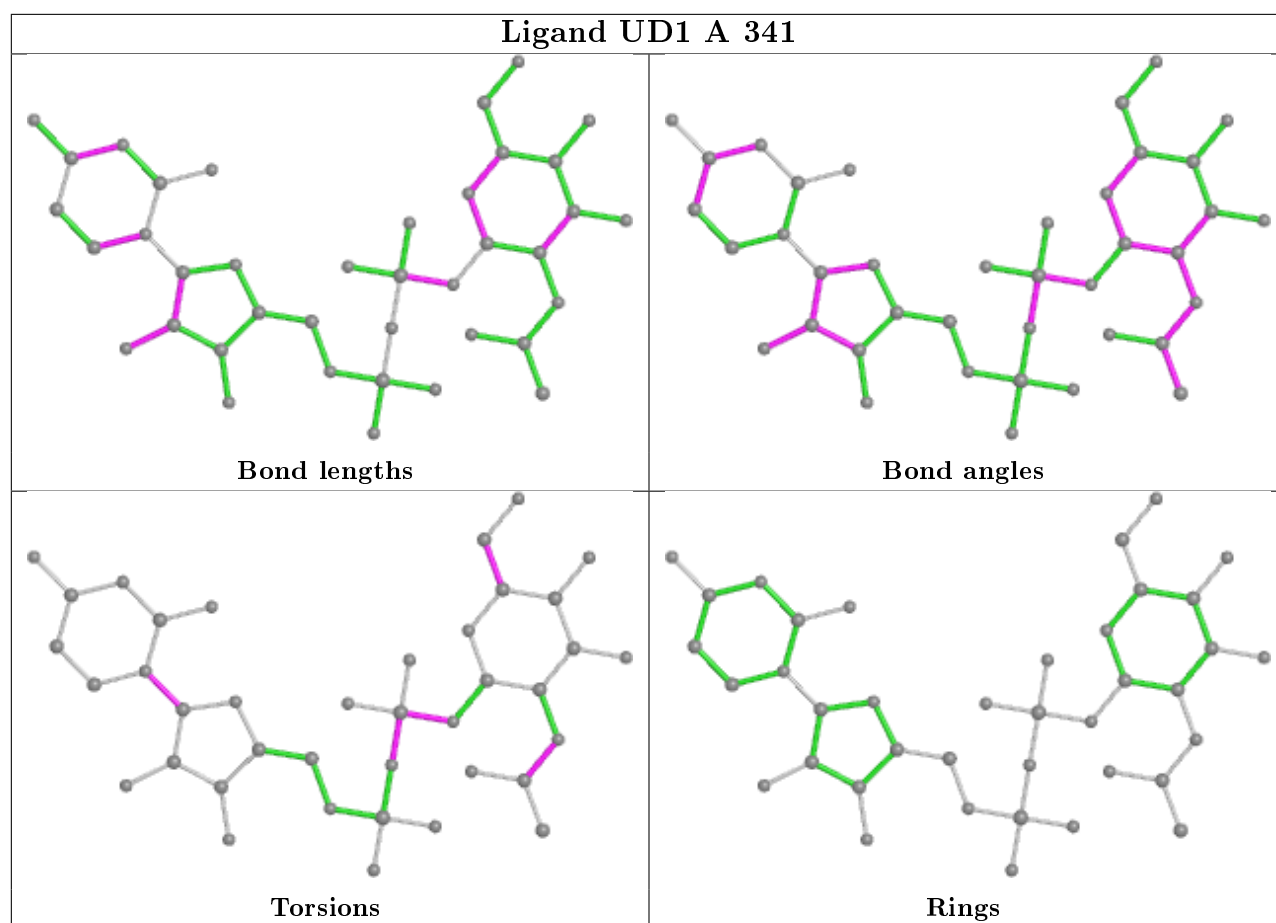
There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	342	PGE	6	0
3	A	340	NAD	2	0
4	A	341	UD1	7	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 [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, 95th 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 ²)	Q<0.9
1	A	338/338 (100%)	-0.48	3 (0%) 84 85	15, 24, 54, 100	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	ASN	3.4
1	A	132	PRO	3.3
1	A	294	GLY	2.7

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](#)

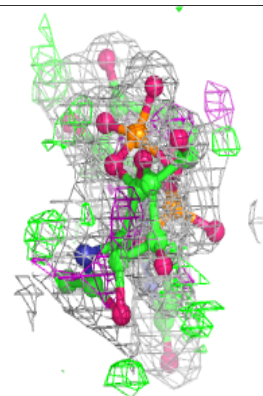
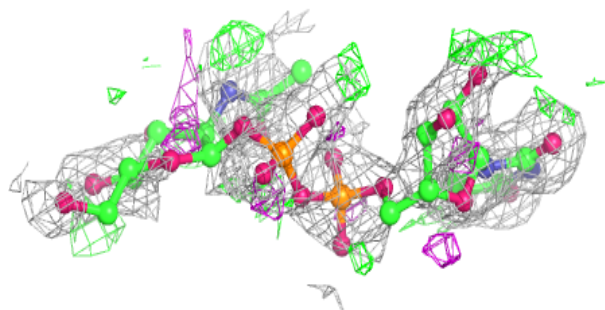
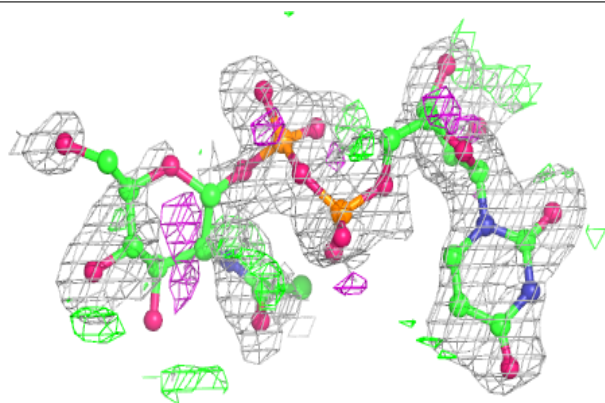
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

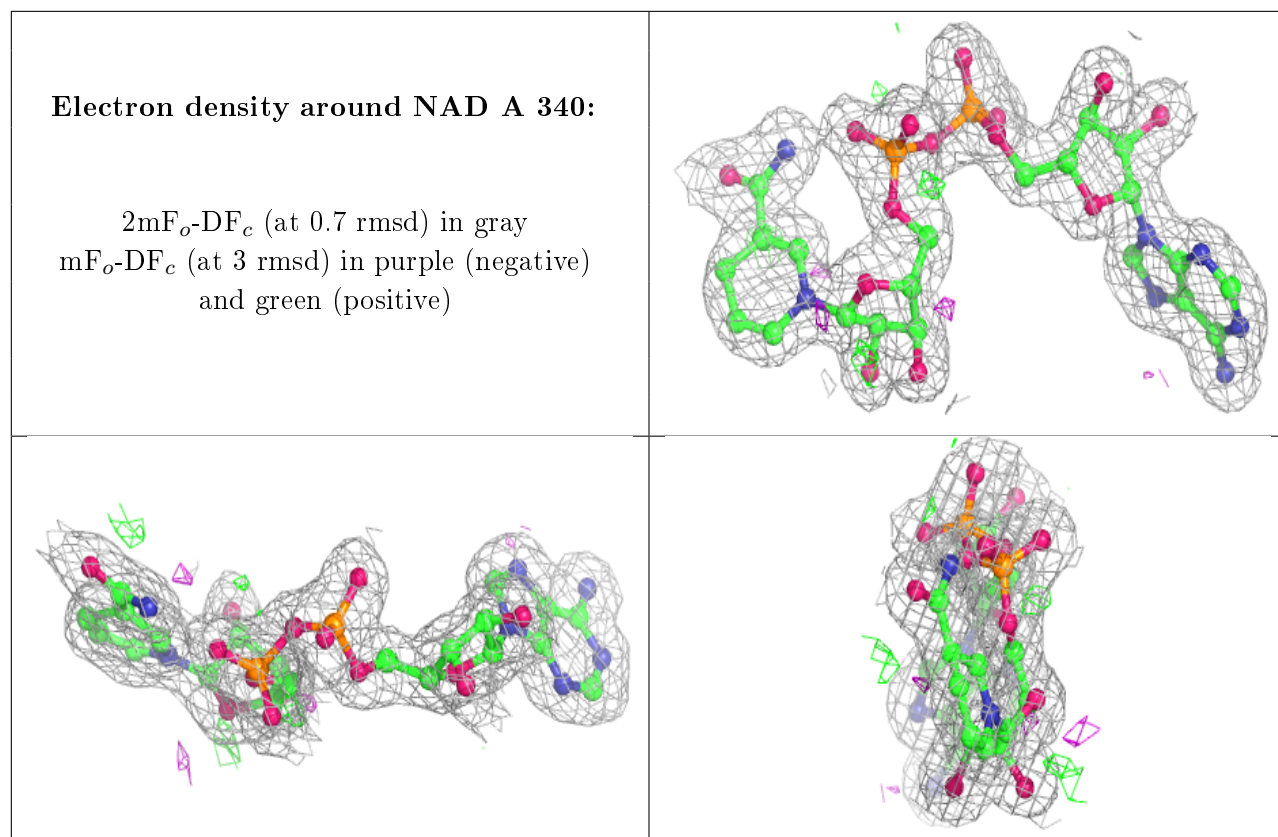
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NA	A	344	1/1	0.85	0.21	41,41,41,41	0
4	UD1	A	341	39/39	0.87	0.19	18,71,100,100	0
5	PGE	A	342	10/10	0.93	0.09	21,32,80,91	0
2	NA	A	343	1/1	0.96	0.06	32,32,32,32	0
3	NAD	A	340	44/44	0.97	0.08	13,22,33,66	0

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 UD1 A 341:

$2mF_o-DF_c$ (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.