



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 06:23 am BST

PDB ID : 4QX6
Title : CRYSTAL STRUCTURE OF GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE FROM STREPTOCOCCUS AGALACTIAE NEM316 at 2.46 ANGSTROM RESOLUTION
Authors : Ayres, C.A.; Schormann, N.; Banerjee, S.; Chattopadhyay, D.
Deposited on : 2014-07-18
Resolution : 2.46 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

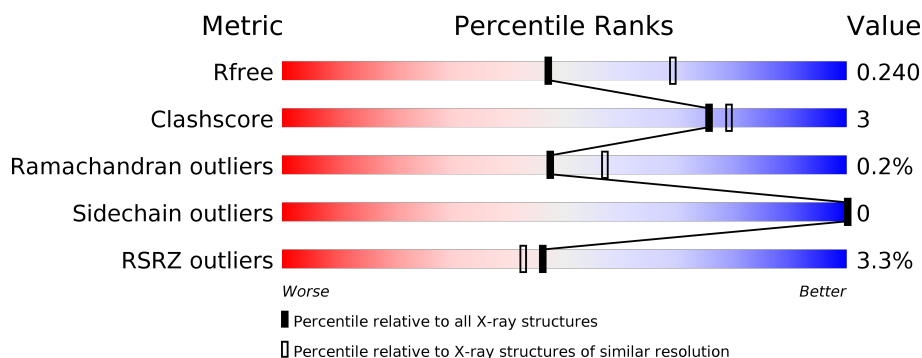
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	356	<div> <div>7%</div> <div>88% 6% 7%</div> </div>
1	B	356	<div> <div>%</div> <div>89% 5% 6%</div> </div>
1	C	356	<div> <div>4%</div> <div>89% 5% 6%</div> </div>
1	D	356	<div> <div>%</div> <div>91% 5% .</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	C	402	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10428 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 Glyceraldehyde 3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	0	0	0
			2498	1561	433	496	8			
1	B	335	Total	C	N	O	S	0	0	0
			2522	1576	438	500	8			
1	C	334	Total	C	N	O	S	0	0	0
			2513	1570	436	499	8			
1	D	344	Total	C	N	O	S	0	0	0
			2587	1616	452	510	9			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q8E3E8
A	-18	GLY	-	EXPRESSION TAG	UNP Q8E3E8
A	-17	SER	-	EXPRESSION TAG	UNP Q8E3E8
A	-16	SER	-	EXPRESSION TAG	UNP Q8E3E8
A	-15	HIS	-	EXPRESSION TAG	UNP Q8E3E8
A	-14	HIS	-	EXPRESSION TAG	UNP Q8E3E8
A	-13	HIS	-	EXPRESSION TAG	UNP Q8E3E8
A	-12	HIS	-	EXPRESSION TAG	UNP Q8E3E8
A	-11	HIS	-	EXPRESSION TAG	UNP Q8E3E8
A	-10	HIS	-	EXPRESSION TAG	UNP Q8E3E8
A	-9	SER	-	EXPRESSION TAG	UNP Q8E3E8
A	-8	SER	-	EXPRESSION TAG	UNP Q8E3E8
A	-7	GLY	-	EXPRESSION TAG	UNP Q8E3E8
A	-6	LEU	-	EXPRESSION TAG	UNP Q8E3E8
A	-5	VAL	-	EXPRESSION TAG	UNP Q8E3E8
A	-4	PRO	-	EXPRESSION TAG	UNP Q8E3E8
A	-3	ARG	-	EXPRESSION TAG	UNP Q8E3E8
A	-2	GLY	-	EXPRESSION TAG	UNP Q8E3E8
A	-1	SER	-	EXPRESSION TAG	UNP Q8E3E8
A	0	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-19	MET	-	EXPRESSION TAG	UNP Q8E3E8

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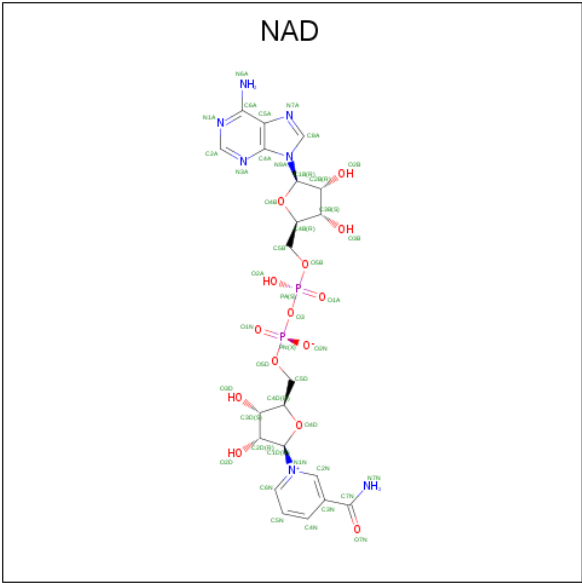
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	EXPRESSION TAG	UNP Q8E3E8
B	-17	SER	-	EXPRESSION TAG	UNP Q8E3E8
B	-16	SER	-	EXPRESSION TAG	UNP Q8E3E8
B	-15	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-14	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-13	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-12	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-11	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-10	HIS	-	EXPRESSION TAG	UNP Q8E3E8
B	-9	SER	-	EXPRESSION TAG	UNP Q8E3E8
B	-8	SER	-	EXPRESSION TAG	UNP Q8E3E8
B	-7	GLY	-	EXPRESSION TAG	UNP Q8E3E8
B	-6	LEU	-	EXPRESSION TAG	UNP Q8E3E8
B	-5	VAL	-	EXPRESSION TAG	UNP Q8E3E8
B	-4	PRO	-	EXPRESSION TAG	UNP Q8E3E8
B	-3	ARG	-	EXPRESSION TAG	UNP Q8E3E8
B	-2	GLY	-	EXPRESSION TAG	UNP Q8E3E8
B	-1	SER	-	EXPRESSION TAG	UNP Q8E3E8
B	0	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-19	MET	-	EXPRESSION TAG	UNP Q8E3E8
C	-18	GLY	-	EXPRESSION TAG	UNP Q8E3E8
C	-17	SER	-	EXPRESSION TAG	UNP Q8E3E8
C	-16	SER	-	EXPRESSION TAG	UNP Q8E3E8
C	-15	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-14	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-13	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-12	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-11	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-10	HIS	-	EXPRESSION TAG	UNP Q8E3E8
C	-9	SER	-	EXPRESSION TAG	UNP Q8E3E8
C	-8	SER	-	EXPRESSION TAG	UNP Q8E3E8
C	-7	GLY	-	EXPRESSION TAG	UNP Q8E3E8
C	-6	LEU	-	EXPRESSION TAG	UNP Q8E3E8
C	-5	VAL	-	EXPRESSION TAG	UNP Q8E3E8
C	-4	PRO	-	EXPRESSION TAG	UNP Q8E3E8
C	-3	ARG	-	EXPRESSION TAG	UNP Q8E3E8
C	-2	GLY	-	EXPRESSION TAG	UNP Q8E3E8
C	-1	SER	-	EXPRESSION TAG	UNP Q8E3E8
C	0	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-19	MET	-	EXPRESSION TAG	UNP Q8E3E8
D	-18	GLY	-	EXPRESSION TAG	UNP Q8E3E8
D	-17	SER	-	EXPRESSION TAG	UNP Q8E3E8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	EXPRESSION TAG	UNP Q8E3E8
D	-15	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-14	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-13	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-12	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-11	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-10	HIS	-	EXPRESSION TAG	UNP Q8E3E8
D	-9	SER	-	EXPRESSION TAG	UNP Q8E3E8
D	-8	SER	-	EXPRESSION TAG	UNP Q8E3E8
D	-7	GLY	-	EXPRESSION TAG	UNP Q8E3E8
D	-6	LEU	-	EXPRESSION TAG	UNP Q8E3E8
D	-5	VAL	-	EXPRESSION TAG	UNP Q8E3E8
D	-4	PRO	-	EXPRESSION TAG	UNP Q8E3E8
D	-3	ARG	-	EXPRESSION TAG	UNP Q8E3E8
D	-2	GLY	-	EXPRESSION TAG	UNP Q8E3E8
D	-1	SER	-	EXPRESSION TAG	UNP Q8E3E8
D	0	HIS	-	EXPRESSION TAG	UNP Q8E3E8

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

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

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		

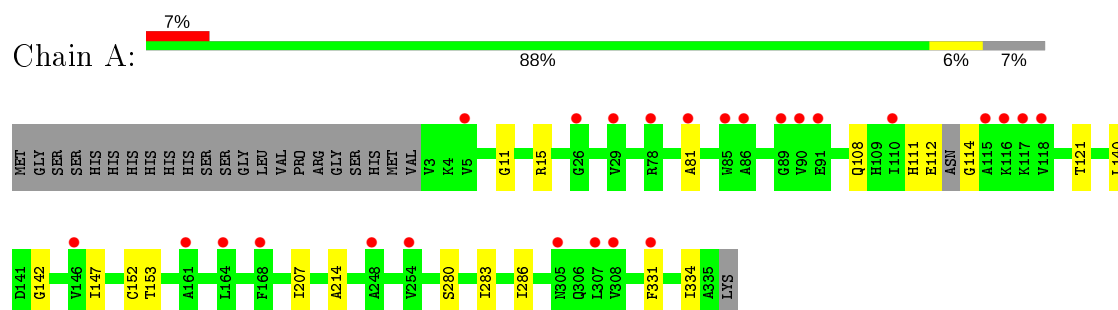
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	32	Total	O	0	0
			32	32		
4	C	21	Total	O	0	0
			21	21		
4	D	39	Total	O	0	0
			39	39		

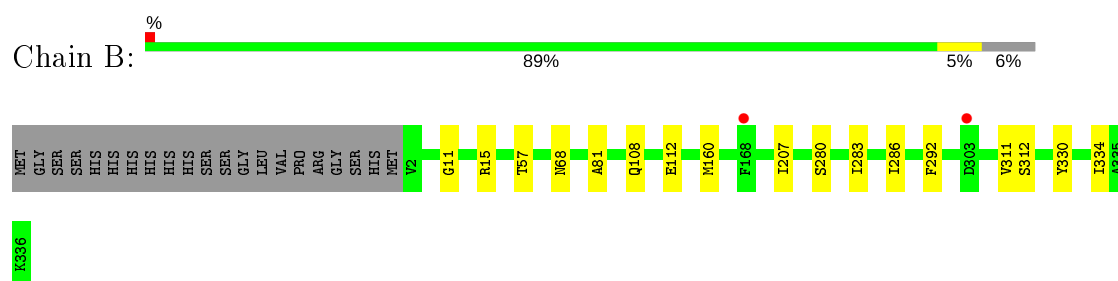
3 Residue-property plots [i](#)

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 ($RSRZ > 2$). 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.

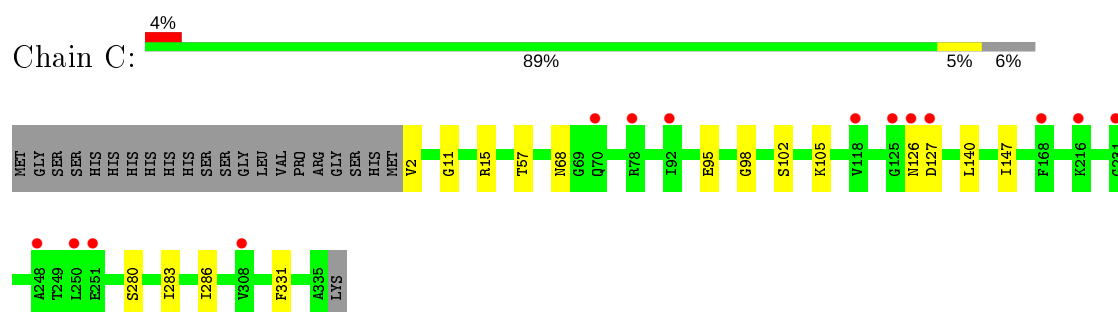
- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase



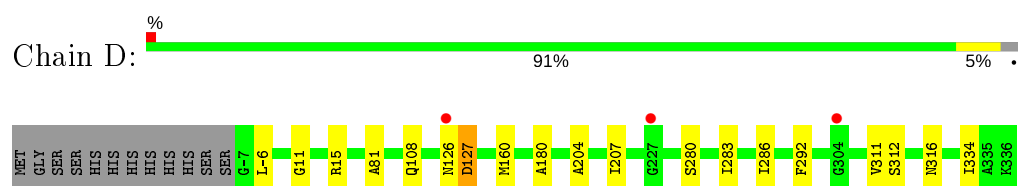
- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase



- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase



- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.77Å 107.86Å 90.93Å 90.00° 106.57° 90.00°	Depositor
Resolution (Å)	64.96 – 2.46 64.96 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.3 (64.96-2.46) 99.3 (64.96-2.46)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.204 , 0.238 0.208 , 0.240	Depositor DCC
R_{free} test set	2279 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	48.4	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10428	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, 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	0.33	0/2535	0.58	0/3439
1	B	0.34	0/2560	0.57	0/3474
1	C	0.34	0/2551	0.57	0/3463
1	D	0.35	0/2627	0.56	0/3564
All	All	0.34	0/10273	0.57	0/13940

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	A	2498	0	2476	16	0
1	B	2522	0	2505	15	0
1	C	2513	0	2492	17	0
1	D	2587	0	2575	18	0
2	A	44	0	26	2	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	D	44	0	26	1	0
3	C	8	0	12	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	32	0	0	0	0
4	B	32	0	0	0	0
4	C	21	0	0	0	0
4	D	39	0	0	1	0
All	All	10428	0	10164	63	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 3.

All (63) 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:D:126:ASN:O	1:D:127:ASP:HB2	1.90	0.71
1:C:98:GLY:CA	3:C:402:EDO:H12	2.23	0.68
1:A:112:GLU:HA	1:A:114:GLY:N	2.10	0.66
1:D:180:ALA:HB3	1:D:316:ASN:HD21	1.65	0.61
1:C:2:VAL:HG12	1:C:2:VAL:O	2.01	0.60
1:D:160:MET:HE1	1:D:311:VAL:N	2.18	0.59
1:B:160:MET:HE1	1:B:311:VAL:N	2.19	0.58
1:C:126:ASN:O	1:C:127:ASP:HB2	2.04	0.57
1:B:160:MET:HE2	1:B:292:PHE:CD1	2.42	0.55
1:D:160:MET:HE2	1:D:292:PHE:CD1	2.42	0.54
1:B:280:SER:O	1:B:283:ILE:HG13	2.07	0.54
1:B:160:MET:HE2	1:B:292:PHE:HD1	1.72	0.54
1:A:280:SER:O	1:A:283:ILE:HG13	2.08	0.54
1:A:147:ILE:HD12	1:A:331:PHE:CE1	2.43	0.53
1:C:280:SER:O	1:C:283:ILE:HG13	2.08	0.53
1:D:280:SER:O	1:D:283:ILE:HG13	2.08	0.53
1:D:160:MET:HE2	1:D:292:PHE:HD1	1.74	0.52
1:A:153:THR:HG23	1:A:214:ALA:CB	2.40	0.52
1:C:147:ILE:HD12	1:C:331:PHE:CE1	2.44	0.52
1:D:-6:LEU:HD13	1:D:334:ILE:HD11	1.92	0.52
1:C:95:GLU:OE2	3:C:402:EDO:O1	2.27	0.52
1:B:330:TYR:O	1:B:334:ILE:HG12	2.11	0.51
1:B:160:MET:CE	1:B:292:PHE:HD1	2.24	0.51
1:D:160:MET:CE	1:D:292:PHE:HD1	2.24	0.50
1:D:180:ALA:HB3	1:D:316:ASN:ND2	2.28	0.49
1:D:160:MET:HE3	1:D:312:SER:HB2	1.95	0.49
1:B:330:TYR:CE2	1:B:334:ILE:HD13	2.48	0.48
1:C:98:GLY:H	3:C:402:EDO:C1	2.27	0.48
1:B:81:ALA:HB2	1:B:108:GLN:HG2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:ILE:HD12	1:A:331:PHE:CZ	2.50	0.47
1:B:160:MET:HE3	1:B:312:SER:HB2	1.95	0.47
1:A:153:THR:HG23	1:A:214:ALA:HB2	1.97	0.47
1:D:160:MET:CE	1:D:292:PHE:CD1	2.98	0.47
1:B:160:MET:CE	1:B:292:PHE:CD1	2.98	0.47
1:D:81:ALA:HB2	1:D:108:GLN:HG2	1.96	0.46
1:A:140:LEU:CD2	1:A:147:ILE:HD13	2.46	0.46
1:C:140:LEU:CD2	1:C:147:ILE:HD13	2.46	0.46
1:A:112:GLU:CA	1:A:114:GLY:N	2.78	0.46
1:C:147:ILE:HD12	1:C:331:PHE:CZ	2.51	0.46
1:C:98:GLY:HA2	3:C:402:EDO:H12	1.96	0.46
1:C:126:ASN:O	1:C:127:ASP:CB	2.63	0.45
1:C:102:SER:OG	1:C:105:LYS:HE3	2.16	0.45
1:D:11:GLY:HA3	2:D:401:NAD:O5B	2.17	0.45
1:A:152:CYS:HB3	2:A:401:NAD:H4N	1.98	0.44
1:D:283:ILE:O	1:D:286:ILE:HG12	2.18	0.43
1:A:81:ALA:HB2	1:A:108:GLN:HG2	1.99	0.43
1:A:283:ILE:O	1:A:286:ILE:HG12	2.19	0.43
1:C:98:GLY:N	3:C:402:EDO:H12	2.34	0.42
1:C:283:ILE:O	1:C:286:ILE:HG12	2.19	0.42
1:B:11:GLY:O	1:B:15:ARG:HG3	2.20	0.42
1:B:283:ILE:O	1:B:286:ILE:HG12	2.19	0.42
1:B:207:ILE:HB	1:D:280:SER:HB3	2.02	0.42
1:B:57:THR:HG22	1:B:68:ASN:HA	2.01	0.42
1:D:11:GLY:O	1:D:15:ARG:HG3	2.20	0.42
1:A:140:LEU:HD22	1:A:147:ILE:HD13	2.03	0.41
1:B:280:SER:HB3	1:D:207:ILE:HB	2.03	0.41
1:C:11:GLY:O	1:C:15:ARG:HG3	2.21	0.41
1:C:57:THR:HG22	1:C:68:ASN:HA	2.02	0.41
1:A:11:GLY:O	1:A:15:ARG:HG3	2.20	0.41
1:A:121:THR:O	2:A:401:NAD:H1D	2.20	0.41
1:D:204:ALA:HB3	4:D:503:HOH:O	2.20	0.41
1:A:142:GLY:HA3	1:A:334:ILE:HG22	2.03	0.40
1:A:207:ILE:HB	1:C:280:SER:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	328/356 (92%)	309 (94%)	18 (6%)	1 (0%)	41	49
1	B	333/356 (94%)	312 (94%)	20 (6%)	1 (0%)	41	49
1	C	332/356 (93%)	313 (94%)	19 (6%)	0	100	100
1	D	342/356 (96%)	321 (94%)	20 (6%)	1 (0%)	41	49
All	All	1335/1424 (94%)	1255 (94%)	77 (6%)	3 (0%)	47	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	127	ASP
1	A	111	HIS
1	B	112	GLU

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	266/287 (93%)	266 (100%)	0	100	100
1	B	269/287 (94%)	269 (100%)	0	100	100
1	C	268/287 (93%)	268 (100%)	0	100	100
1	D	276/287 (96%)	276 (100%)	0	100	100
All	All	1079/1148 (94%)	1079 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$
2	NAD	A	401	-	42,48,48	0.78	1 (2%)	50,73,73	1.12	4 (8%)
3	EDO	C	403	-	3,3,3	0.46	0	2,2,2	0.34	0
2	NAD	D	401	-	42,48,48	0.90	3 (7%)	50,73,73	1.24	5 (10%)
3	EDO	C	402	-	3,3,3	0.45	0	2,2,2	0.34	0
2	NAD	B	401	-	42,48,48	0.79	2 (4%)	50,73,73	1.13	4 (8%)
2	NAD	C	401	-	42,48,48	0.83	2 (4%)	50,73,73	1.24	4 (8%)

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
2	NAD	A	401	-	-	5/26/62/62	0/5/5/5
3	EDO	C	403	-	-	1/1/1/1	-
2	NAD	D	401	-	-	6/26/62/62	0/5/5/5
3	EDO	C	402	-	-	1/1/1/1	-
2	NAD	B	401	-	-	7/26/62/62	0/5/5/5
2	NAD	C	401	-	-	3/26/62/62	0/5/5/5

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	NAD	O4D-C1D	2.70	1.44	1.41
2	C	401	NAD	C5A-C4A	2.22	1.46	1.40
2	D	401	NAD	C5A-C4A	2.19	1.46	1.40
2	A	401	NAD	C5A-C4A	2.19	1.46	1.40
2	B	401	NAD	C5A-C4A	2.16	1.46	1.40
2	B	401	NAD	O4D-C1D	2.10	1.44	1.41
2	C	401	NAD	C2A-N3A	2.05	1.35	1.32
2	D	401	NAD	C2A-N3A	2.02	1.35	1.32

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	NAD	N3A-C2A-N1A	-4.02	122.40	128.68
2	B	401	NAD	N3A-C2A-N1A	-3.90	122.58	128.68
2	A	401	NAD	N3A-C2A-N1A	-3.87	122.63	128.68
2	C	401	NAD	N3A-C2A-N1A	-3.65	122.98	128.68
2	C	401	NAD	C3N-C7N-N7N	3.49	121.94	117.75
2	C	401	NAD	PN-O3-PA	-3.25	121.69	132.83
2	C	401	NAD	C4A-C5A-N7A	-2.72	106.57	109.40
2	D	401	NAD	C4A-C5A-N7A	-2.54	106.76	109.40
2	B	401	NAD	C4A-C5A-N7A	-2.51	106.78	109.40
2	D	401	NAD	PN-O3-PA	-2.49	124.30	132.83
2	B	401	NAD	C2A-N1A-C6A	2.36	122.80	118.75
2	A	401	NAD	C4A-C5A-N7A	-2.25	107.06	109.40
2	A	401	NAD	C2A-N1A-C6A	2.23	122.56	118.75
2	B	401	NAD	PN-O3-PA	-2.21	125.23	132.83
2	D	401	NAD	O7N-C7N-C3N	-2.14	117.07	119.63
2	A	401	NAD	O2A-PA-O1A	2.12	122.73	112.24
2	D	401	NAD	C1B-N9A-C4A	-2.09	122.97	126.64

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAD	C5B-O5B-PA-O1A
2	A	401	NAD	O4D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	C2D-C1D-N1N-C6N
2	B	401	NAD	O4D-C1D-N1N-C2N
2	B	401	NAD	O4D-C1D-N1N-C6N
2	C	401	NAD	O4B-C4B-C5B-O5B
2	C	401	NAD	O4D-C1D-N1N-C2N
2	A	401	NAD	O4B-C4B-C5B-O5B
2	B	401	NAD	O4B-C4B-C5B-O5B
2	C	401	NAD	C3B-C4B-C5B-O5B
2	A	401	NAD	C3B-C4B-C5B-O5B
3	C	403	EDO	O1-C1-C2-O2
3	C	402	EDO	O1-C1-C2-O2
2	B	401	NAD	C3B-C4B-C5B-O5B
2	A	401	NAD	C5B-O5B-PA-O3
2	D	401	NAD	O4B-C4B-C5B-O5B
2	B	401	NAD	C2N-C3N-C7N-N7N
2	D	401	NAD	C2D-C1D-N1N-C2N
2	B	401	NAD	C2D-C1D-N1N-C6N
2	D	401	NAD	PN-O3-PA-O1A
2	B	401	NAD	C2N-C3N-C7N-O7N

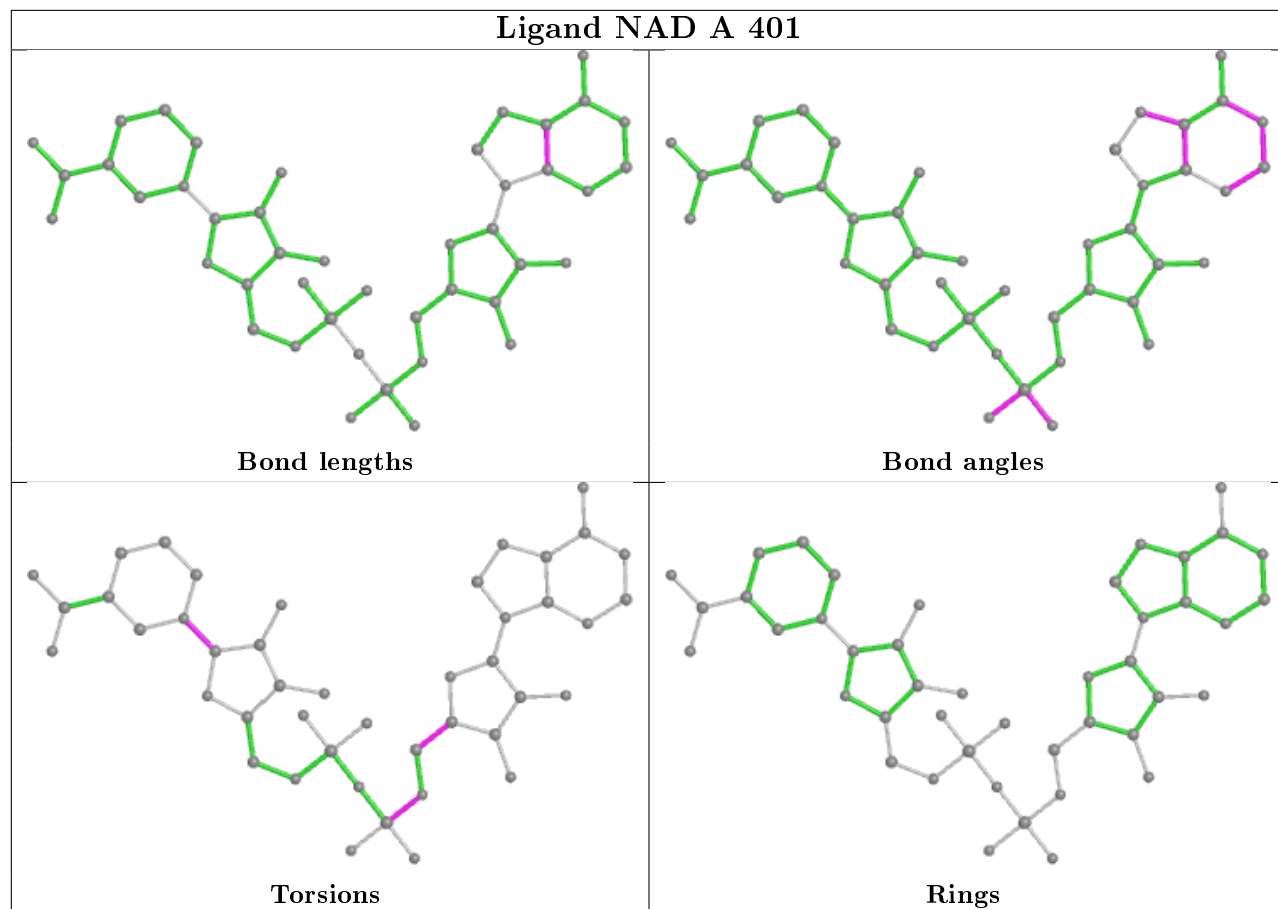
There are no ring outliers.

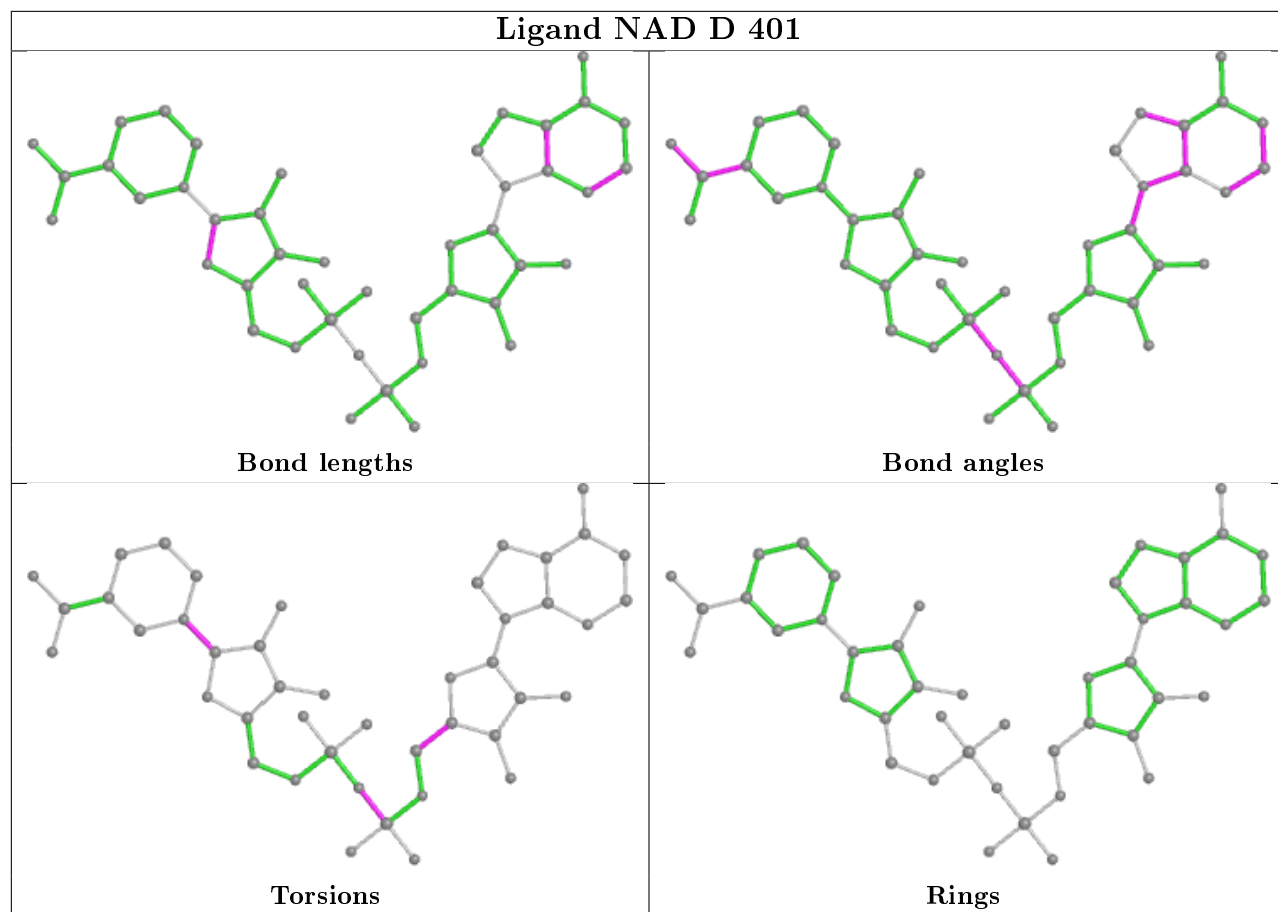
3 monomers are involved in 8 short contacts:

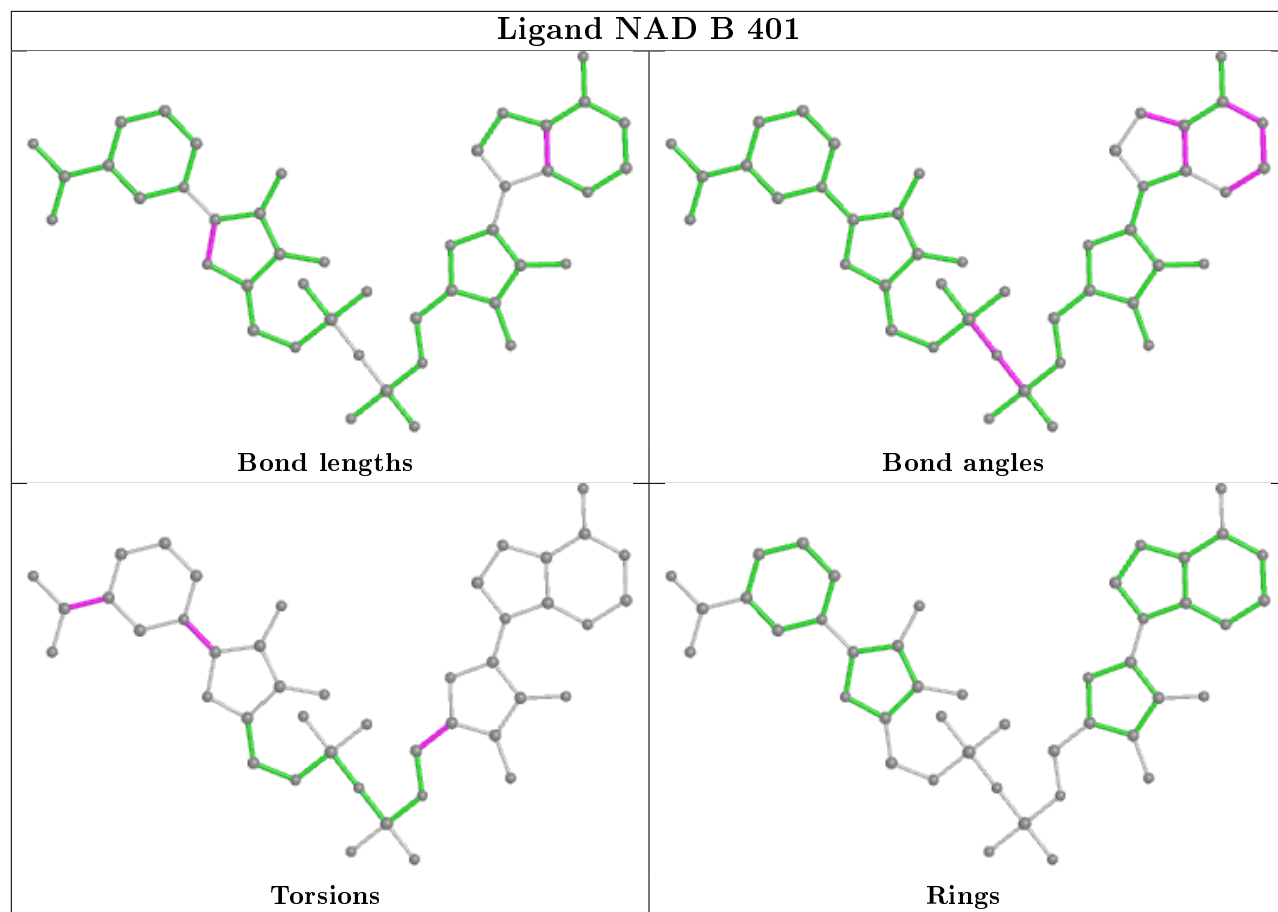
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAD	2	0
2	D	401	NAD	1	0
3	C	402	EDO	5	0

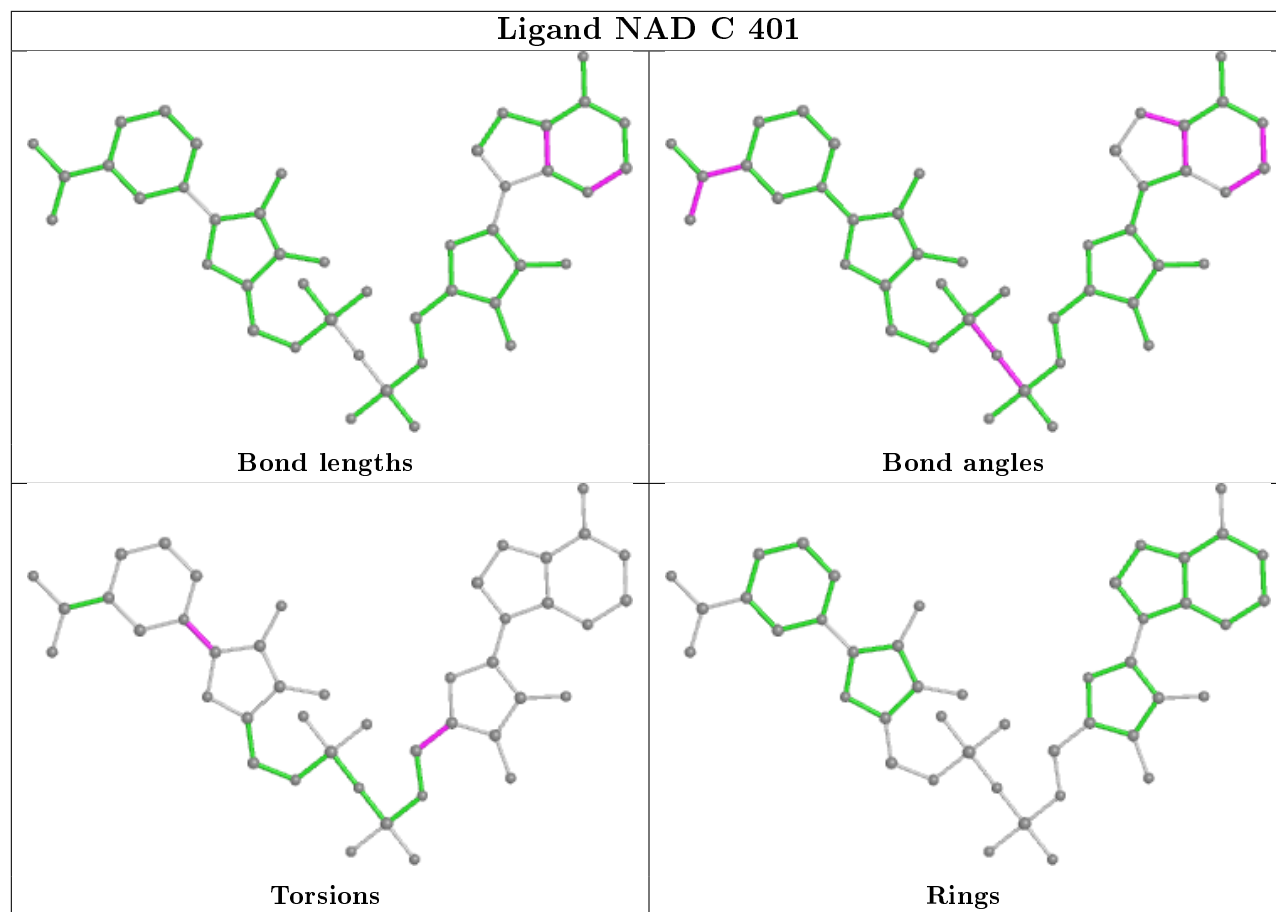
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	332/356 (93%)	0.50	25 (7%) 14 11	32, 65, 90, 102	0
1	B	335/356 (94%)	0.06	2 (0%) 89 89	29, 52, 75, 86	0
1	C	334/356 (93%)	0.30	14 (4%) 36 33	32, 59, 85, 119	0
1	D	344/356 (96%)	0.11	3 (0%) 84 85	32, 52, 73, 93	0
All	All	1345/1424 (94%)	0.24	44 (3%) 46 43	29, 56, 84, 119	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	125	GLY	6.5
1	A	117	LYS	4.4
1	A	116	LYS	4.1
1	A	85	TRP	3.9
1	D	304	GLY	3.7
1	C	127	ASP	3.6
1	A	118	VAL	3.4
1	A	81	ALA	3.1
1	A	86	ALA	3.1
1	A	90	VAL	3.0
1	A	164	LEU	2.9
1	A	115	ALA	2.8
1	A	26	GLY	2.8
1	C	70	GLN	2.8
1	A	5	VAL	2.8
1	A	168	PHE	2.8
1	A	89	GLY	2.8
1	A	305	ASN	2.7
1	C	126	ASN	2.7
1	C	250	LEU	2.7
1	B	303	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	146	VAL	2.6
1	C	248	ALA	2.6
1	A	29	VAL	2.5
1	C	168	PHE	2.5
1	A	307	LEU	2.5
1	C	118	VAL	2.5
1	C	251	GLU	2.5
1	C	231	GLY	2.4
1	B	168	PHE	2.4
1	A	91	GLU	2.4
1	C	78	ARG	2.4
1	C	92	ILE	2.4
1	D	227	GLY	2.3
1	A	254	VAL	2.2
1	D	126	ASN	2.1
1	C	216	LYS	2.1
1	A	78	ARG	2.1
1	A	248	ALA	2.1
1	A	308	VAL	2.1
1	A	331	PHE	2.0
1	A	110	ILE	2.0
1	A	161	ALA	2.0
1	C	308	VAL	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 carbohydrates in this entry.

6.4 Ligands [i](#)

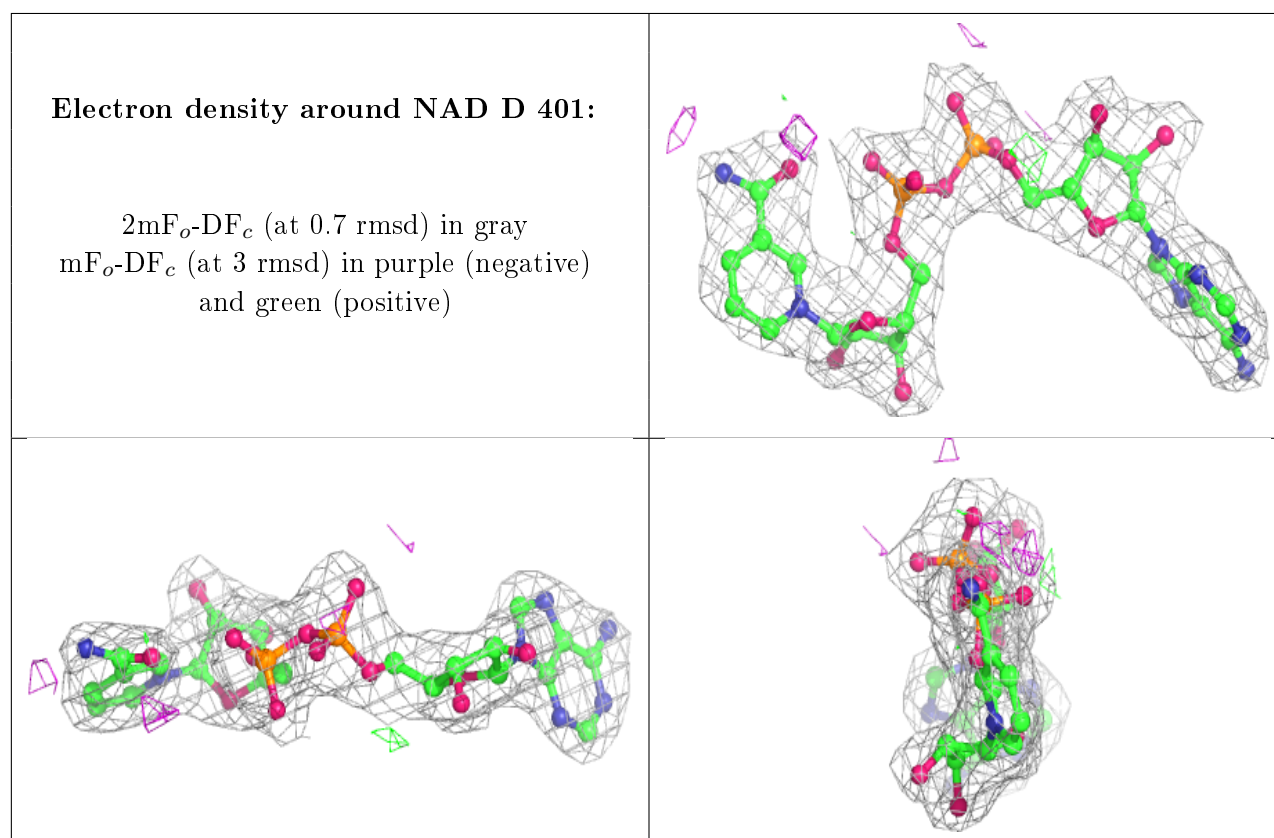
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

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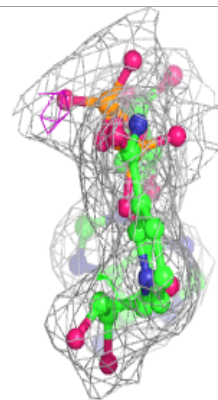
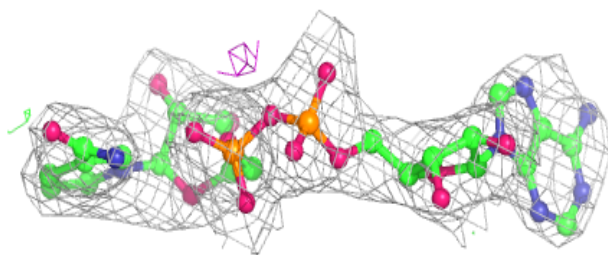
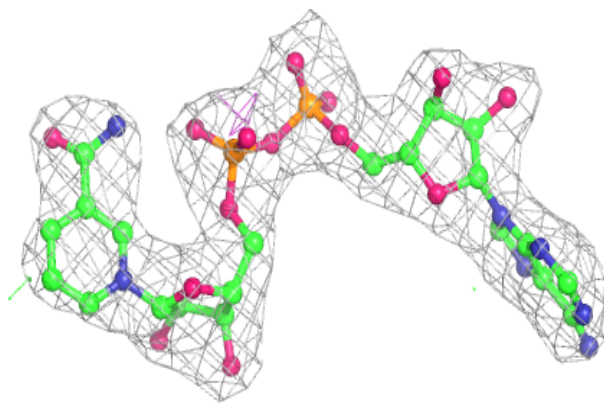
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	EDO	C	403	4/4	0.84	0.21	55,55,56,58	0
3	EDO	C	402	4/4	0.90	0.32	66,66,68,69	0
2	NAD	D	401	44/44	0.97	0.13	32,39,46,50	0
2	NAD	A	401	44/44	0.97	0.13	39,50,63,64	0
2	NAD	C	401	44/44	0.97	0.13	40,50,54,55	0
2	NAD	B	401	44/44	0.98	0.13	35,41,52,54	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.

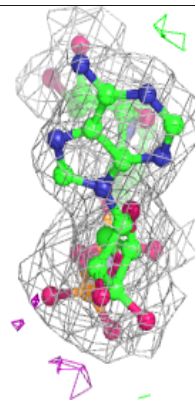
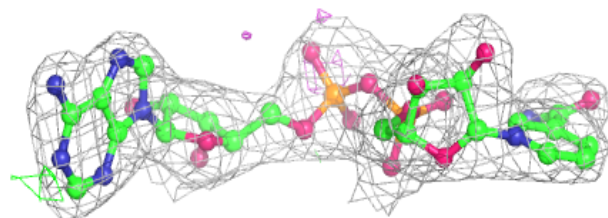
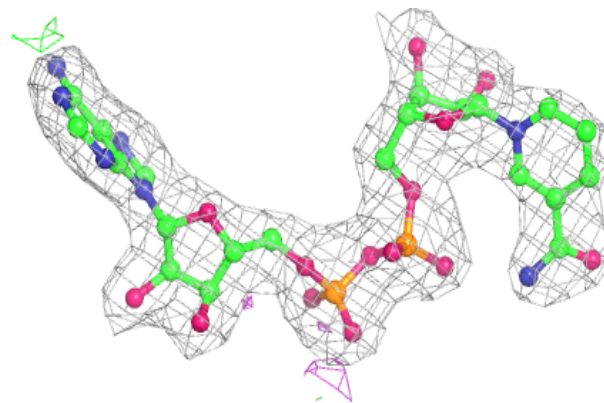


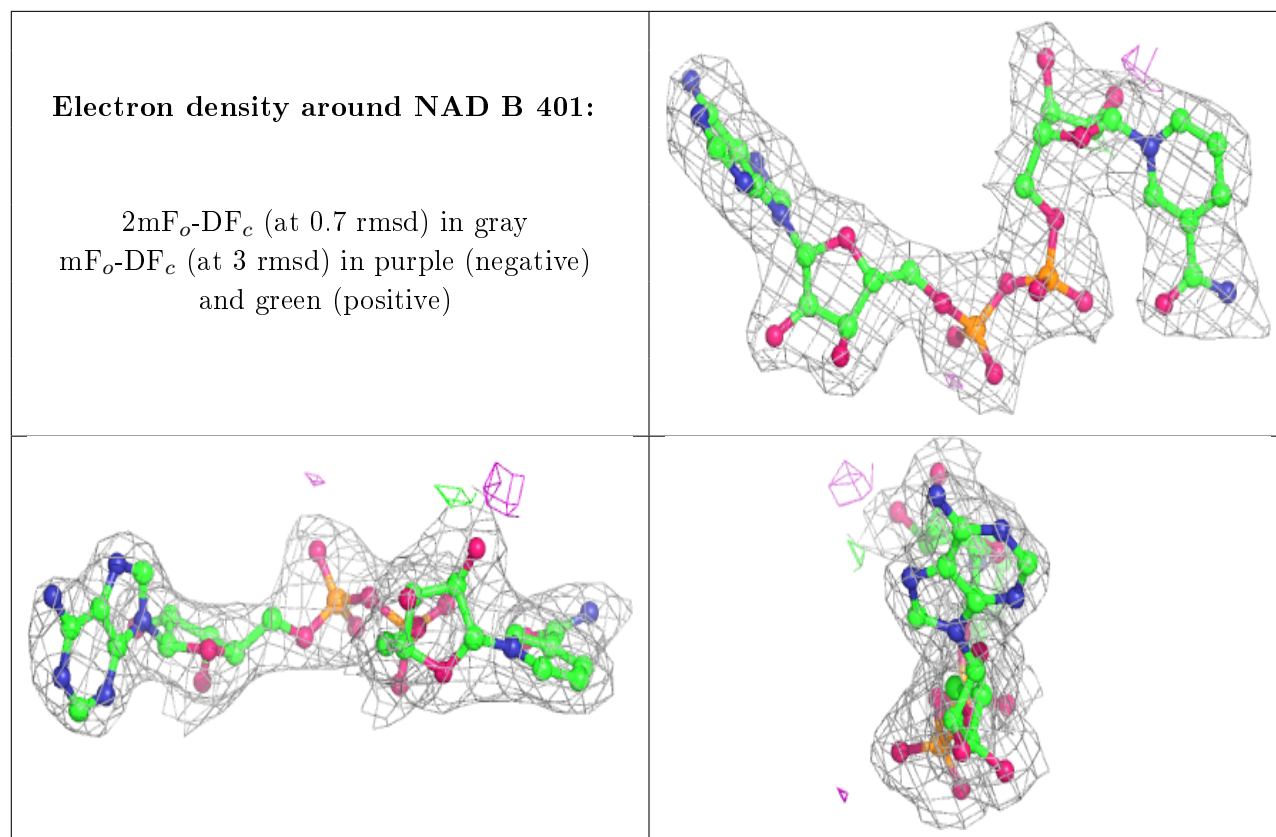
Electron density around NAD A 401:

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

**Electron density around NAD C 401:**

$2mF_o-DF_c$ (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.