

Full wwPDB X-ray Structure Validation Report (i)

Aug 14, 2023 – 05:14 PM EDT

PDB ID	:	1SB8
Title	:	Crystal structure of Pseudomonas aeruginosa UDP-N-acetylglucosamine
		4-epimerase complexed with UDP-N-acetylgalactosamine
Authors	:	Ishiyama, N.; Creuzenet, C.; Lam, J.S.; Berghuis, A.M.
Deposited on		
Resolution	:	2.10 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

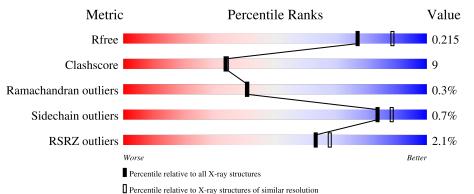
Xtriage (Phenix) EDS buster-report Percentile statistics Refmac CCP4 Ideal geometry (proteins) Ideal geometry (DNA, RNA)	::	20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove) Engh & Huber (2001) Parkinson et al. (1996)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.10 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5197(2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 of 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 <=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	А	352	2% 79 %	18%	·



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2927 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 wbpP.

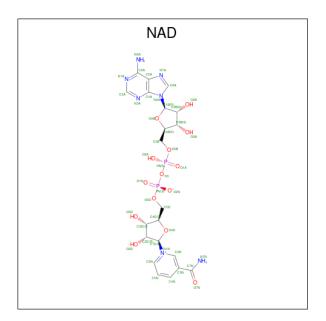
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	341	Total 2662	C 1688	N 460	O 504	S 10	0	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	-10	MET	-	initiating methionine	GB 20560072
А	-9	HIS	-	expression tag	GB 20560072
А	-8	HIS	-	expression tag	GB 20560072
А	-7	HIS	-	expression tag	GB 20560072
А	-6	HIS	-	expression tag	GB 20560072
А	-5	HIS	-	expression tag	GB 20560072
А	-4	HIS	-	expression tag	GB 20560072
А	-3	GLY	-	cloning artifact	GB 20560072
А	-2	SER	-	cloning artifact	GB 20560072
А	-1	MET	-	cloning artifact	GB 20560072
А	0	GLY	-	cloning artifact	GB 20560072

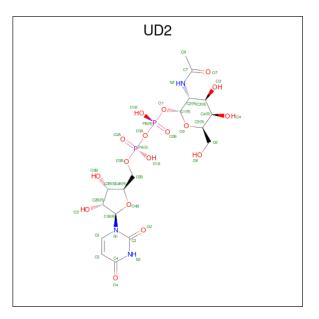
• Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
9	Δ	1	Total	С	Ν	Ο	Р	0	0
	A	1	44	21	7	14	2	0	0

• Molecule 3 is URIDINE-DIPHOSPHATE-N-ACETYLGALACTOSAMINE (three-letter code: UD2) (formula: $C_{17}H_{27}N_3O_{17}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	А	1	Total 39				Р 2	0	0

• Molecule 4 is water.

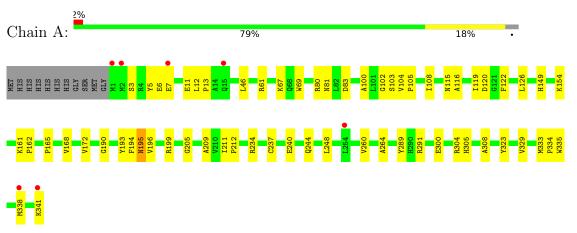


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	182	Total O 182 182	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 (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: wbpP



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	60.44Å 95.88Å 139.64Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.25 - 2.10	Depositor
Resolution (A)	48.01 - 2.10	EDS
% Data completeness	96.3 (41.25-2.10)	Depositor
(in resolution range)	96.4 (48.01-2.10)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	0.08	Depositor
$< I/\sigma(I) > 1$	$5.79 (at 2.10 \text{\AA})$	Xtriage
Refinement program	CNS 1.1	Depositor
D D.	0.190 , 0.228	Depositor
R, R_{free}	0.176 , 0.215	DCC
R_{free} test set	2421 reflections (10.06%)	wwPDB-VP
Wilson B-factor $(Å^2)$	18.3	Xtriage
Anisotropy	1.066	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.40, 65.0	EDS
L-test for twinning ²	$ < L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2927	wwPDB-VP
Average B, all atoms $(Å^2)$	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.41% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: NAD, UD2

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).

Mal	Chain	Bond	lengths	Bond angles		
	Unain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.32	0/2720	0.55	0/3691	

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	А	2662	0	2607	45	0
2	А	44	0	26	3	0
3	А	39	0	25	1	0
4	А	182	0	0	7	0
All	All	2927	0	2658	46	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 9.

All (46) 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:122:PHE:O	1:A:126:LEU:HD13	1.93	0.69
1:A:199:ARG:HG3	4:A:456:HOH:O	1.91	0.69
1:A:300:GLU:HG2	4:A:460:HOH:O	1.93	0.68
1:A:122:PHE:CZ	1:A:126:LEU:HD11	2.32	0.64
1:A:3:SER:OG	1:A:6:GLU:HG3	1.96	0.64
1:A:338:MET:O	1:A:341:LYS:HG2	1.98	0.63
1:A:80:ARG:HD2	4:A:435:HOH:O	2.00	0.61
1:A:193:TYR:HB2	2:A:342:NAD:C5N	2.30	0.61
1:A:115:ASN:OD1	1:A:119:ILE:HD12	2.02	0.60
1:A:338:MET:HA	1:A:341:LYS:HD3	1.84	0.59
1:A:194:PHE:CZ	1:A:308:ALA:HB2	2.38	0.58
1:A:104:VAL:HB	1:A:105:PRO:HD3	1.85	0.58
2:A:342:NAD:C4N	3:A:343:UD2:H4'	2.34	0.58
1:A:116:ALA:O	1:A:120:ASP:HB2	2.08	0.52
1:A:108:ILE:HD13	1:A:165:PRO:HD2	1.91	0.52
1:A:168:VAL:O	1:A:172:VAL:HG23	2.11	0.51
1:A:108:ILE:HD13	1:A:165:PRO:CD	2.41	0.50
1:A:194:PHE:CE2	1:A:308:ALA:HB2	2.45	0.50
1:A:211:ILE:HB	1:A:212:PRO:HD3	1.93	0.50
1:A:102:GLY:O	1:A:103:SER:HB3	2.12	0.49
1:A:194:PHE:HB2	1:A:264:ALA:HB2	1.95	0.49
1:A:7:GLU:O	1:A:11:GLU:HG3	2.13	0.48
1:A:289:TYR:CZ	1:A:291:ARG:HB2	2.49	0.48
1:A:199:ARG:HG2	1:A:335:TRP:CB	2.45	0.46
1:A:329:VAL:O	1:A:333:MET:HG2	2.16	0.45
1:A:108:ILE:HG21	1:A:300:GLU:OE2	2.16	0.45
1:A:67:LYS:HG2	4:A:369:HOH:O	2.16	0.45
1:A:196:VAL:HA	1:A:237:CYS:O	2.18	0.43
1:A:199:ARG:HG2	1:A:335:TRP:HB2	2.00	0.43
1:A:61:ARG:HD3	1:A:69:TRP:CE2	2.54	0.43
1:A:5:TYR:CG	1:A:244:GLN:HG3	2.54	0.42
1:A:81:ASN:OD1	1:A:83:ASP:HB2	2.17	0.42
1:A:122:PHE:CE2	1:A:126:LEU:HD11	2.54	0.42
1:A:161:LYS:HA	1:A:162:PRO:HD3	1.93	0.42
1:A:190:GLY:O	1:A:260:VAL:HA	2.20	0.42
1:A:100:ALA:HB3	2:A:342:NAD:O3D	2.20	0.41
1:A:240:GLU:HG3	4:A:447:HOH:O	2.19	0.41
1:A:304:ARG:HG3	1:A:305:HIS:CD2	2.55	0.41
1:A:205:GLY:HA3	4:A:485:HOH:O	2.19	0.41
1:A:209:ALA:O	1:A:212:PRO:HD2	2.20	0.41
1:A:333:MET:HB2	1:A:334:PRO:HD3	2.03	0.41
1:A:12:LEU:HB2	1:A:13:PRO:HD3	2.01	0.41

Continued on next page...



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:ASN:OD1	1:A:234:ARG:HD2	2.21	0.41
1:A:260:VAL:HG22	4:A:377:HOH:O	2.20	0.41
1:A:149:HIS:O	1:A:154:LYS:NZ	2.52	0.40
1:A:240:GLU:HG2	1:A:323:TYR:CE2	2.55	0.40

Continued from previous page...

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	А	339/352~(96%)	328~(97%)	10 (3%)	1 (0%)	41 41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	195	ASN

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.

Μ	ol	Chain	Analysed	Analysed Rotameric		Percentiles	
1	L	А	281/290~(97%)	279~(99%)	2(1%)	84 88	

All (2) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	А	46	LEU
1	А	248	LEU

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

Mol	Chain	Res	Type
1	А	149	HIS

5.3.3 RNA (i)

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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).

Mal	Turne	Chain	Dec	Link	B	ond leng	gths	B	ond ang	les
Mol	Type	Chain	Res	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	NAD	А	342	-	42,48,48	1.60	11 (26%)	50,73,73	1.82	8 (16%)
3	UD2	А	343	-	38,41,41	1.07	3 (7%)	57,62,62	1.32	2 (3%)

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.



Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	А	342	-	-	8/26/62/62	0/5/5/5
3	UD2	А	343	-	-	5/26/63/63	0/3/3/3

'-' means no outliers of that kind were identified.

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	342	NAD	C2D-C1D	-3.46	1.48	1.53
3	А	343	UD2	C5-C4	3.23	1.50	1.43
2	А	342	NAD	C2A-N1A	3.22	1.39	1.33
3	А	343	UD2	PA-O2A	3.08	1.61	1.50
2	А	342	NAD	PN-O1N	3.02	1.61	1.50
3	А	343	UD2	PB-O2B	2.96	1.61	1.50
2	А	342	NAD	PA-O1A	2.88	1.61	1.50
2	А	342	NAD	C2N-N1N	2.62	1.38	1.35
2	А	342	NAD	C4N-C3N	2.62	1.43	1.39
2	А	342	NAD	C7N-N7N	2.58	1.37	1.33
2	А	342	NAD	C4A-N3A	2.52	1.39	1.35
2	А	342	NAD	C2A-N3A	2.45	1.36	1.32
2	А	342	NAD	C5N-C4N	2.27	1.43	1.38
2	А	342	NAD	C2B-C1B	-2.13	1.50	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
3	А	343	UD2	O5'-C1'-O1'	-7.56	101.48	111.36
2	А	342	NAD	C6N-N1N-C2N	7.42	128.74	121.97
2	А	342	NAD	C3N-C2N-N1N	-5.27	115.28	120.43
2	А	342	NAD	C3N-C7N-N7N	-3.18	113.93	117.75
2	А	342	NAD	O7N-C7N-C3N	3.06	123.29	119.63
2	А	342	NAD	N6A-C6A-N1A	2.71	124.20	118.57
2	А	342	NAD	C5N-C6N-N1N	-2.69	116.54	120.40
2	А	342	NAD	C1B-N9A-C4A	-2.66	121.97	126.64
3	А	343	UD2	O1B-PB-O1'	2.21	115.49	106.78
2	А	342	NAD	O3D-C3D-C4D	-2.06	105.09	111.05

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms			
2	А	342	NAD	C5D-O5D-PN-O1N			
 Continued on most maga							

Continued on next page...



Mol	Chain	Res	Type	Atoms
3	А	343	UD2	C8'-C7'-N2'-C2'
3	А	343	UD2	O7'-C7'-N2'-C2'
2	А	342	NAD	PN-O3-PA-O5B
3	А	343	UD2	PB-O3A-PA-O5B
2	А	342	NAD	C5D-O5D-PN-O3
2	А	342	NAD	C5B-O5B-PA-O1A
2	А	342	NAD	C5B-O5B-PA-O2A
2	А	342	NAD	C5D-O5D-PN-O2N
3	А	343	UD2	O5'-C1'-O1'-PB
2	А	342	NAD	O4B-C4B-C5B-O5B
2	А	342	NAD	PA-O3-PN-O2N
3	А	343	UD2	C4'-C5'-C6'-O6'

Continued from previous page...

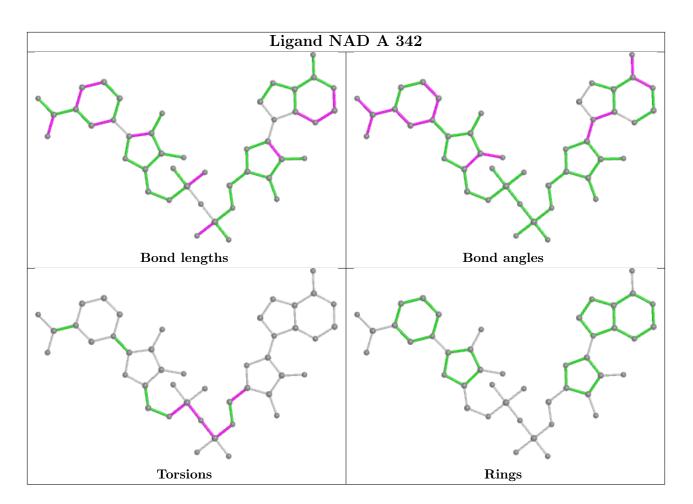
There are no ring outliers.

2 monomers are involved in 3 short contacts:

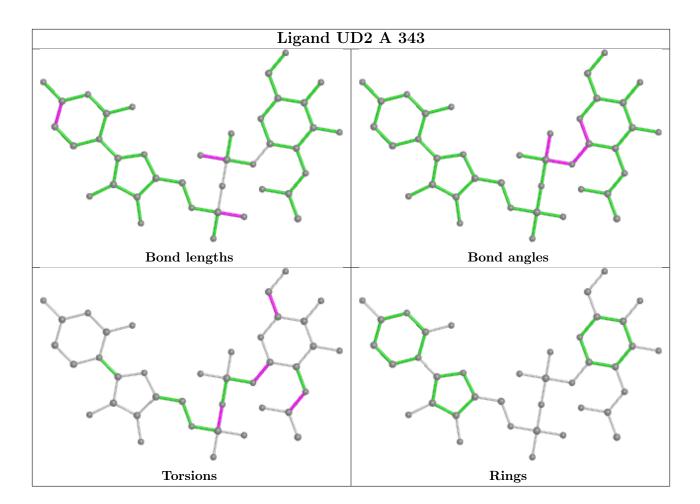
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	342	NAD	3	0
3	А	343	UD2	1	0

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 then 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 sufficient 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, 95^{th} 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	А	341/352~(96%)	0.17	7 (2%)	63	68	13, 20, 35, 65	0

All (7) RSRZ outliers are listed below:

Mol	Chain Re		Type	RSRZ	
1	А	1	MET	5.8	
1	А	254	LEU	3.8	
1	А	341	LYS	3.2	
1	А	15	GLN	2.4	
1	А	338	MET	2.4	
1	А	2	MET	2.3	
1	А	7	GLU	2.0	

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 monosaccharides 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, 95^{th} 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	NAD	А	342	44/44	0.95	0.17	$13,\!17,\!40,\!40$	0

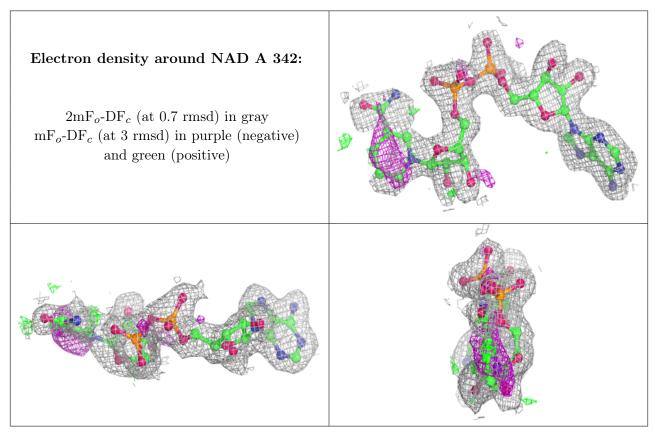
Continued on next page...



Continued from previous page...

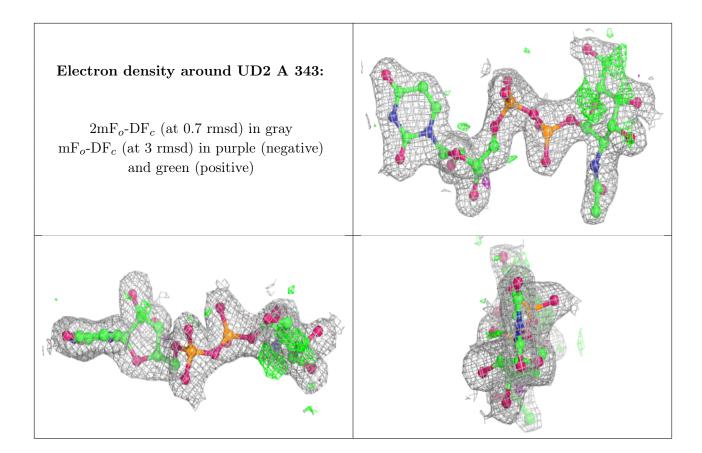
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
3	UD2	А	343	39/39	0.95	0.16	15,20,25,25	14

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.









6.5 Other polymers (i)

There are no such residues in this entry.

