



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2023 – 12:21 PM JST

PDB ID : 8HYE
Title : Structure of amino acid dehydrogenase-2752 with ligand
Authors : Sakuraba, H.; Ohshima, T.
Deposited on : 2023-01-06
Resolution : 2.20 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.32.2
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) : 2.32.2

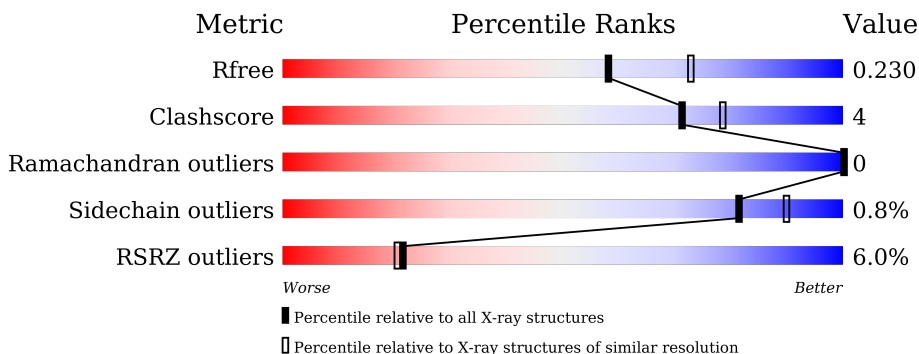
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	372	
1	B	372	
1	C	372	

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
4	EPE	A	1003	-	-	X	-

2 Entry composition [i](#)

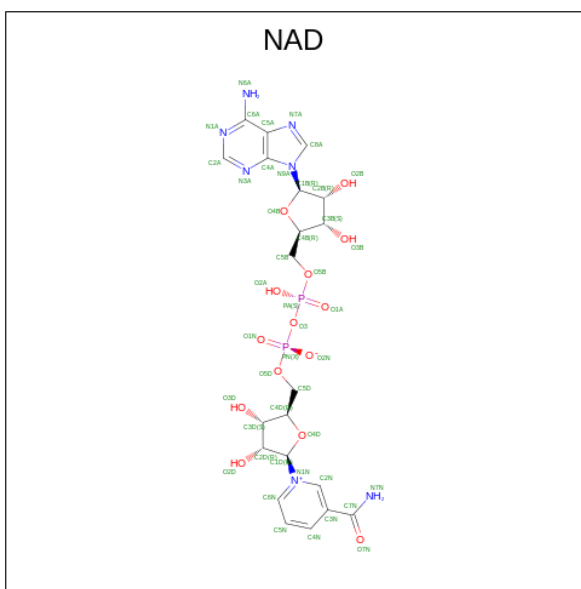
There are 6 unique types of molecules in this entry. The entry contains 8944 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alanine dehydrogenase.

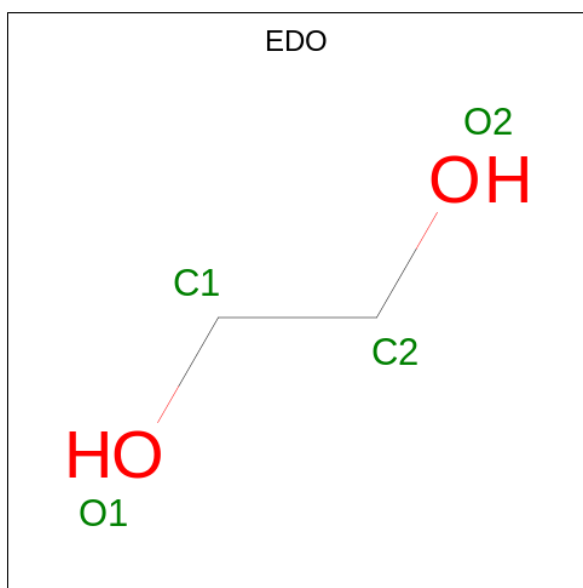
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	366	Total 2729	C 1731	N 465	O 521	S 12	0	0	0
1	B	367	Total 2738	C 1737	N 467	O 522	S 12	0	0	0
1	C	371	Total 2762	C 1753	N 471	O 526	S 12	0	0	0

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



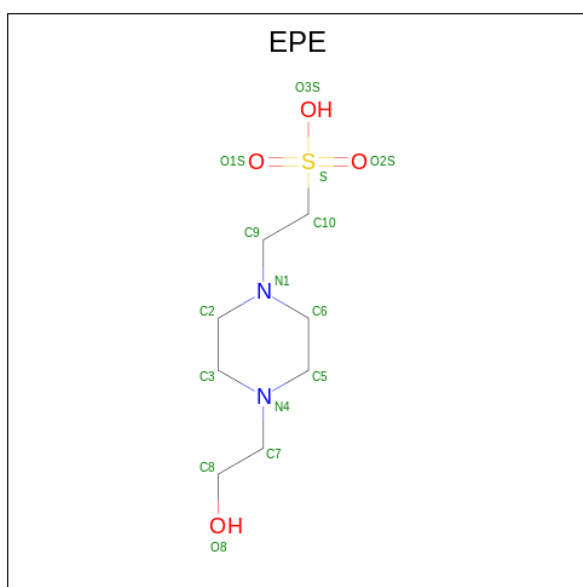
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	B	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	C	1	Total 44	C 21	N 7	O 14	P 2	0	0

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



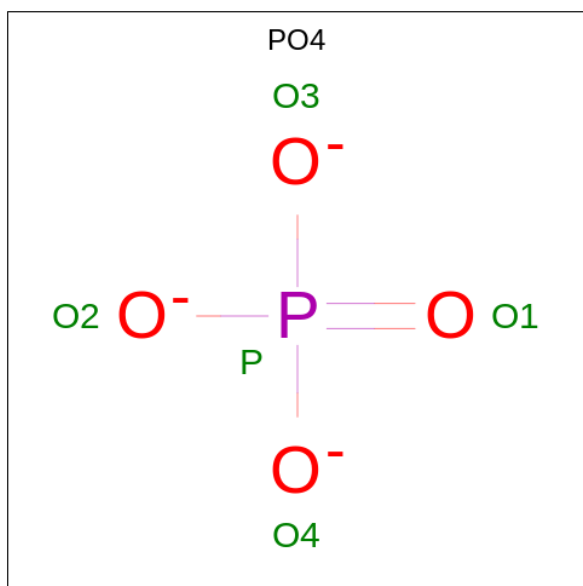
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	A	1	15	8	2	4	1	0	0

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
5	C	1	5	4	1	0	0

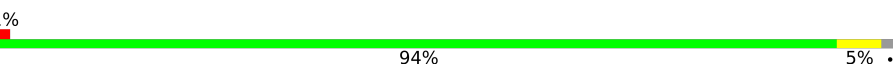
- Molecule 6 is water.

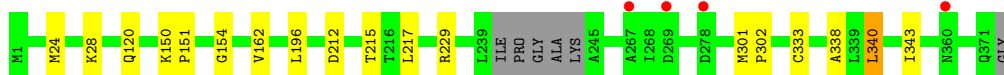
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	213	Total	O	0	0
			213	213		
6	B	228	Total	O	0	0
			228	228		
6	C	114	Total	O	0	0
			114	114		

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: Alanine dehydrogenase

Chain A:  94% 5%




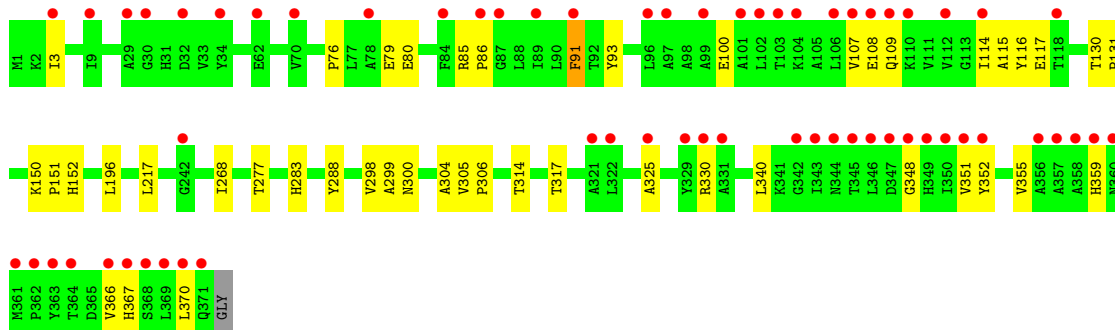
- Molecule 1: Alanine dehydrogenase

Chain B:  93% 6%



- Molecule 1: Alanine dehydrogenase

Chain C:  17% 87% 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	194.60Å 194.60Å 158.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.69 – 2.20 49.65 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.69-2.20) 99.9 (49.65-2.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.183 , 0.225 0.194 , 0.230	Depositor DCC
R_{free} test set	4494 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtrriage
Anisotropy	0.491	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8944	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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: PO4, EDO, NAD, EPE

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	0.76	0/2775	0.86	0/3774
1	B	0.77	0/2784	0.87	0/3785
1	C	0.79	0/2810	0.89	0/3823
All	All	0.77	0/8369	0.87	0/11382

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	2729	0	2772	12	0
1	B	2738	0	2785	12	0
1	C	2762	0	2812	28	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	1	0
3	A	4	0	6	1	0
3	C	4	0	6	0	0
4	A	15	0	18	8	0
5	C	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	213	0	0	8	0
6	B	228	0	0	3	0
6	C	114	0	0	1	0
All	All	8944	0	8477	59	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 4.

All (59) 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:212:ASP:O	4:A:1003:EPE:H21	1.63	0.97
4:A:1003:EPE:H52	6:A:1270:HOH:O	1.70	0.90
4:A:1003:EPE:H71	6:A:1206:HOH:O	1.72	0.89
1:C:100:GLU:HG2	1:C:359:HIS:CE1	2.10	0.85
4:A:1003:EPE:H22	6:A:1207:HOH:O	1.83	0.77
4:A:1003:EPE:C2	6:A:1207:HOH:O	2.36	0.71
1:C:340:LEU:HD21	1:C:367:HIS:CE1	2.33	0.63
1:B:349:HIS:HE1	6:B:1298:HOH:O	1.81	0.62
1:C:108:GLU:OE1	1:C:109:GLN:NE2	2.33	0.61
1:C:298:VAL:HG12	1:C:299:ALA:O	2.04	0.58
1:C:300:ASN:H	2:C:1001:NAD:H72N	1.53	0.57
1:C:93:TYR:CD1	1:C:116:TYR:HB2	2.43	0.54
1:C:352:TYR:HB3	1:C:355:VAL:HG12	1.90	0.53
1:C:85:ARG:HB3	1:C:86:PRO:HD2	1.91	0.53
1:C:115:ALA:HB1	1:C:117:GLU:OE2	2.10	0.52
1:A:340:LEU:HD12	1:A:340:LEU:O	2.10	0.52
1:B:169:LYS:HE2	6:B:1117:HOH:O	2.09	0.52
1:A:229:ARG:NH1	6:A:1107:HOH:O	2.43	0.50
1:B:283:HIS:CE1	1:B:304:ALA:HB2	2.46	0.50
1:A:24:MET:SD	1:A:28:LYS:HE3	2.52	0.49
1:B:74:LYS:NZ	6:B:1114:HOH:O	2.45	0.49
1:C:283:HIS:CE1	1:C:304:ALA:HB2	2.48	0.49
1:C:107:VAL:O	1:C:348:GLY:HA2	2.13	0.48
1:C:150:LYS:HB3	1:C:151:PRO:HD3	1.96	0.48
1:C:130:THR:N	1:C:131:PRO:HD2	2.29	0.48
1:C:304:ALA:C	1:C:306:PRO:HD3	2.34	0.48
1:C:330:ARG:NH1	1:C:370:LEU:O	2.43	0.48
1:C:76:PRO:HA	1:C:80:GLU:OE1	2.15	0.47
1:A:301:MET:N	1:A:302:PRO:CD	2.79	0.46
1:B:277:THR:HA	1:B:288:TYR:OH	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:THR:OG1	4:A:1003:EPE:H102	2.16	0.46
1:C:152:HIS:HE1	6:C:1136:HOH:O	1.97	0.46
1:A:24:MET:CG	1:A:28:LYS:HE3	2.46	0.46
1:B:150:LYS:HB3	1:B:151:PRO:HD3	1.97	0.46
1:C:91:PHE:CE1	1:C:114:ILE:HG22	2.51	0.46
1:A:150:LYS:HD2	1:A:154:GLY:O	2.15	0.45
1:B:196:LEU:HA	1:B:217:LEU:O	2.16	0.45
1:A:333:CYS:SG	1:A:343:ILE:HD11	2.56	0.45
1:B:10:LYS:HE3	1:B:75:GLU:OE2	2.16	0.45
1:C:314:THR:HA	1:C:317:THR:OG1	2.17	0.45
1:A:150:LYS:HB3	1:A:151:PRO:HD3	1.99	0.45
1:C:268:ILE:HD13	1:C:277:THR:HG21	1.99	0.44
1:B:150:LYS:N	1:B:151:PRO:CD	2.81	0.44
1:B:148:LEU:HD23	1:B:148:LEU:HA	1.90	0.43
1:C:79:GLU:H	1:C:79:GLU:HG2	1.56	0.43
4:A:1003:EPE:H72	6:A:1270:HOH:O	2.18	0.43
1:C:150:LYS:N	1:C:151:PRO:CD	2.81	0.43
1:C:351:VAL:HB	1:C:366:VAL:HB	2.00	0.43
1:B:239:LEU:HG	1:B:239:LEU:O	2.19	0.42
1:C:277:THR:HA	1:C:288:TYR:OH	2.20	0.42
1:A:196:LEU:HA	1:A:217:LEU:O	2.18	0.42
1:B:162:VAL:HB	1:B:163:PRO:CD	2.50	0.41
1:C:305:VAL:N	1:C:306:PRO:HD3	2.35	0.41
1:C:3:ILE:HD11	1:C:325:ALA:HB2	2.03	0.41
4:A:1003:EPE:H92	6:A:1159:HOH:O	2.20	0.40
1:C:3:ILE:HD11	1:C:325:ALA:CB	2.51	0.40
1:C:196:LEU:HA	1:C:217:LEU:O	2.21	0.40
1:A:120:GLN:O	1:A:338:ALA:HA	2.22	0.40
3:A:1002:EDO:H22	6:A:1165:HOH:O	2.21	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	362/372 (97%)	357 (99%)	5 (1%)	0	100	100
1	B	363/372 (98%)	361 (99%)	2 (1%)	0	100	100
1	C	369/372 (99%)	359 (97%)	10 (3%)	0	100	100
All	All	1094/1116 (98%)	1077 (98%)	17 (2%)	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	287/290 (99%)	285 (99%)	2 (1%)	84	91
1	B	288/290 (99%)	284 (99%)	4 (1%)	67	80
1	C	290/290 (100%)	289 (100%)	1 (0%)	92	97
All	All	865/870 (99%)	858 (99%)	7 (1%)	81	90

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	VAL
1	A	340	LEU
1	B	195	ILE
1	B	340	LEU
1	B	344	ASN
1	B	368	SER
1	C	91	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	152	HIS
1	A	270	GLN
1	A	371	GLN

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Mol	Chain	Res	Type
1	B	283	HIS
1	B	291	HIS
1	C	152	HIS
1	C	270	GLN
1	C	291	HIS
1	C	344	ASN
1	C	367	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](#)

7 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	1001	-	42,48,48	1.00	2 (4%)	50,73,73	1.69	13 (26%)
2	NAD	B	1001	-	42,48,48	1.16	4 (9%)	50,73,73	1.47	10 (20%)
3	EDO	C	1003	-	3,3,3	1.02	0	2,2,2	0.61	0
4	EPE	A	1003	-	15,15,15	2.03	1 (6%)	18,20,20	1.13	3 (16%)
2	NAD	C	1001	-	42,48,48	0.94	1 (2%)	50,73,73	2.35	12 (24%)
5	PO4	C	1002	-	4,4,4	0.96	0	6,6,6	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	A	1002	-	3,3,3	0.39	0	2,2,2	0.61	0

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	1001	-	-	12/26/62/62	0/5/5/5
2	NAD	B	1001	-	-	12/26/62/62	0/5/5/5
3	EDO	C	1003	-	-	1/1/1/1	-
4	EPE	A	1003	-	-	7/9/19/19	0/1/1/1
2	NAD	C	1001	-	-	9/26/62/62	0/5/5/5
3	EDO	A	1002	-	-	0/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1003	EPE	C10-S	-7.62	1.66	1.77
2	B	1001	NAD	C2A-N3A	2.97	1.36	1.32
2	A	1001	NAD	O4B-C1B	2.90	1.45	1.41
2	B	1001	NAD	O2B-C2B	2.90	1.49	1.43
2	B	1001	NAD	C5A-C4A	2.84	1.48	1.40
2	C	1001	NAD	C5A-C4A	2.55	1.47	1.40
2	A	1001	NAD	C5A-C4A	2.49	1.47	1.40
2	B	1001	NAD	O4D-C1D	2.24	1.44	1.41

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	NAD	C3N-C7N-N7N	-8.50	107.55	117.75
2	C	1001	NAD	O7N-C7N-C3N	7.00	128.00	119.63
2	C	1001	NAD	C6N-N1N-C2N	-4.88	117.52	121.97
2	C	1001	NAD	O4D-C1D-C2D	-4.43	100.45	106.93
2	A	1001	NAD	C4A-C5A-N7A	-4.41	104.81	109.40
2	C	1001	NAD	C1B-N9A-C4A	-4.35	119.00	126.64
2	C	1001	NAD	N3A-C2A-N1A	-3.91	122.56	128.68
2	B	1001	NAD	C1B-N9A-C4A	-3.87	119.84	126.64
2	A	1001	NAD	N3A-C2A-N1A	-3.70	122.90	128.68
2	A	1001	NAD	C1B-N9A-C4A	-3.40	120.68	126.64
2	C	1001	NAD	PN-O3-PA	-3.26	121.63	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	NAD	C3N-C2N-N1N	3.21	123.57	120.43
2	C	1001	NAD	O4B-C1B-C2B	-3.16	102.31	106.93
2	B	1001	NAD	C2N-N1N-C1D	-3.09	112.25	119.14
2	C	1001	NAD	C5D-C4D-C3D	-2.96	104.08	115.18
2	B	1001	NAD	N3A-C2A-N1A	-2.82	124.27	128.68
2	A	1001	NAD	C2N-N1N-C1D	-2.63	113.28	119.14
2	B	1001	NAD	O2N-PN-O1N	2.57	124.94	112.24
2	A	1001	NAD	O3D-C3D-C2D	2.53	120.02	111.82
2	B	1001	NAD	O4B-C1B-C2B	2.51	110.59	106.93
2	C	1001	NAD	C5N-C4N-C3N	-2.34	117.57	120.34
2	A	1001	NAD	O5B-PA-O1A	-2.32	100.01	109.07
4	A	1003	EPE	O3S-S-C10	2.31	109.50	105.77
2	B	1001	NAD	N6A-C6A-N1A	2.29	123.33	118.57
2	A	1001	NAD	C6N-N1N-C2N	-2.25	119.92	121.97
2	B	1001	NAD	C5A-C6A-N6A	-2.23	116.97	120.35
2	A	1001	NAD	O2A-PA-O1A	2.22	123.21	112.24
2	A	1001	NAD	C3N-C2N-N1N	2.19	122.57	120.43
2	A	1001	NAD	C2A-N1A-C6A	2.18	122.49	118.75
2	B	1001	NAD	C3N-C2N-N1N	2.18	122.56	120.43
2	A	1001	NAD	O2N-PN-O1N	2.15	122.86	112.24
4	A	1003	EPE	O2S-S-C10	2.14	109.49	106.92
4	A	1003	EPE	O1S-S-C10	2.12	109.46	106.92
2	A	1001	NAD	C5A-C6A-N6A	2.10	123.55	120.35
2	B	1001	NAD	C2B-C3B-C4B	2.08	106.68	102.64
2	C	1001	NAD	C4A-C5A-N7A	-2.07	107.25	109.40
2	A	1001	NAD	C4N-C3N-C7N	2.04	126.49	121.04
2	B	1001	NAD	C2D-C3D-C4D	2.03	106.59	102.64

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	NAD	C5D-O5D-PN-O3
2	A	1001	NAD	C5D-O5D-PN-O1N
2	A	1001	NAD	C5D-O5D-PN-O2N
2	A	1001	NAD	O4D-C4D-C5D-O5D
2	A	1001	NAD	C3D-C4D-C5D-O5D
2	A	1001	NAD	O4D-C1D-N1N-C6N
2	B	1001	NAD	C5B-O5B-PA-O3
2	B	1001	NAD	O4D-C4D-C5D-O5D
2	B	1001	NAD	O4D-C1D-N1N-C2N
2	B	1001	NAD	O4D-C1D-N1N-C6N

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Mol	Chain	Res	Type	Atoms
2	B	1001	NAD	C2D-C1D-N1N-C2N
2	B	1001	NAD	C2D-C1D-N1N-C6N
2	C	1001	NAD	C5D-O5D-PN-O2N
2	C	1001	NAD	O4D-C1D-N1N-C2N
2	C	1001	NAD	O4D-C1D-N1N-C6N
2	C	1001	NAD	C2D-C1D-N1N-C6N
4	A	1003	EPE	S-C10-C9-N1
4	A	1003	EPE	C9-C10-S-O1S
4	A	1003	EPE	C9-C10-S-O3S
2	B	1001	NAD	C3D-C4D-C5D-O5D
2	B	1001	NAD	O4B-C4B-C5B-O5B
2	A	1001	NAD	C2N-C3N-C7N-N7N
3	C	1003	EDO	O1-C1-C2-O2
2	C	1001	NAD	C5D-O5D-PN-O3
2	A	1001	NAD	PN-O3-PA-O1A
2	C	1001	NAD	C2N-C3N-C7N-N7N
4	A	1003	EPE	C8-C7-N4-C5
2	B	1001	NAD	C5B-O5B-PA-O1A
2	B	1001	NAD	C5B-O5B-PA-O2A
2	C	1001	NAD	C5D-O5D-PN-O1N
4	A	1003	EPE	C9-C10-S-O2S
2	A	1001	NAD	C2N-C3N-C7N-O7N
2	A	1001	NAD	PN-O3-PA-O2A
2	A	1001	NAD	C4N-C3N-C7N-N7N
2	C	1001	NAD	C2N-C3N-C7N-O7N
4	A	1003	EPE	C10-C9-N1-C2
4	A	1003	EPE	C10-C9-N1-C6
2	B	1001	NAD	PN-O3-PA-O1A
2	C	1001	NAD	O4B-C4B-C5B-O5B
2	B	1001	NAD	PN-O3-PA-O2A
2	A	1001	NAD	O4B-C4B-C5B-O5B

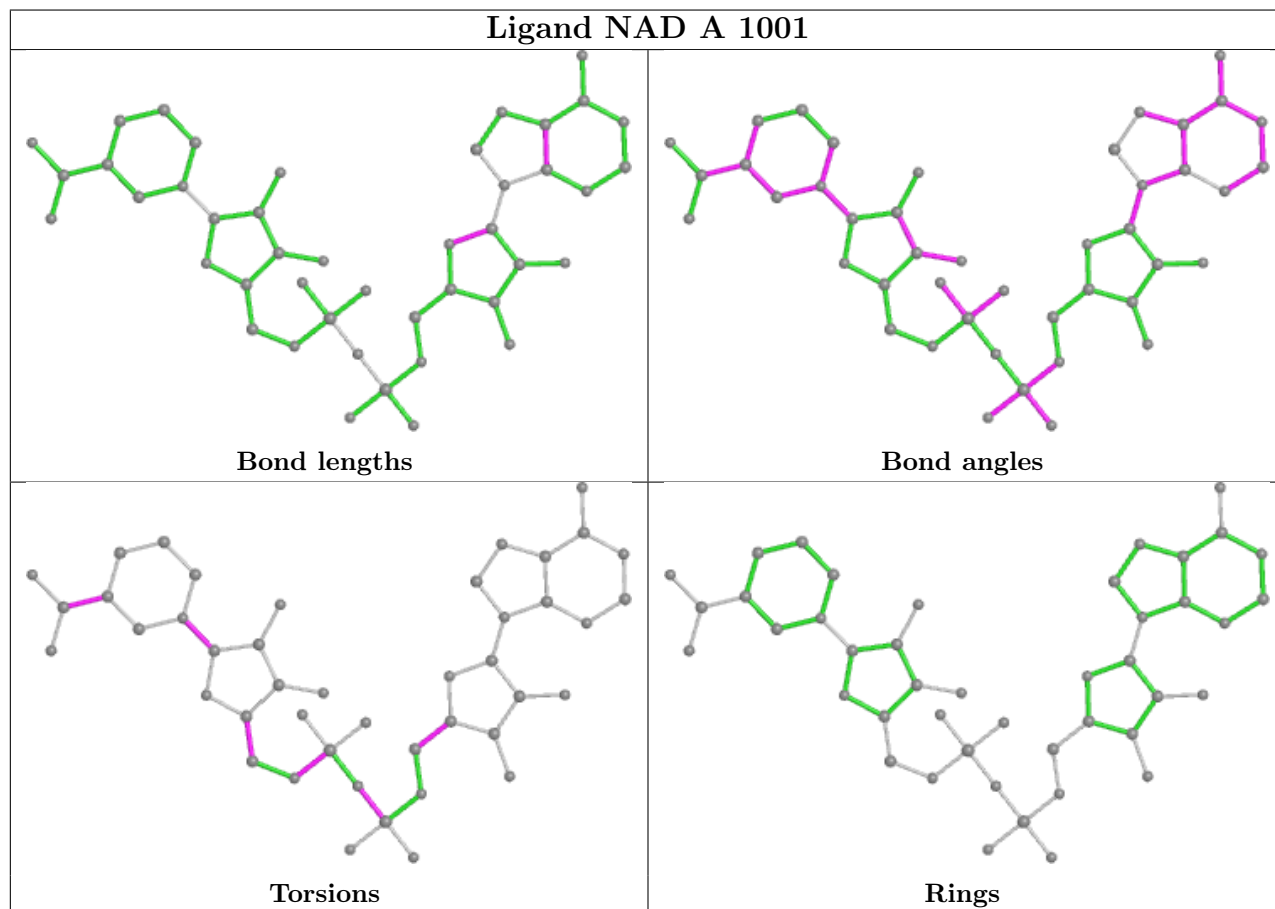
There are no ring outliers.

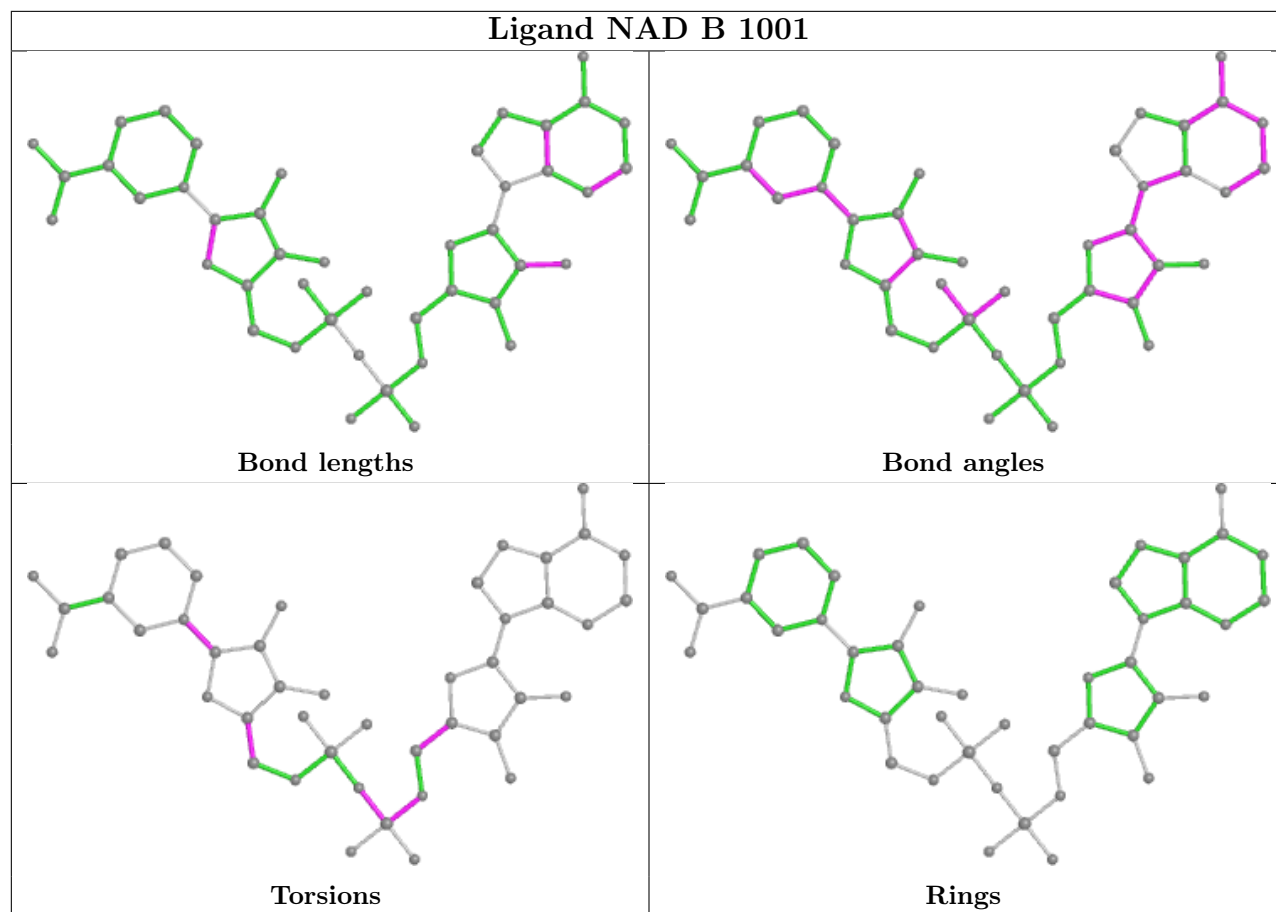
3 monomers are involved in 10 short contacts:

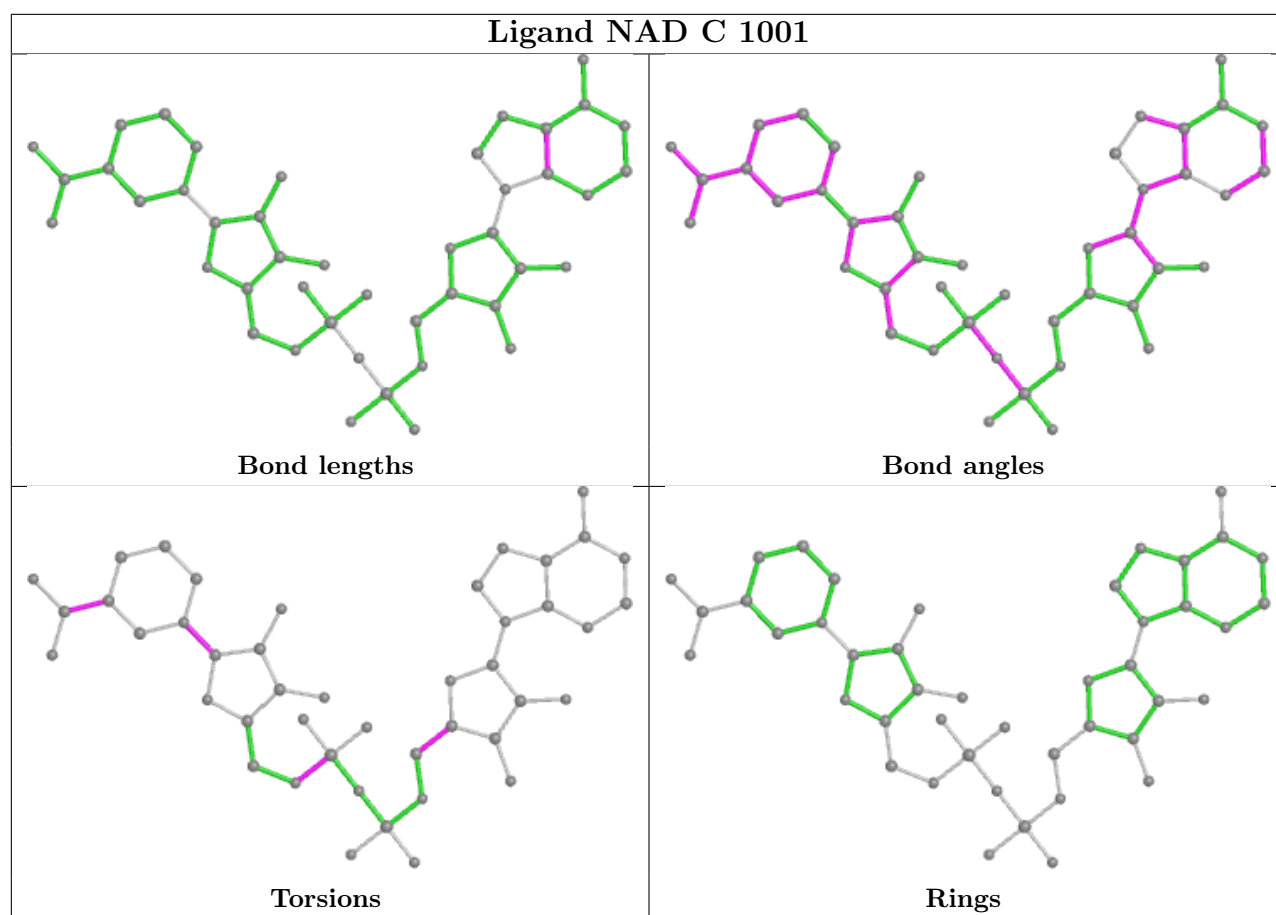
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1003	EPE	8	0
2	C	1001	NAD	1	0
3	A	1002	EDO	1	0

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

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	366/372 (98%)	-0.50	4 (1%) 80 79	25, 35, 61, 80	0
1	B	367/372 (98%)	-0.54	0 100 100	25, 35, 57, 94	0
1	C	371/372 (99%)	0.59	62 (16%) 1 1	25, 54, 102, 117	0
All	All	1104/1116 (98%)	-0.15	66 (5%) 21 20	25, 37, 83, 117	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	351	VAL	9.1
1	C	364	THR	8.7
1	C	368	SER	6.5
1	C	360	ASN	5.6
1	C	346	LEU	5.5
1	C	349	HIS	5.4
1	C	370	LEU	5.1
1	C	369	LEU	4.9
1	C	347	ASP	4.9
1	C	366	VAL	4.7
1	C	358	ALA	4.7
1	C	102	LEU	4.5
1	C	114	ILE	4.3
1	C	99	ALA	4.2
1	C	363	TYR	4.2
1	C	371	GLN	4.2
1	C	107	VAL	4.2
1	C	357	ALA	4.1
1	C	367	HIS	4.1
1	C	97	ALA	4.0
1	C	343	ILE	3.9
1	C	112	VAL	3.8
1	C	352	TYR	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	96	LEU	3.5
1	C	86	PRO	3.3
1	C	118	THR	3.3
1	C	329	TYR	3.1
1	C	110	LYS	3.0
1	C	78	ALA	3.0
1	C	331	ALA	3.0
1	C	350	ILE	3.0
1	C	108	GLU	3.0
1	C	104	LYS	3.0
1	C	348	GLY	2.9
1	C	70	VAL	2.9
1	C	361	MET	2.9
1	C	91	PHE	2.8
1	C	242	GLY	2.8
1	C	345	THR	2.8
1	C	106	LEU	2.7
1	C	322	LEU	2.7
1	C	30	GLY	2.6
1	C	84	PHE	2.6
1	C	356	ALA	2.6
1	C	103	THR	2.6
1	C	330	ARG	2.6
1	C	344	ASN	2.6
1	C	359	HIS	2.5
1	C	342	GLY	2.5
1	C	29	ALA	2.5
1	C	34	TYR	2.5
1	C	89	ILE	2.4
1	C	362	PRO	2.4
1	C	325	ALA	2.4
1	C	3	ILE	2.4
1	A	360	ASN	2.3
1	A	278	ASP	2.3
1	C	101	ALA	2.3
1	C	32	ASP	2.2
1	C	9	ILE	2.2
1	C	87	GLY	2.2
1	C	109	GLN	2.2
1	C	62	GLU	2.1
1	A	267	ALA	2.1
1	C	321	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	269	ASP	2.1

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](#)

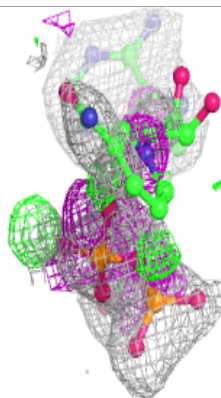
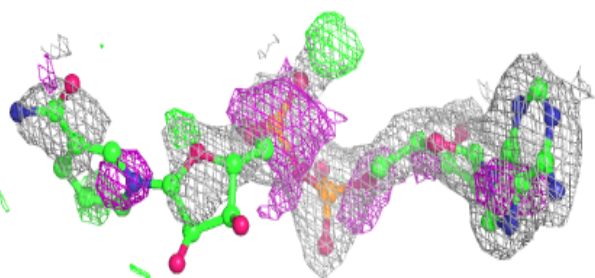
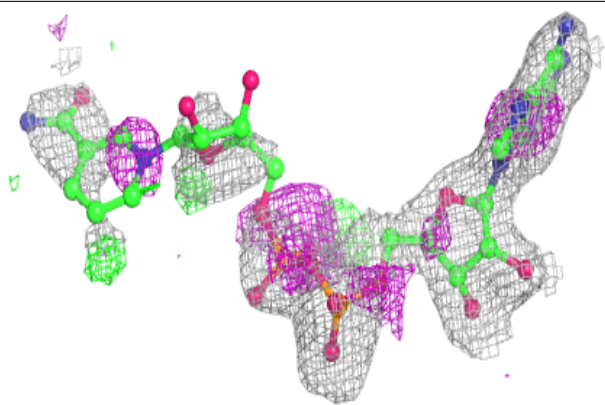
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EPE	A	1003	15/15	0.69	0.31	29,40,108,117	0
3	EDO	C	1003	4/4	0.79	0.22	44,57,59,64	0
5	PO4	C	1002	5/5	0.79	0.14	46,53,64,91	0
2	NAD	B	1001	44/44	0.86	0.24	38,77,139,153	0
3	EDO	A	1002	4/4	0.88	0.31	63,66,67,68	0
2	NAD	A	1001	44/44	0.88	0.24	46,74,126,136	0
2	NAD	C	1001	44/44	0.94	0.13	32,40,103,115	0

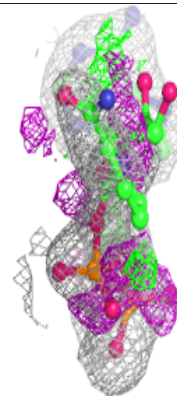
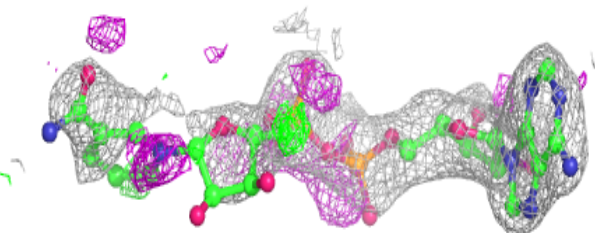
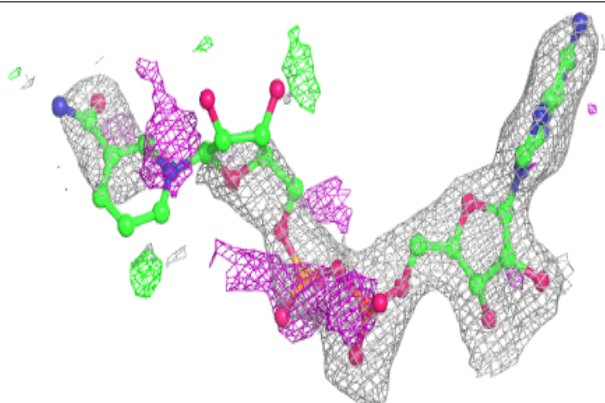
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

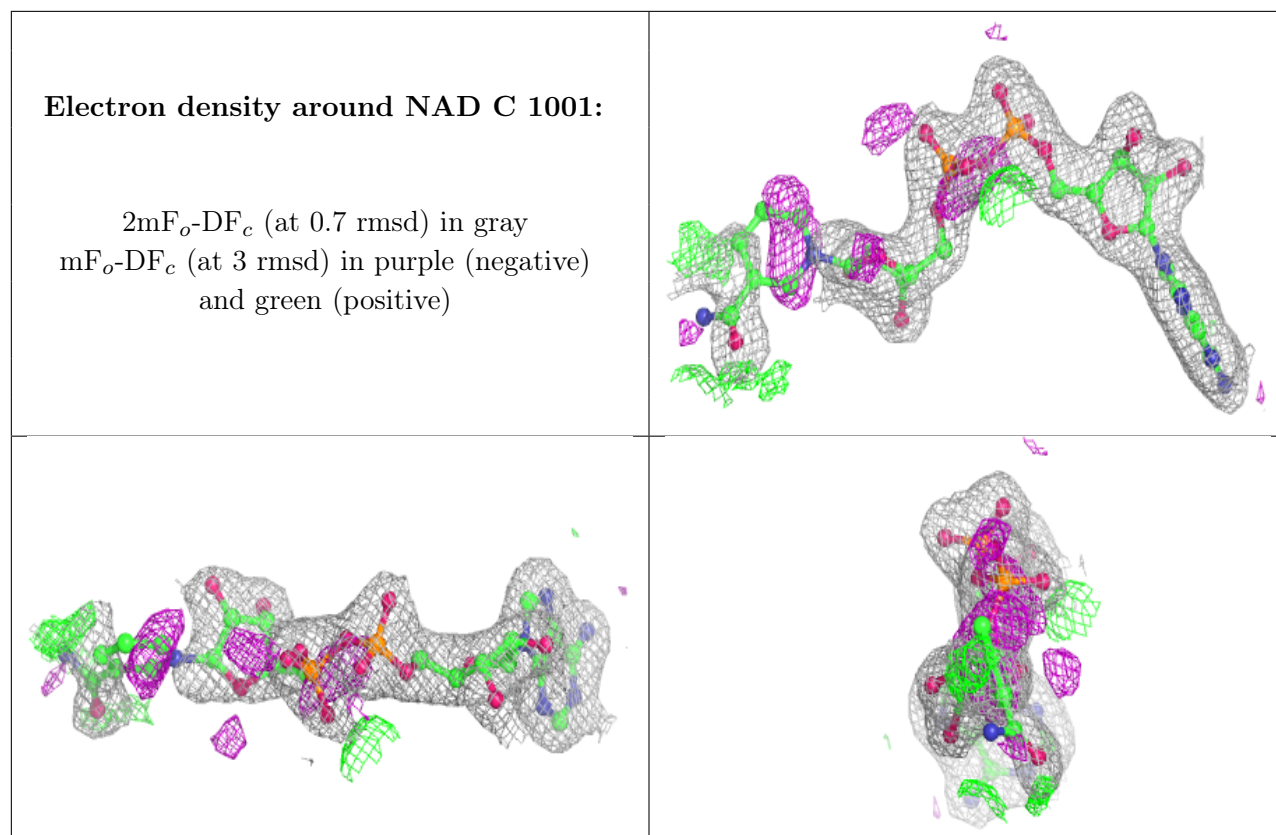
Electron density around NAD B 1001:

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

**Electron density around NAD A 1001:**

$2mF_o-DF_c$ (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.