



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 1, 2026 – 04:08 PM EDT

PDB ID : 36CA / pdb_000036ca
Title : Dehydrogenase AprD5 complexed with NAD and UDP-beta-D-Glucose
Authors : Kim, W.; Zhang, Y.J.
Deposited on : 2026-05-29
Resolution : 1.87 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.2
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.2
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.51

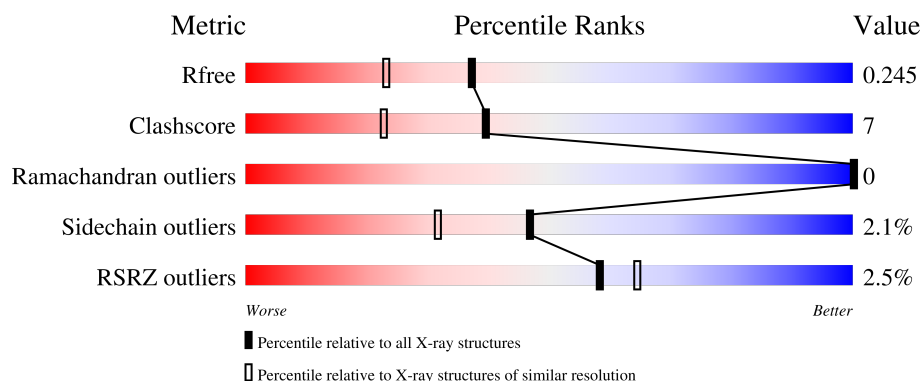
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

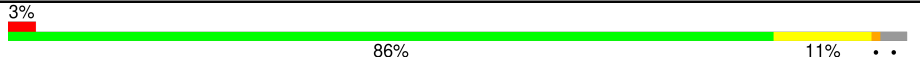
The reported resolution of this entry is 1.87 Å.

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}	180053	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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 $\leq 5\%$. 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	348	
1	B	348	

2 Entry composition [i](#)

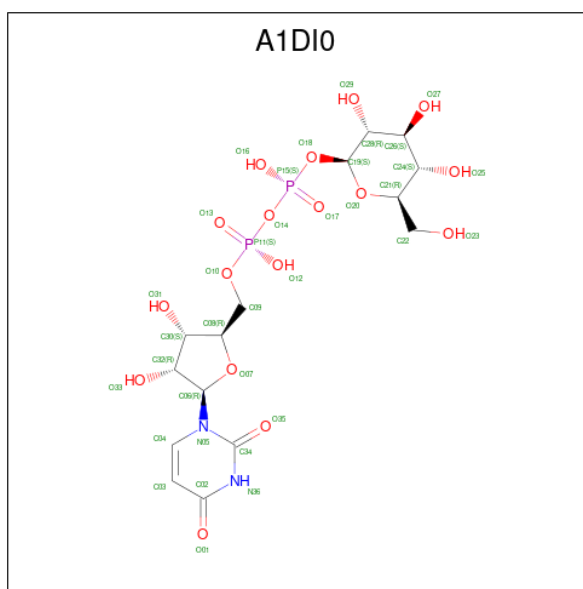
There are 4 unique types of molecules in this entry. The entry contains 6031 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nucleoside-diphosphate-sugar epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	0	0
			2642	1642	486	509	5			
1	B	339	Total	C	N	O	S	0	0	0
			2642	1642	486	509	5			

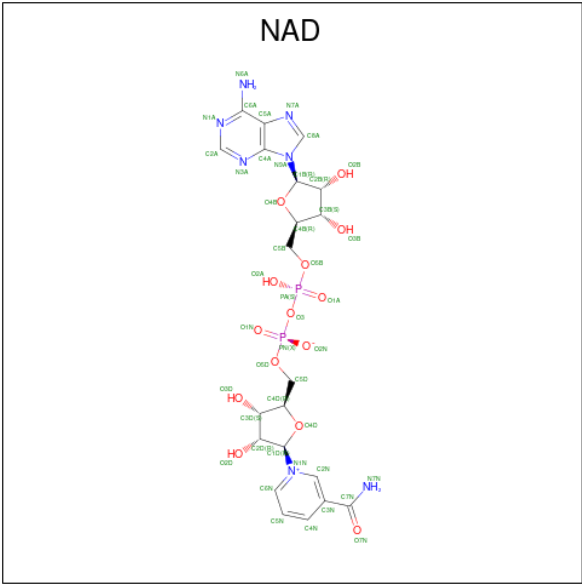
- Molecule 2 is [(2R,3S,4R,5R)-5-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)-3,4-dihydroxyoxolan-2-yl]methyl (2S,3R,4S,5S,6R)-3,4,5-trihydroxy-6-(hydroxymethyl)oxan-2-yl dihydrogen diphosphate (CCD ID: A1DI0) (formula: C₁₅H₂₄N₂O₁₇P₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula:

C₂₁H₂₇N₇O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



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

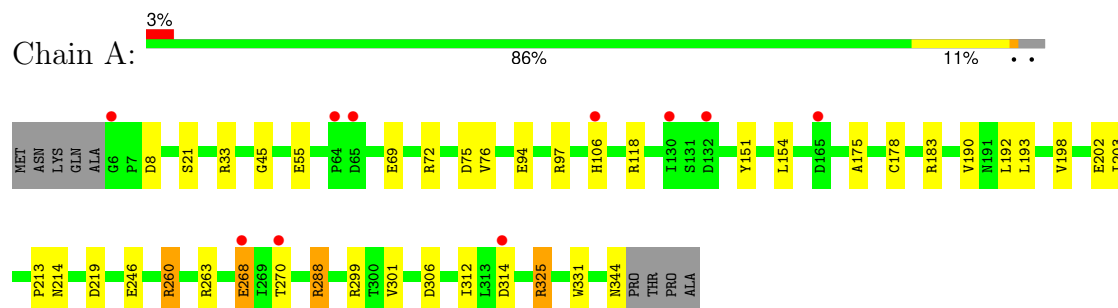
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	298	Total	O	0	0
			298	298		
4	B	289	Total	O	0	0
			289	289		

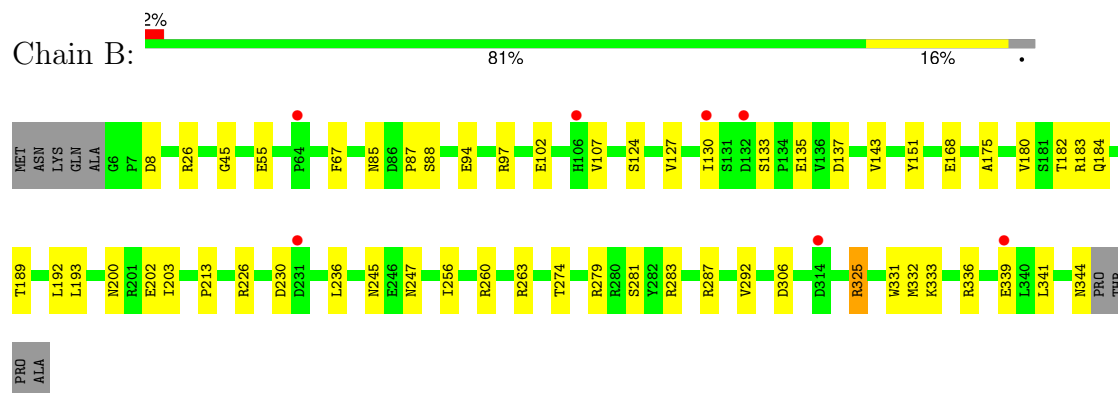
3 Residue-property plots [i](#)

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: Nucleoside-diphosphate-sugar epimerase



- Molecule 1: Nucleoside-diphosphate-sugar epimerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.10Å 60.33Å 71.03Å 70.40° 72.30° 89.93°	Depositor
Resolution (Å)	35.46 – 1.87 35.46 – 1.87	Depositor EDS
% Data completeness (in resolution range)	97.4 (35.46-1.87) 97.4 (35.46-1.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 1.87Å)	Xtriage
Refinement program	PHENIX (1.18rc3_3805: ???)	Depositor
R, R_{free}	0.215 , 0.245 0.215 , 0.245	Depositor DCC
R_{free} test set	2009 reflections (2.93%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.328	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6031	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NAD, A1DI0

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	0.32	0/2685	0.53	0/3656
1	B	0.37	1/2685 (0.0%)	0.52	0/3656
All	All	0.35	1/5370 (0.0%)	0.53	0/7312

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	245	ASN	CA-C	5.92	1.55	1.52

There are no bond angle outliers.

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	2642	0	2636	34	0
1	B	2642	0	2636	41	0
2	A	36	0	0	0	0
2	B	36	0	0	1	0
3	A	44	0	25	0	0
3	B	44	0	25	0	0
4	A	298	0	0	19	3
4	B	289	0	0	17	3

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6031	0	5322	72	3

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 7.

All (72) 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:214:ASN:HB2	4:A:515:HOH:O	1.76	0.85
1:A:314:ASP:OD1	4:A:501:HOH:O	1.97	0.83
1:A:75:ASP:O	4:A:502:HOH:O	2.02	0.78
1:A:33:ARG:NH1	4:A:504:HOH:O	2.20	0.74
1:A:268:GLU:OE2	4:A:503:HOH:O	2.07	0.73
1:A:76:VAL:HG22	4:A:502:HOH:O	1.88	0.73
1:B:189:THR:HG23	2:B:401:A1DI0:O17	1.89	0.72
1:B:133:SER:O	1:B:283:ARG:NH1	2.23	0.71
1:A:55:GLU:OE2	4:A:504:HOH:O	2.09	0.70
1:A:288:ARG:NH2	4:A:508:HOH:O	2.23	0.70
1:B:97:ARG:HA	4:B:509:HOH:O	1.91	0.70
1:A:94:GLU:OE2	4:A:505:HOH:O	2.11	0.68
1:B:102:GLU:OE1	4:B:502:HOH:O	2.10	0.68
1:B:168:GLU:CD	4:B:508:HOH:O	2.42	0.63
1:B:247:ASN:HB2	4:B:717:HOH:O	1.99	0.63
1:B:151:TYR:HB3	4:B:512:HOH:O	2.01	0.61
1:B:333:LYS:HD2	4:B:523:HOH:O	1.99	0.61
1:B:130:ILE:HD12	1:B:281:SER:H	1.65	0.60
1:A:76:VAL:HA	4:A:502:HOH:O	2.01	0.59
1:B:8:ASP:OD2	4:B:503:HOH:O	2.16	0.58
1:A:193:LEU:HD22	1:A:203:ILE:HG23	1.87	0.57
1:B:168:GLU:OE1	4:B:504:HOH:O	2.17	0.57
1:B:263:ARG:NH1	4:B:510:HOH:O	2.31	0.57
1:B:135:GLU:OE1	1:B:287:ARG:NH2	2.37	0.57
1:B:135:GLU:CD	1:B:287:ARG:HH22	2.13	0.55
1:A:214:ASN:ND2	4:A:515:HOH:O	2.39	0.55
1:A:118:ARG:N	4:A:502:HOH:O	2.40	0.54
1:A:246:GLU:HB2	4:A:515:HOH:O	2.08	0.54
1:B:85:ASN:CG	1:B:87:PRO:HD2	2.33	0.52
1:B:325:ARG:HG2	1:B:331:TRP:CD2	2.44	0.52
1:A:325:ARG:HG2	1:A:331:TRP:CD2	2.45	0.52
1:A:106:HIS:CD2	4:A:517:HOH:O	2.63	0.52
1:A:118:ARG:HB3	4:A:502:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ASP:OD1	1:A:299:ARG:HD2	2.10	0.51
1:B:137:ASP:OD2	4:B:506:HOH:O	2.19	0.51
1:B:88:SER:HB3	1:B:332:MET:SD	2.50	0.51
1:B:193:LEU:HD22	1:B:203:ILE:HG23	1.94	0.50
1:B:344:ASN:C	4:B:678:HOH:O	2.54	0.50
1:B:274:THR:HG22	4:B:640:HOH:O	2.12	0.49
1:A:97:ARG:HD2	1:B:143:VAL:HG12	1.94	0.49
1:B:279:ARG:NH1	4:B:517:HOH:O	2.43	0.48
1:B:256:ILE:O	1:B:260:ARG:HG2	2.13	0.48
1:A:299:ARG:NH2	4:A:537:HOH:O	2.48	0.47
1:A:151:TYR:OH	1:B:143:VAL:HA	2.16	0.46
1:B:226:ARG:HE	1:B:230:ASP:CG	2.23	0.46
1:B:124:SER:O	1:B:127:VAL:HG22	2.16	0.46
1:B:175:ALA:HB1	1:B:213:PRO:C	2.41	0.45
1:B:274:THR:HG23	4:B:669:HOH:O	2.15	0.45
1:B:332:MET:HG2	1:B:341:LEU:HD11	1.97	0.45
1:A:263:ARG:NH1	4:A:526:HOH:O	2.44	0.45
1:B:45:GLY:O	1:B:182:THR:OG1	2.32	0.45
1:A:178:CYS:HB3	1:A:190:VAL:HB	1.99	0.45
1:A:8:ASP:O	1:A:33:ARG:HB3	2.18	0.44
1:B:339:GLU:CD	1:B:339:GLU:H	2.26	0.44
1:A:72:ARG:HA	1:A:72:ARG:HE	1.83	0.43
1:A:175:ALA:HB1	1:A:213:PRO:C	2.43	0.43
1:B:200:ASN:ND2	4:B:501:HOH:O	1.87	0.43
1:A:198:VAL:HG11	1:A:312:ILE:HG21	2.01	0.43
1:A:344:ASN:C	4:A:518:HOH:O	2.61	0.43
1:A:97:ARG:HD2	1:B:143:VAL:CG1	2.49	0.42
1:A:301:VAL:CG2	4:A:622:HOH:O	2.67	0.42
1:B:94:GLU:HB2	4:B:555:HOH:O	2.19	0.42
1:A:260:ARG:NH1	1:A:306:ASP:OD2	2.53	0.42
1:B:260:ARG:NH2	1:B:306:ASP:OD2	2.52	0.42
1:B:67:PHE:CZ	1:B:107:VAL:HG13	2.55	0.41
1:B:236:LEU:HG	1:B:292:VAL:HG11	2.02	0.41
1:B:55:GLU:HB2	4:B:706:HOH:O	2.20	0.41
1:A:21:SER:OG	1:A:45:GLY:HA3	2.21	0.41
1:A:154:LEU:HD23	1:A:154:LEU:HA	1.76	0.41
1:A:202:GLU:HG3	1:A:270:THR:HB	2.02	0.41
1:B:180:VAL:HA	1:B:184:GLN:OE1	2.21	0.40
1:B:192:LEU:C	1:B:192:LEU:HD23	2.46	0.40

All (3) 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)
4:A:506:HOH:O	4:B:684:HOH:O[1_664]	1.94	0.26
4:A:552:HOH:O	4:B:694:HOH:O[1_564]	1.94	0.26
4:A:655:HOH:O	4:B:730:HOH:O[1_664]	2.11	0.09

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	337/348 (97%)	328 (97%)	9 (3%)	0	100	100
1	B	337/348 (97%)	326 (97%)	11 (3%)	0	100	100
All	All	674/696 (97%)	654 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	291/298 (98%)	284 (98%)	7 (2%)	43	27
1	B	291/298 (98%)	286 (98%)	5 (2%)	53	40
All	All	582/596 (98%)	570 (98%)	12 (2%)	47	32

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

Mol	Chain	Res	Type
1	A	69	GLU

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Mol	Chain	Res	Type
1	A	183	ARG
1	A	192	LEU
1	A	260	ARG
1	A	268	GLU
1	A	288	ARG
1	A	325	ARG
1	B	26	ARG
1	B	183	ARG
1	B	202	GLU
1	B	325	ARG
1	B	336	ARG

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

Mol	Chain	Res	Type
1	A	106	HIS
1	A	200	ASN
1	B	214	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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
3	NAD	A	402	-	46,48,48	3.53	18 (39%)	64,73,73	1.89	17 (26%)
2	A1DI0	B	401	-	37,38,38	3.13	16 (43%)	55,58,58	1.31	6 (10%)
3	NAD	B	402	-	46,48,48	3.48	20 (43%)	64,73,73	1.69	14 (21%)
2	A1DI0	A	401	-	37,38,38	3.13	16 (43%)	55,58,58	1.56	10 (18%)

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	402	-	-	5/30/62/62	0/5/5/5
2	A1DI0	B	401	-	-	6/23/59/59	0/3/3/3
3	NAD	B	402	-	-	4/30/62/62	0/5/5/5
2	A1DI0	A	401	-	-	1/23/59/59	0/3/3/3

All (70) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	NAD	C2D-C3D	-9.76	1.26	1.53
3	B	402	NAD	C2D-C3D	-9.44	1.27	1.53
2	B	401	A1DI0	C30-C08	-8.63	1.31	1.53
3	B	402	NAD	C7N-N7N	8.39	1.48	1.33
3	A	402	NAD	O4D-C1D	8.24	1.51	1.40
3	A	402	NAD	C7N-N7N	8.24	1.48	1.33
2	A	401	A1DI0	C30-C08	-8.13	1.32	1.53
3	A	402	NAD	O4B-C1B	8.03	1.60	1.42
3	B	402	NAD	O4B-C1B	7.91	1.60	1.42
2	A	401	A1DI0	O07-C08	7.89	1.62	1.45
2	B	401	A1DI0	O07-C08	7.84	1.62	1.45
2	A	401	A1DI0	P15-O14	7.77	1.67	1.59
2	B	401	A1DI0	P15-O14	7.64	1.67	1.59
3	B	402	NAD	C2B-C1B	-7.39	1.30	1.53
3	A	402	NAD	C2B-C1B	-7.21	1.30	1.53
3	B	402	NAD	O4D-C1D	7.01	1.50	1.40
2	A	401	A1DI0	P11-O14	6.60	1.66	1.59
2	B	401	A1DI0	P11-O14	6.40	1.66	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	NAD	O4B-C4B	-6.17	1.31	1.45
3	A	402	NAD	O4B-C4B	-6.14	1.31	1.45
2	A	401	A1DI0	O07-C06	-5.63	1.29	1.42
2	B	401	A1DI0	O07-C06	-5.63	1.29	1.42
3	A	402	NAD	O4D-C4D	-5.60	1.32	1.45
3	B	402	NAD	O4D-C4D	-5.48	1.32	1.45
3	A	402	NAD	C6A-N6A	5.29	1.47	1.34
3	B	402	NAD	C6A-N6A	4.92	1.46	1.34
3	B	402	NAD	C3D-C4D	4.80	1.65	1.53
3	A	402	NAD	C3D-C4D	4.79	1.65	1.53
3	B	402	NAD	C3N-C7N	4.68	1.57	1.50
3	A	402	NAD	C3N-C7N	3.98	1.56	1.50
2	A	401	A1DI0	O31-C30	3.76	1.52	1.43
3	B	402	NAD	O2D-C2D	3.51	1.51	1.43
2	B	401	A1DI0	O31-C30	3.49	1.51	1.43
2	B	401	A1DI0	C02-N36	-3.32	1.32	1.38
2	B	401	A1DI0	C34-N05	-3.30	1.33	1.38
3	A	402	NAD	O2D-C2D	3.22	1.50	1.43
2	A	401	A1DI0	C02-N36	-3.06	1.33	1.38
3	A	402	NAD	O2B-C2B	3.04	1.50	1.43
2	B	401	A1DI0	C34-N36	-3.04	1.32	1.38
2	A	401	A1DI0	C34-N36	-2.91	1.32	1.38
2	A	401	A1DI0	C34-N05	-2.85	1.34	1.38
2	B	401	A1DI0	P15-O18	2.76	1.67	1.59
2	A	401	A1DI0	O20-C19	2.71	1.48	1.41
2	A	401	A1DI0	P15-O18	2.64	1.67	1.59
3	B	402	NAD	O3D-C3D	2.57	1.49	1.43
3	B	402	NAD	O2B-C2B	2.47	1.49	1.43
3	B	402	NAD	C5A-C4A	-2.47	1.34	1.39
3	B	402	NAD	O3B-C3B	-2.46	1.36	1.43
2	A	401	A1DI0	O01-C02	-2.39	1.19	1.24
2	B	401	A1DI0	O20-C19	2.39	1.48	1.41
3	A	402	NAD	C2N-N1N	2.37	1.37	1.35
2	A	401	A1DI0	P11-O10	2.37	1.68	1.59
3	B	402	NAD	C2N-N1N	2.33	1.37	1.35
3	B	402	NAD	PA-O3	2.31	1.62	1.59
3	A	402	NAD	C5A-C4A	-2.30	1.35	1.39
2	B	401	A1DI0	O35-C34	-2.26	1.19	1.23
3	A	402	NAD	C8A-N7A	2.25	1.36	1.31
2	B	401	A1DI0	P11-O10	2.24	1.68	1.59
2	A	401	A1DI0	O27-C26	2.24	1.48	1.43
2	A	401	A1DI0	O33-C32	-2.23	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	NAD	C4N-C3N	-2.19	1.36	1.39
2	B	401	A1DI0	O01-C02	-2.19	1.20	1.24
3	B	402	NAD	C8A-N7A	2.18	1.35	1.31
3	A	402	NAD	C5A-N7A	-2.17	1.35	1.39
2	A	401	A1DI0	C04-N05	-2.16	1.32	1.38
3	B	402	NAD	C5B-C4B	2.15	1.58	1.51
3	A	402	NAD	C4N-C3N	-2.06	1.36	1.39
2	B	401	A1DI0	O33-C32	-2.03	1.37	1.43
3	A	402	NAD	PN-O3	2.02	1.61	1.59
2	B	401	A1DI0	C04-C03	2.00	1.39	1.35

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	NAD	N3A-C2A-N1A	-5.60	120.10	128.58
3	A	402	NAD	N3A-C2A-N1A	-5.58	120.13	128.58
3	A	402	NAD	C5A-C4A-N3A	-5.16	119.61	126.72
2	A	401	A1DI0	C02-N36-C34	-4.43	121.11	126.61
3	B	402	NAD	C5A-C4A-N3A	-4.37	120.69	126.72
3	A	402	NAD	N9A-C8A-N7A	-4.28	107.86	113.94
3	A	402	NAD	C6N-N1N-C2N	-4.27	118.25	121.88
2	A	401	A1DI0	N36-C34-N05	4.24	120.41	114.89
2	B	401	A1DI0	N36-C34-N05	4.21	120.37	114.89
3	B	402	NAD	N9A-C8A-N7A	-4.05	108.20	113.94
2	A	401	A1DI0	O20-C19-O18	-4.04	106.08	111.36
2	B	401	A1DI0	C02-N36-C34	-3.64	122.09	126.61
3	A	402	NAD	C2A-N3A-C4A	3.54	120.48	111.83
3	A	402	NAD	C1B-N9A-C8A	-3.37	119.62	127.09
3	B	402	NAD	C2A-N3A-C4A	3.32	119.94	111.83
3	B	402	NAD	C6N-N1N-C2N	-3.28	119.09	121.88
2	A	401	A1DI0	C03-C02-N36	3.23	119.32	114.80
3	A	402	NAD	N3A-C4A-N9A	3.07	132.40	127.17
3	B	402	NAD	O3D-C3D-C4D	-3.07	102.26	111.08
2	A	401	A1DI0	C26-C24-C21	3.06	115.78	110.23
2	A	401	A1DI0	O01-C02-C03	-2.91	120.15	125.16
2	B	401	A1DI0	C03-C02-N36	2.86	118.80	114.80
3	A	402	NAD	C5A-N7A-C8A	2.76	107.79	103.45
3	A	402	NAD	O3D-C3D-C4D	-2.63	103.52	111.08
3	B	402	NAD	C1B-N9A-C8A	-2.52	121.50	127.09
3	A	402	NAD	C4B-O4B-C1B	-2.52	103.90	109.47
3	A	402	NAD	N6A-C6A-N1A	-2.46	112.89	118.38
3	A	402	NAD	O4B-C1B-C2B	-2.43	101.42	106.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	NAD	N3A-C4A-N9A	2.40	131.25	127.17
3	A	402	NAD	C4A-N9A-C1B	2.37	132.18	126.63
3	B	402	NAD	C5A-N7A-C8A	2.37	107.17	103.45
3	B	402	NAD	N6A-C6A-N1A	-2.36	113.12	118.38
3	A	402	NAD	C4A-N9A-C8A	2.35	108.20	105.74
3	A	402	NAD	O2A-PA-O3	2.27	113.42	107.27
2	A	401	A1DI0	O20-C21-C24	2.26	113.78	109.70
3	B	402	NAD	C4A-N9A-C8A	2.24	108.08	105.74
2	A	401	A1DI0	C22-C21-C24	-2.19	107.64	113.02
3	A	402	NAD	C4A-C5A-N7A	-2.13	108.14	110.58
2	B	401	A1DI0	O20-C19-O18	-2.12	108.59	111.36
2	A	401	A1DI0	O27-C26-C28	-2.09	105.44	110.38
2	B	401	A1DI0	O35-C34-N05	-2.07	120.10	122.80
2	B	401	A1DI0	O01-C02-C03	-2.07	121.60	125.16
3	B	402	NAD	C5A-C4A-N9A	2.05	108.05	105.81
2	A	401	A1DI0	C03-C04-N05	-2.05	118.51	121.84
3	B	402	NAD	O5B-C5B-C4B	-2.03	102.08	108.99
3	B	402	NAD	O7N-C7N-N7N	-2.00	119.72	122.62
3	A	402	NAD	C5A-C4A-N9A	2.00	107.99	105.81

There are no chirality outliers.

All (16) torsion outliers are listed below:

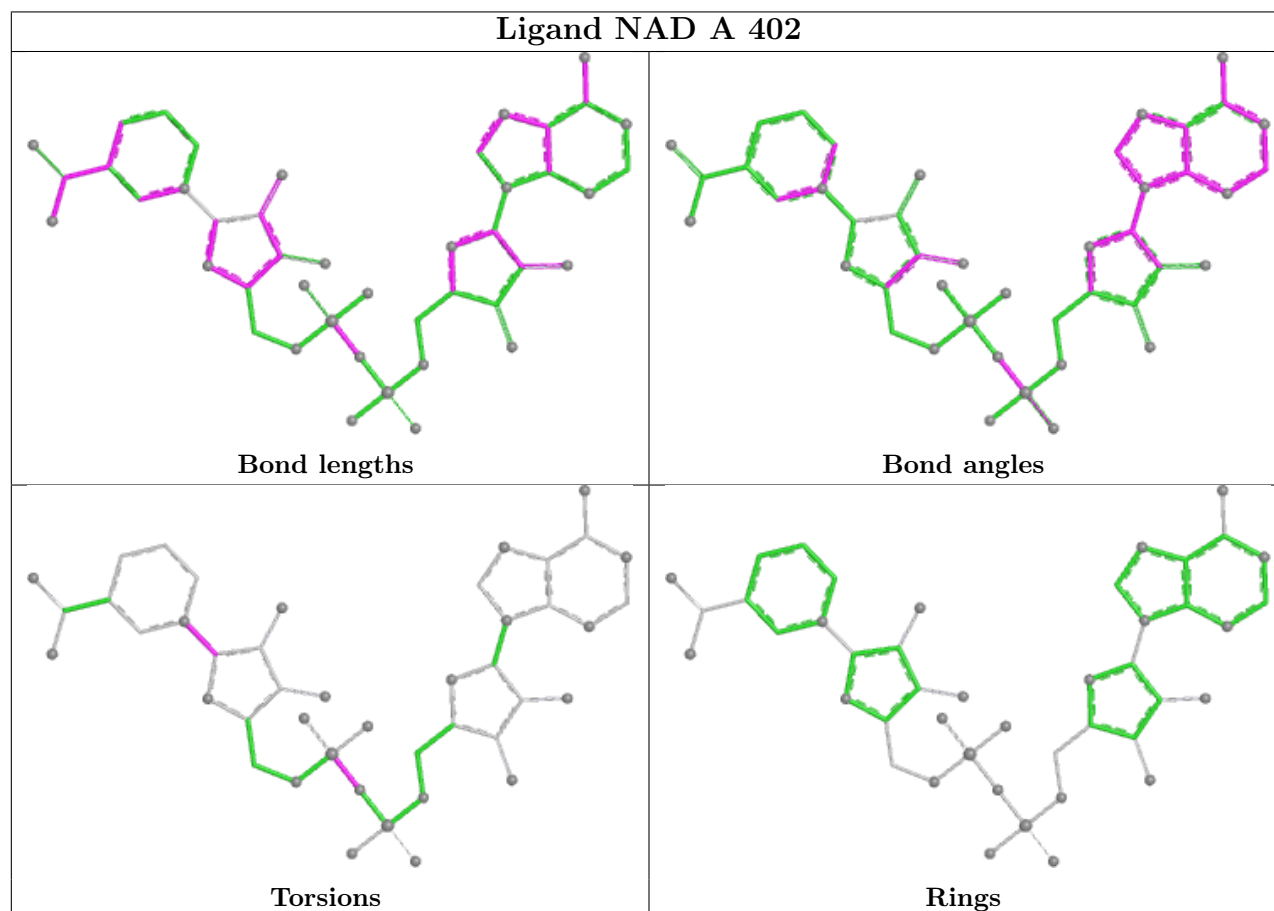
Mol	Chain	Res	Type	Atoms
2	B	401	A1DI0	C09-O10-P11-O13
3	A	402	NAD	O4D-C1D-N1N-C2N
3	A	402	NAD	O4D-C1D-N1N-C6N
3	B	402	NAD	O4D-C1D-N1N-C2N
3	B	402	NAD	O4D-C1D-N1N-C6N
3	B	402	NAD	C2D-C1D-N1N-C6N
2	B	401	A1DI0	O07-C08-C09-O10
2	B	401	A1DI0	C30-C08-C09-O10
2	A	401	A1DI0	P15-O14-P11-O10
2	B	401	A1DI0	P15-O14-P11-O10
2	B	401	A1DI0	C09-O10-P11-O12
2	B	401	A1DI0	C09-O10-P11-O14
3	A	402	NAD	PA-O3-PN-O2N
3	A	402	NAD	C2D-C1D-N1N-C6N
3	A	402	NAD	PA-O3-PN-O1N
3	B	402	NAD	O4B-C4B-C5B-O5B

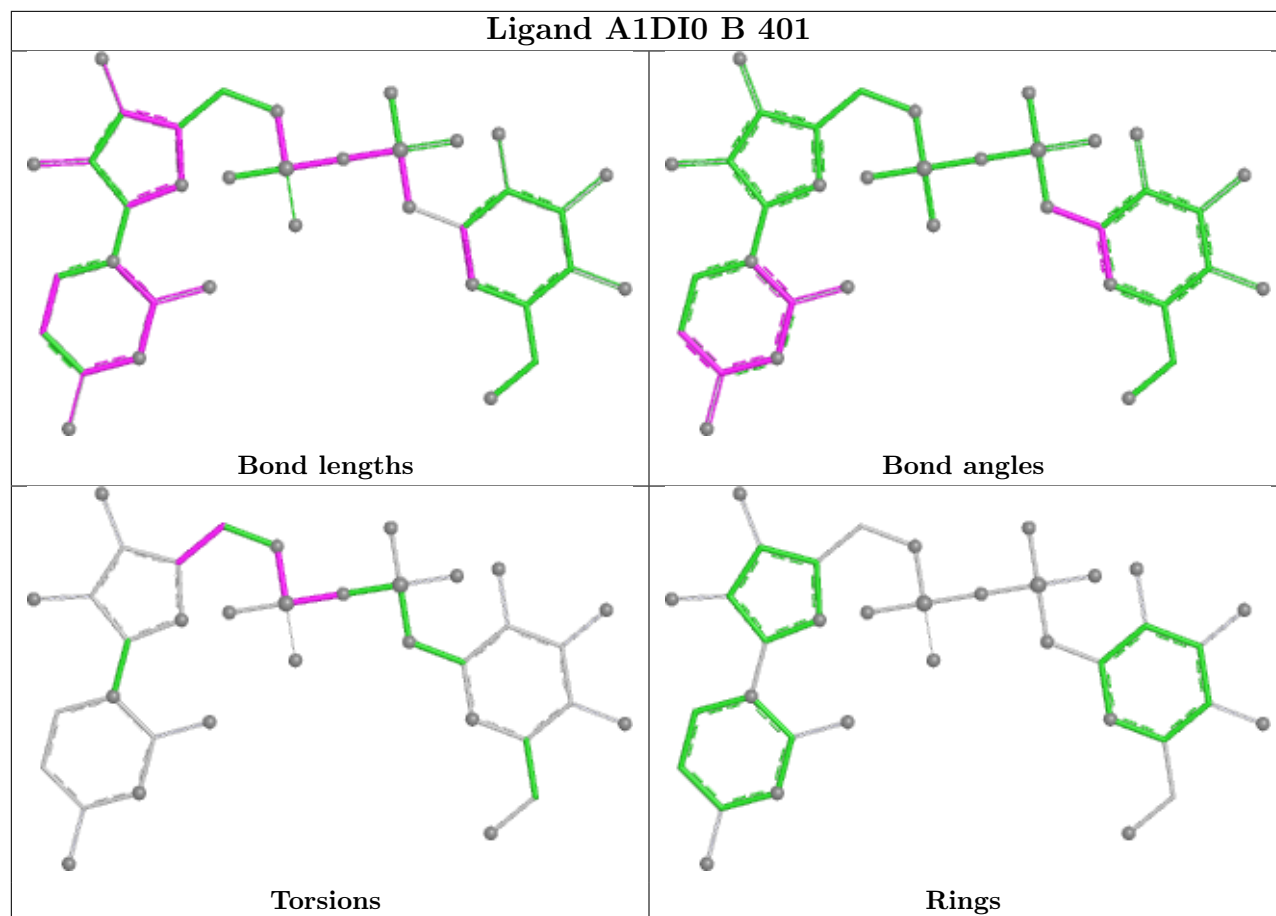
There are no ring outliers.

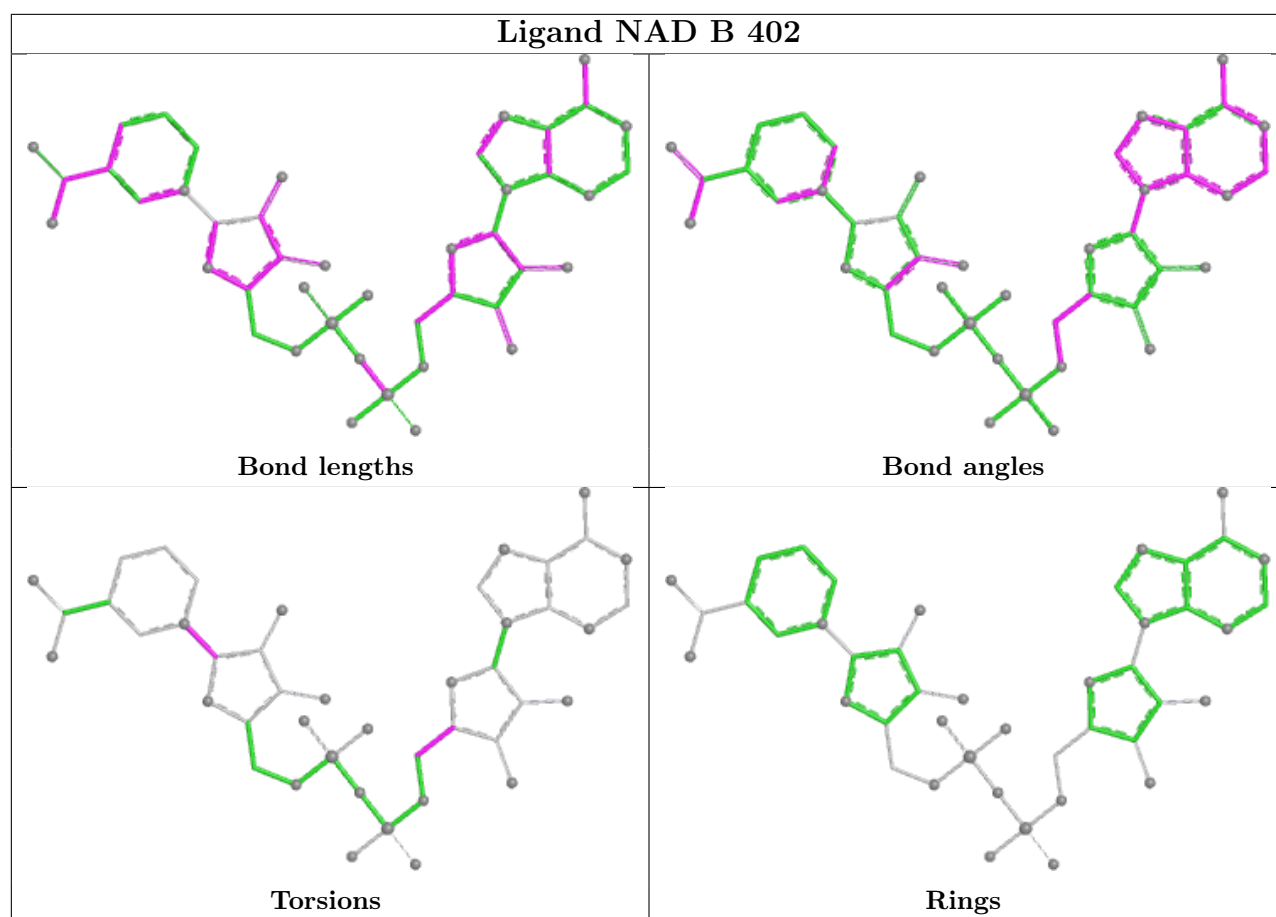
1 monomer is involved in 1 short contact:

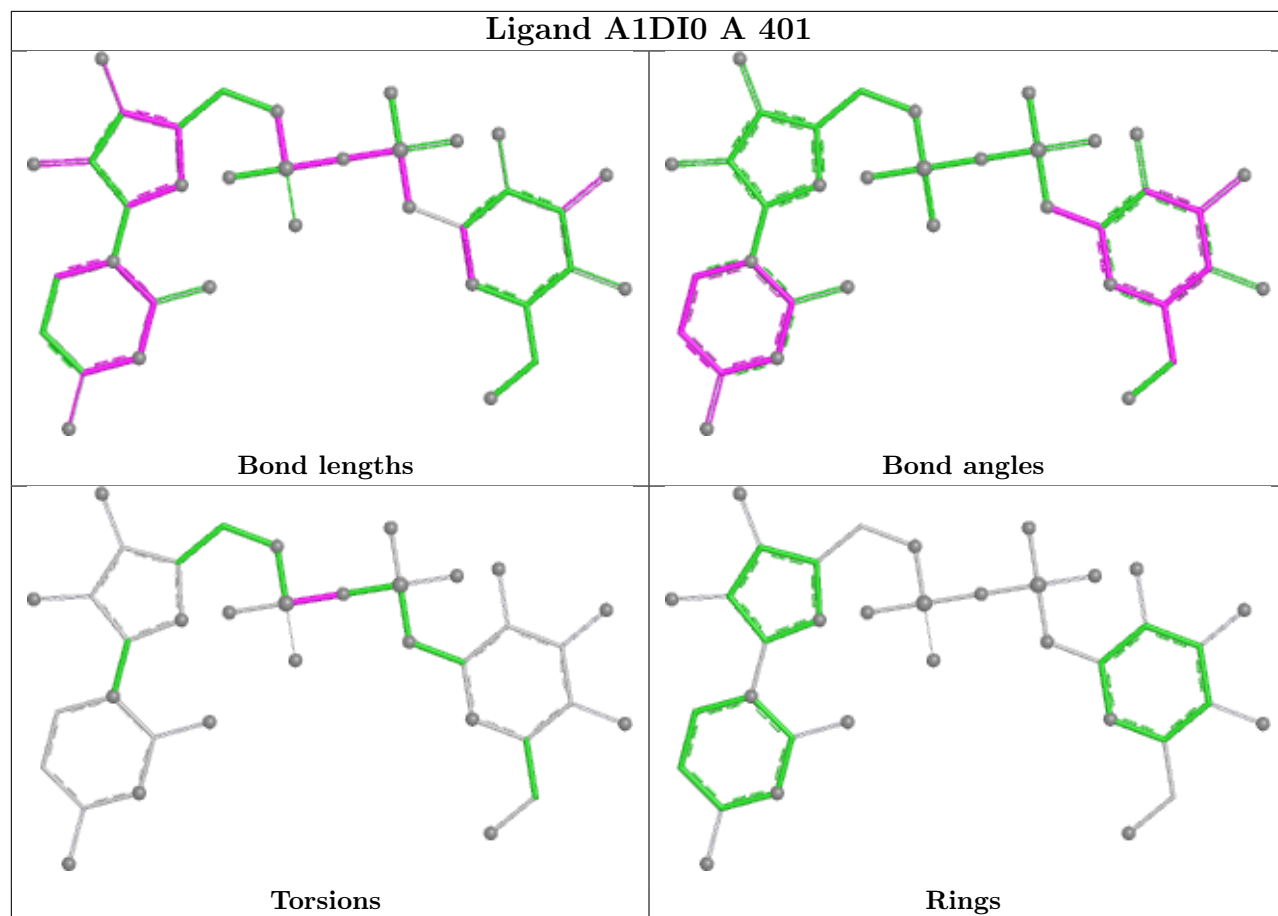
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	A1DI0	1	0

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 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	339/348 (97%)	0.39	10 (2%) 53 58	9, 18, 33, 51	0
1	B	339/348 (97%)	0.35	7 (2%) 63 70	10, 17, 32, 51	0
All	All	678/696 (97%)	0.37	17 (2%) 58 64	9, 18, 32, 51	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	ASP	3.7
1	A	65	ASP	3.1
1	A	6	GLY	2.9
1	B	106	HIS	2.8
1	A	314	ASP	2.8
1	A	106	HIS	2.7
1	A	130	ILE	2.4
1	B	64	PRO	2.4
1	A	64	PRO	2.4
1	A	270	THR	2.3
1	B	314	ASP	2.3
1	B	339	GLU	2.2
1	B	130	ILE	2.2
1	A	268	GLU	2.2
1	A	132	ASP	2.2
1	A	165	ASP	2.0
1	B	231	ASP	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](#)

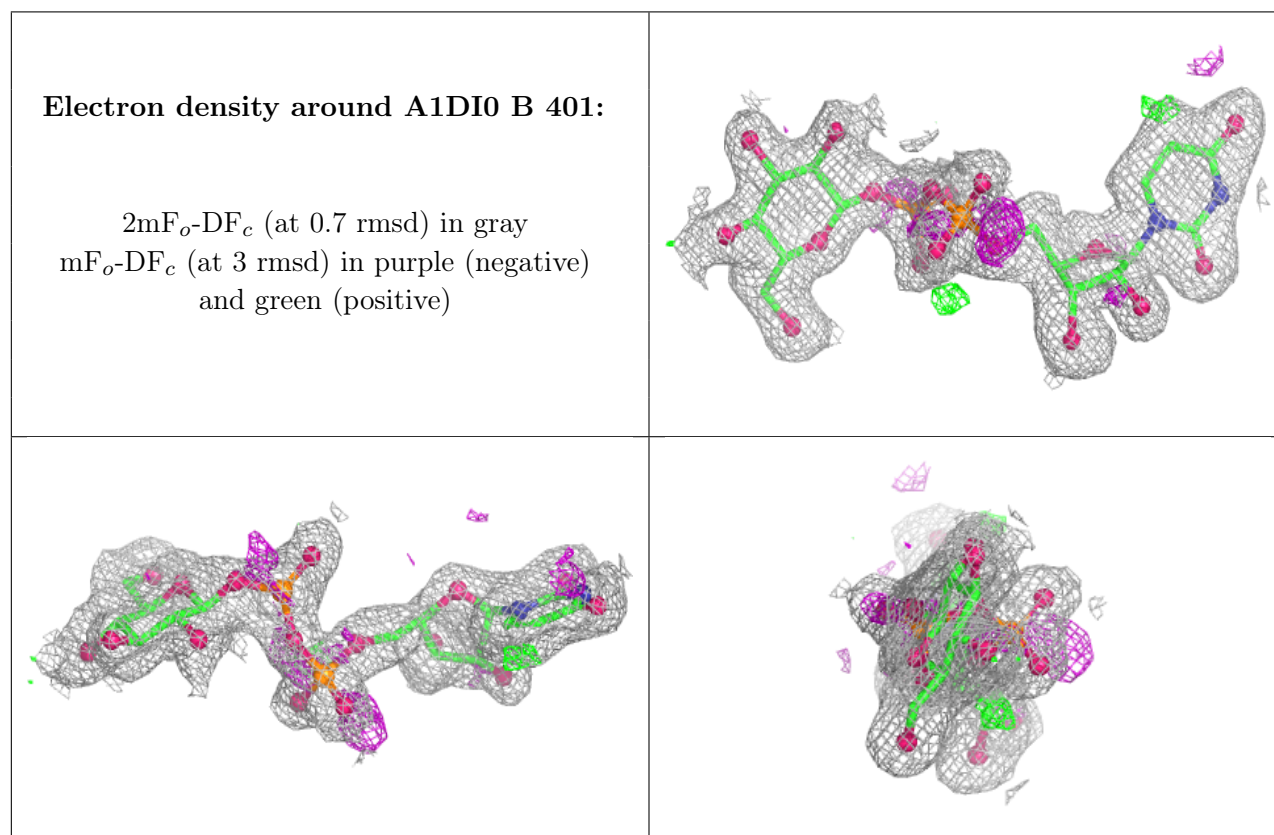
There are no oligosaccharides in this entry.

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.

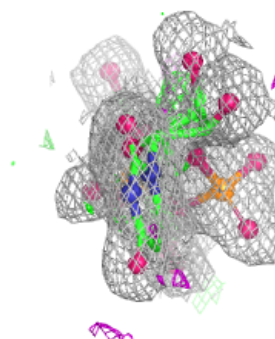
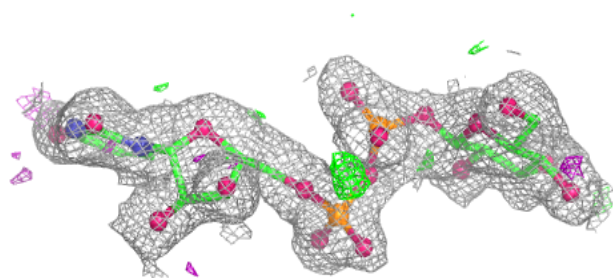
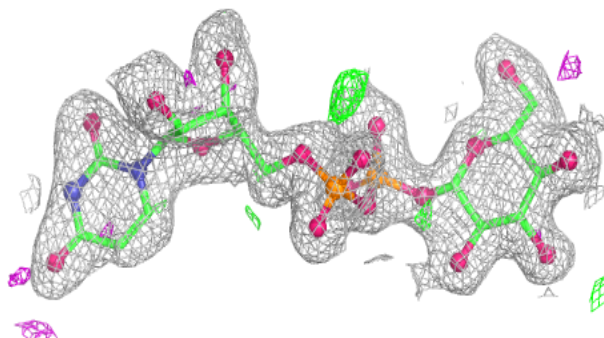
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	A1DI0	B	401	36/36	0.92	0.10	13,20,24,25	0
2	A1DI0	A	401	36/36	0.94	0.08	13,21,27,28	0
3	NAD	A	402	44/44	0.95	0.07	6,12,14,15	0
3	NAD	B	402	44/44	0.96	0.06	6,12,15,18	0

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.

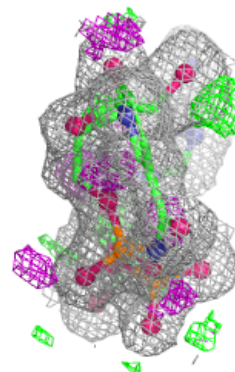
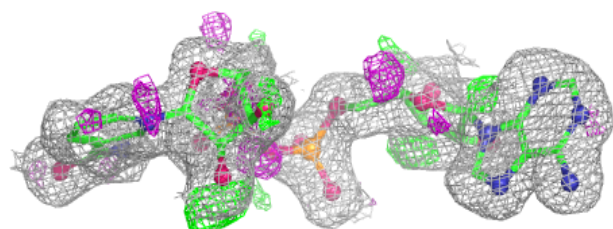
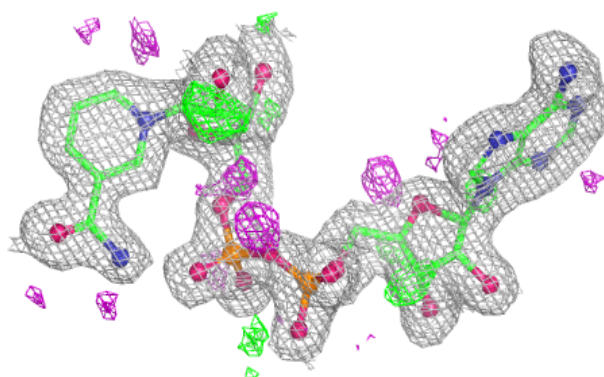


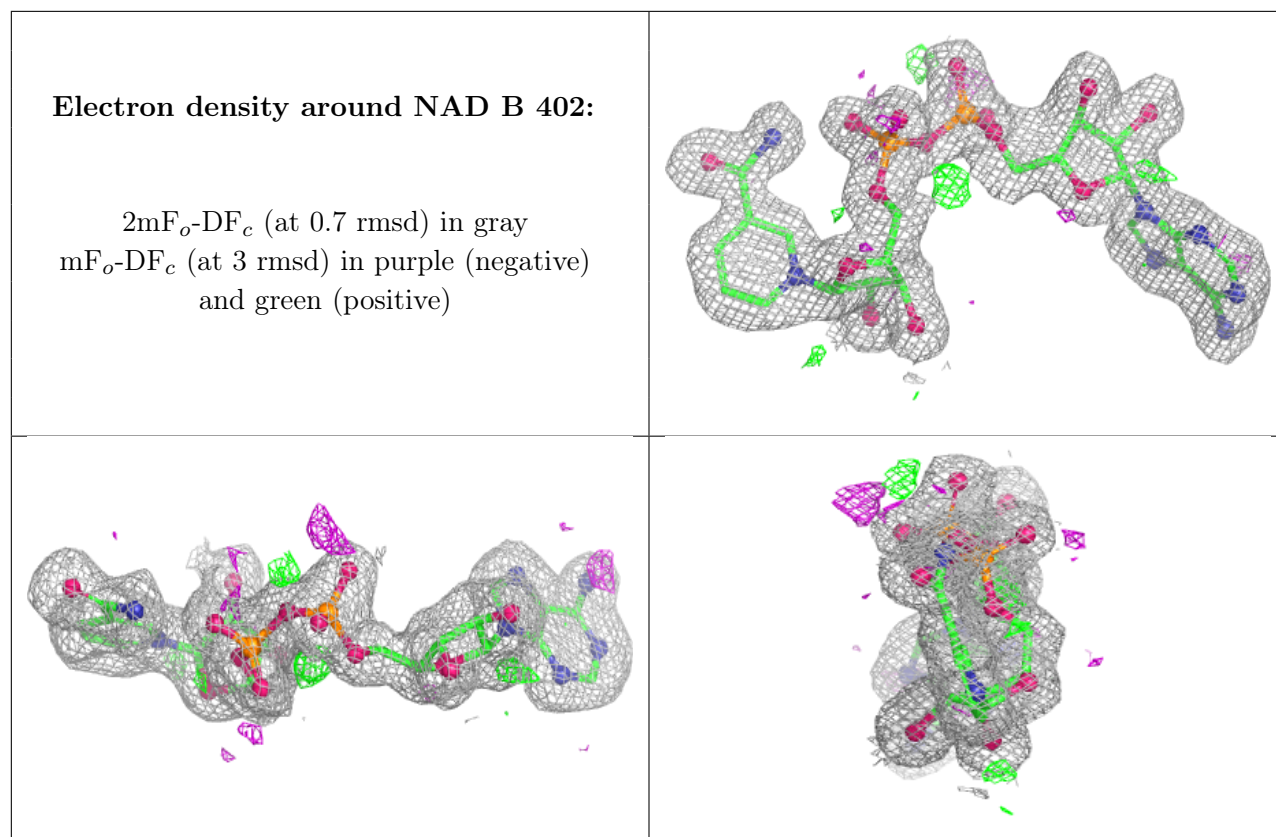
Electron density around A1DI0 A 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around NAD A 402:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers ⓘ

There are no such residues in this entry.