



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:59 AM UTC

PDB ID : 4QN2 / pdb_00004qn2
Title : 2.6 Angstrom resolution crystal structure of betaine aldehyde dehydrogenase (betB) G234S mutant from *Staphylococcus aureus* (IDP00699) in complex with NAD⁺ and BME-free Cys289
Authors : Halavaty, A.S.; Minasov, G.; Chen, C.; Joo, J.C.; Yakunin, A.F.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2014-06-17
Resolution : 2.60 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](#) ①) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

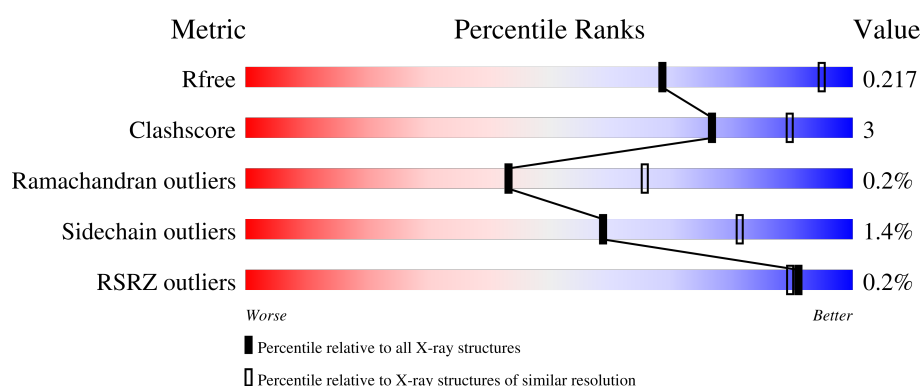
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	517	87% 9% . .
1	B	517	87% 8% . .
1	C	517	87% 8% .
1	D	517	88% 8% .
1	E	517	87% 9% . .

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Mol	Chain	Length	Quality of chain
1	F	517	 88% 7% . .
1	G	517	 88% 7% . .
1	H	517	 86% 9% . .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 31934 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 Betaine aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	502	Total	C	N	O	S	0	0	0
			3899	2462	659	764	14			
1	B	496	Total	C	N	O	S	0	4	0
			3881	2445	655	767	14			
1	C	496	Total	C	N	O	S	0	0	0
			3847	2427	651	755	14			
1	D	496	Total	C	N	O	S	0	1	0
			3856	2432	652	758	14			
1	E	496	Total	C	N	O	S	0	0	0
			3847	2427	651	755	14			
1	F	496	Total	C	N	O	S	0	0	0
			3847	2427	651	755	14			
1	G	496	Total	C	N	O	S	0	0	0
			3847	2427	651	755	14			
1	H	496	Total	C	N	O	S	0	0	0
			3847	2427	651	755	14			

There are 176 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q5HCU0
A	-19	GLY	-	expression tag	UNP Q5HCU0
A	-18	SER	-	expression tag	UNP Q5HCU0
A	-17	SER	-	expression tag	UNP Q5HCU0
A	-16	HIS	-	expression tag	UNP Q5HCU0
A	-15	HIS	-	expression tag	UNP Q5HCU0
A	-14	HIS	-	expression tag	UNP Q5HCU0
A	-13	HIS	-	expression tag	UNP Q5HCU0
A	-12	HIS	-	expression tag	UNP Q5HCU0
A	-11	HIS	-	expression tag	UNP Q5HCU0
A	-10	SER	-	expression tag	UNP Q5HCU0
A	-9	SER	-	expression tag	UNP Q5HCU0
A	-8	GLY	-	expression tag	UNP Q5HCU0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	ARG	-	expression tag	UNP Q5HCU0
A	-6	GLU	-	expression tag	UNP Q5HCU0
A	-5	ASN	-	expression tag	UNP Q5HCU0
A	-4	LEU	-	expression tag	UNP Q5HCU0
A	-3	TYR	-	expression tag	UNP Q5HCU0
A	-2	PHE	-	expression tag	UNP Q5HCU0
A	-1	GLN	-	expression tag	UNP Q5HCU0
A	0	GLY	-	expression tag	UNP Q5HCU0
A	234	SER	GLY	engineered mutation	UNP Q5HCU0
B	-20	MET	-	expression tag	UNP Q5HCU0
B	-19	GLY	-	expression tag	UNP Q5HCU0
B	-18	SER	-	expression tag	UNP Q5HCU0
B	-17	SER	-	expression tag	UNP Q5HCU0
B	-16	HIS	-	expression tag	UNP Q5HCU0
B	-15	HIS	-	expression tag	UNP Q5HCU0
B	-14	HIS	-	expression tag	UNP Q5HCU0
B	-13	HIS	-	expression tag	UNP Q5HCU0
B	-12	HIS	-	expression tag	UNP Q5HCU0
B	-11	HIS	-	expression tag	UNP Q5HCU0
B	-10	SER	-	expression tag	UNP Q5HCU0
B	-9	SER	-	expression tag	UNP Q5HCU0
B	-8	GLY	-	expression tag	UNP Q5HCU0
B	-7	ARG	-	expression tag	UNP Q5HCU0
B	-6	GLU	-	expression tag	UNP Q5HCU0
B	-5	ASN	-	expression tag	UNP Q5HCU0
B	-4	LEU	-	expression tag	UNP Q5HCU0
B	-3	TYR	-	expression tag	UNP Q5HCU0
B	-2	PHE	-	expression tag	UNP Q5HCU0
B	-1	GLN	-	expression tag	UNP Q5HCU0
B	0	GLY	-	expression tag	UNP Q5HCU0
B	234	SER	GLY	engineered mutation	UNP Q5HCU0
C	-20	MET	-	expression tag	UNP Q5HCU0
C	-19	GLY	-	expression tag	UNP Q5HCU0
C	-18	SER	-	expression tag	UNP Q5HCU0
C	-17	SER	-	expression tag	UNP Q5HCU0
C	-16	HIS	-	expression tag	UNP Q5HCU0
C	-15	HIS	-	expression tag	UNP Q5HCU0
C	-14	HIS	-	expression tag	UNP Q5HCU0
C	-13	HIS	-	expression tag	UNP Q5HCU0
C	-12	HIS	-	expression tag	UNP Q5HCU0
C	-11	HIS	-	expression tag	UNP Q5HCU0
C	-10	SER	-	expression tag	UNP Q5HCU0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	SER	-	expression tag	UNP Q5HCU0
C	-8	GLY	-	expression tag	UNP Q5HCU0
C	-7	ARG	-	expression tag	UNP Q5HCU0
C	-6	GLU	-	expression tag	UNP Q5HCU0
C	-5	ASN	-	expression tag	UNP Q5HCU0
C	-4	LEU	-	expression tag	UNP Q5HCU0
C	-3	TYR	-	expression tag	UNP Q5HCU0
C	-2	PHE	-	expression tag	UNP Q5HCU0
C	-1	GLN	-	expression tag	UNP Q5HCU0
C	0	GLY	-	expression tag	UNP Q5HCU0
C	234	SER	GLY	engineered mutation	UNP Q5HCU0
D	-20	MET	-	expression tag	UNP Q5HCU0
D	-19	GLY	-	expression tag	UNP Q5HCU0
D	-18	SER	-	expression tag	UNP Q5HCU0
D	-17	SER	-	expression tag	UNP Q5HCU0
D	-16	HIS	-	expression tag	UNP Q5HCU0
D	-15	HIS	-	expression tag	UNP Q5HCU0
D	-14	HIS	-	expression tag	UNP Q5HCU0
D	-13	HIS	-	expression tag	UNP Q5HCU0
D	-12	HIS	-	expression tag	UNP Q5HCU0
D	-11	HIS	-	expression tag	UNP Q5HCU0
D	-10	SER	-	expression tag	UNP Q5HCU0
D	-9	SER	-	expression tag	UNP Q5HCU0
D	-8	GLY	-	expression tag	UNP Q5HCU0
D	-7	ARG	-	expression tag	UNP Q5HCU0
D	-6	GLU	-	expression tag	UNP Q5HCU0
D	-5	ASN	-	expression tag	UNP Q5HCU0
D	-4	LEU	-	expression tag	UNP Q5HCU0
D	-3	TYR	-	expression tag	UNP Q5HCU0
D	-2	PHE	-	expression tag	UNP Q5HCU0
D	-1	GLN	-	expression tag	UNP Q5HCU0
D	0	GLY	-	expression tag	UNP Q5HCU0
D	234	SER	GLY	engineered mutation	UNP Q5HCU0
E	-20	MET	-	expression tag	UNP Q5HCU0
E	-19	GLY	-	expression tag	UNP Q5HCU0
E	-18	SER	-	expression tag	UNP Q5HCU0
E	-17	SER	-	expression tag	UNP Q5HCU0
E	-16	HIS	-	expression tag	UNP Q5HCU0
E	-15	HIS	-	expression tag	UNP Q5HCU0
E	-14	HIS	-	expression tag	UNP Q5HCU0
E	-13	HIS	-	expression tag	UNP Q5HCU0
E	-12	HIS	-	expression tag	UNP Q5HCU0

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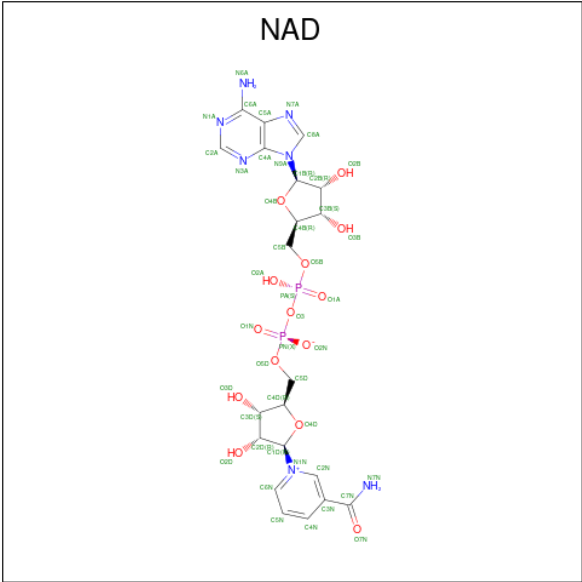
Chain	Residue	Modelled	Actual	Comment	Reference
E	-11	HIS	-	expression tag	UNP Q5HCU0
E	-10	SER	-	expression tag	UNP Q5HCU0
E	-9	SER	-	expression tag	UNP Q5HCU0
E	-8	GLY	-	expression tag	UNP Q5HCU0
E	-7	ARG	-	expression tag	UNP Q5HCU0
E	-6	GLU	-	expression tag	UNP Q5HCU0
E	-5	ASN	-	expression tag	UNP Q5HCU0
E	-4	LEU	-	expression tag	UNP Q5HCU0
E	-3	TYR	-	expression tag	UNP Q5HCU0
E	-2	PHE	-	expression tag	UNP Q5HCU0
E	-1	GLN	-	expression tag	UNP Q5HCU0
E	0	GLY	-	expression tag	UNP Q5HCU0
E	234	SER	GLY	engineered mutation	UNP Q5HCU0
F	-20	MET	-	expression tag	UNP Q5HCU0
F	-19	GLY	-	expression tag	UNP Q5HCU0
F	-18	SER	-	expression tag	UNP Q5HCU0
F	-17	SER	-	expression tag	UNP Q5HCU0
F	-16	HIS	-	expression tag	UNP Q5HCU0
F	-15	HIS	-	expression tag	UNP Q5HCU0
F	-14	HIS	-	expression tag	UNP Q5HCU0
F	-13	HIS	-	expression tag	UNP Q5HCU0
F	-12	HIS	-	expression tag	UNP Q5HCU0
F	-11	HIS	-	expression tag	UNP Q5HCU0
F	-10	SER	-	expression tag	UNP Q5HCU0
F	-9	SER	-	expression tag	UNP Q5HCU0
F	-8	GLY	-	expression tag	UNP Q5HCU0
F	-7	ARG	-	expression tag	UNP Q5HCU0
F	-6	GLU	-	expression tag	UNP Q5HCU0
F	-5	ASN	-	expression tag	UNP Q5HCU0
F	-4	LEU	-	expression tag	UNP Q5HCU0
F	-3	TYR	-	expression tag	UNP Q5HCU0
F	-2	PHE	-	expression tag	UNP Q5HCU0
F	-1	GLN	-	expression tag	UNP Q5HCU0
F	0	GLY	-	expression tag	UNP Q5HCU0
F	234	SER	GLY	engineered mutation	UNP Q5HCU0
G	-20	MET	-	expression tag	UNP Q5HCU0
G	-19	GLY	-	expression tag	UNP Q5HCU0
G	-18	SER	-	expression tag	UNP Q5HCU0
G	-17	SER	-	expression tag	UNP Q5HCU0
G	-16	HIS	-	expression tag	UNP Q5HCU0
G	-15	HIS	-	expression tag	UNP Q5HCU0
G	-14	HIS	-	expression tag	UNP Q5HCU0

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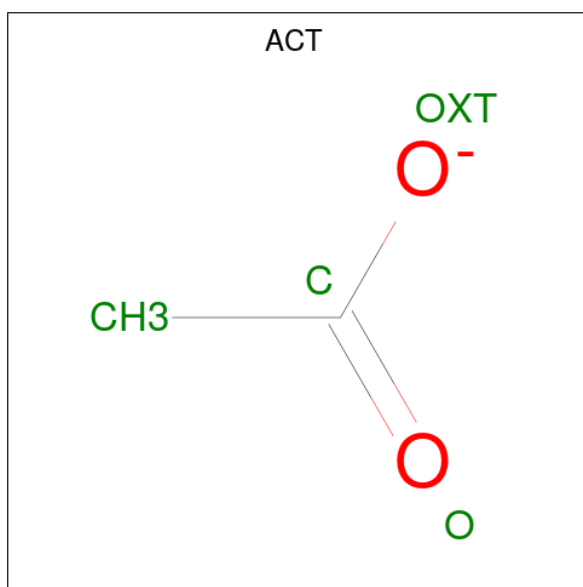
Chain	Residue	Modelled	Actual	Comment	Reference
G	-13	HIS	-	expression tag	UNP Q5HCU0
G	-12	HIS	-	expression tag	UNP Q5HCU0
G	-11	HIS	-	expression tag	UNP Q5HCU0
G	-10	SER	-	expression tag	UNP Q5HCU0
G	-9	SER	-	expression tag	UNP Q5HCU0
G	-8	GLY	-	expression tag	UNP Q5HCU0
G	-7	ARG	-	expression tag	UNP Q5HCU0
G	-6	GLU	-	expression tag	UNP Q5HCU0
G	-5	ASN	-	expression tag	UNP Q5HCU0
G	-4	LEU	-	expression tag	UNP Q5HCU0
G	-3	TYR	-	expression tag	UNP Q5HCU0
G	-2	PHE	-	expression tag	UNP Q5HCU0
G	-1	GLN	-	expression tag	UNP Q5HCU0
G	0	GLY	-	expression tag	UNP Q5HCU0
G	234	SER	GLY	engineered mutation	UNP Q5HCU0
H	-20	MET	-	expression tag	UNP Q5HCU0
H	-19	GLY	-	expression tag	UNP Q5HCU0
H	-18	SER	-	expression tag	UNP Q5HCU0
H	-17	SER	-	expression tag	UNP Q5HCU0
H	-16	HIS	-	expression tag	UNP Q5HCU0
H	-15	HIS	-	expression tag	UNP Q5HCU0
H	-14	HIS	-	expression tag	UNP Q5HCU0
H	-13	HIS	-	expression tag	UNP Q5HCU0
H	-12	HIS	-	expression tag	UNP Q5HCU0
H	-11	HIS	-	expression tag	UNP Q5HCU0
H	-10	SER	-	expression tag	UNP Q5HCU0
H	-9	SER	-	expression tag	UNP Q5HCU0
H	-8	GLY	-	expression tag	UNP Q5HCU0
H	-7	ARG	-	expression tag	UNP Q5HCU0
H	-6	GLU	-	expression tag	UNP Q5HCU0
H	-5	ASN	-	expression tag	UNP Q5HCU0
H	-4	LEU	-	expression tag	UNP Q5HCU0
H	-3	TYR	-	expression tag	UNP Q5HCU0
H	-2	PHE	-	expression tag	UNP Q5HCU0
H	-1	GLN	-	expression tag	UNP Q5HCU0
H	0	GLY	-	expression tag	UNP Q5HCU0
H	234	SER	GLY	engineered mutation	UNP Q5HCU0

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

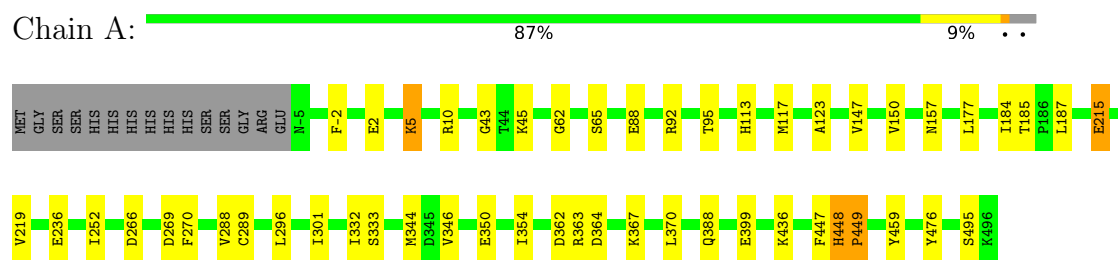
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	109	Total	O	0	0
			109	109		
4	B	115	Total	O	0	2
			116	116		
4	C	77	Total	O	0	1
			78	78		
4	D	108	Total	O	0	1
			109	109		
4	E	58	Total	O	0	0
			58	58		
4	F	90	Total	O	0	2
			92	92		
4	G	86	Total	O	0	0
			86	86		
4	H	59	Total	O	0	0
			59	59		

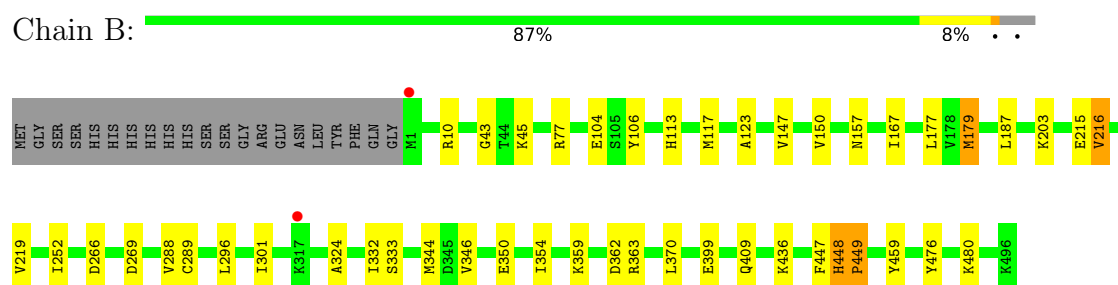
3 Residue-property plots

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.

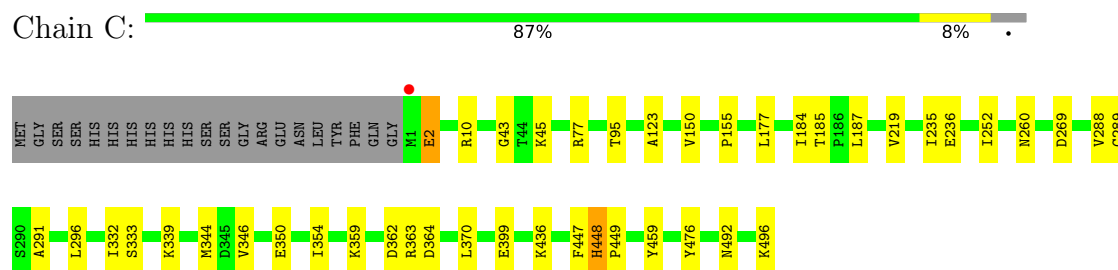
- Molecule 1: Betaine aldehyde dehydrogenase



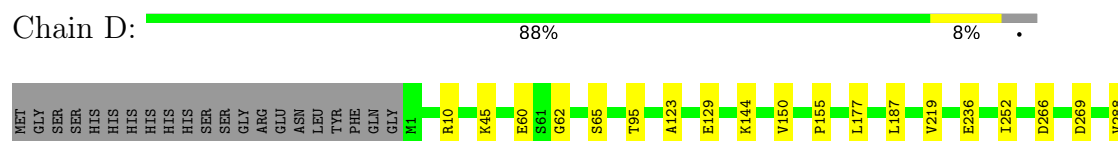
- Molecule 1: Betaine aldehyde dehydrogenase



- Molecule 1: Betaine aldehyde dehydrogenase



- Molecule 1: Betaine aldehyde dehydrogenase





• Molecule 1: Betaine aldehyde dehydrogenase

Chain E: 87% 9% . .



• Molecule 1: Betaine aldehyde dehydrogenase

Chain F: 88% 7% . .



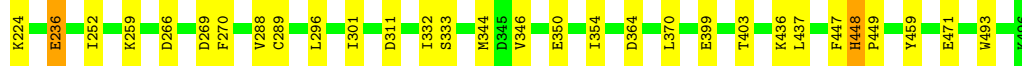
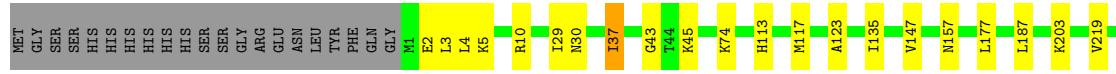
• Molecule 1: Betaine aldehyde dehydrogenase

Chain G: 88% 7% . .



• Molecule 1: Betaine aldehyde dehydrogenase

Chain H: 86% 9% . .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	88.59Å 168.37Å 144.52Å 90.00° 104.92° 90.00°	Depositor
Resolution (Å)	29.78 – 2.60 29.78 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.6 (29.78-2.60) 93.6 (29.78-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.184 , 0.219 0.184 , 0.217	Depositor DCC
R_{free} test set	5916 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.584	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 29.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	31934	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% 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, ACT

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	0.92	0/3972	0.99	14/5371 (0.3%)
1	B	0.83	0/3952	0.98	11/5344 (0.2%)
1	C	0.83	0/3918	0.99	13/5298 (0.2%)
1	D	0.86	0/3927	0.97	9/5310 (0.2%)
1	E	0.78	0/3918	0.98	13/5298 (0.2%)
1	F	0.86	1/3918 (0.0%)	0.98	7/5298 (0.1%)
1	G	0.84	0/3918	0.98	14/5298 (0.3%)
1	H	0.84	2/3918 (0.1%)	0.98	14/5298 (0.3%)
All	All	0.84	3/31441 (0.0%)	0.98	95/42515 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	492	ASN	CA-CB	5.25	1.61	1.53
1	H	135	ILE	N-CA	-5.25	1.41	1.46
1	H	29	ILE	C-O	-5.03	1.18	1.24

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	332	ILE	N-CA-C	10.49	121.32	110.62
1	C	332	ILE	N-CA-C	10.45	121.28	110.62
1	E	332	ILE	N-CA-C	10.29	121.12	110.62
1	B	332	ILE	N-CA-C	10.26	121.08	110.62
1	H	332	ILE	N-CA-C	10.24	121.07	110.62
1	D	332	ILE	N-CA-C	10.24	121.07	110.62
1	F	332	ILE	N-CA-C	10.20	121.02	110.62
1	G	332	ILE	N-CA-C	10.16	120.99	110.62
1	A	459	TYR	N-CA-C	-8.17	98.36	110.46
1	F	459	TYR	N-CA-C	-8.03	97.98	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	87	ARG	CG-CD-NE	7.95	129.49	112.00
1	C	459	TYR	N-CA-C	-7.90	98.18	110.42
1	G	212	ALA	N-CA-C	-7.84	94.37	108.48
1	B	179	MET	CA-CB-CG	7.82	129.74	114.10
1	D	459	TYR	N-CA-C	-7.80	98.32	110.42
1	B	459	TYR	N-CA-C	-7.76	98.39	110.42
1	E	459	TYR	N-CA-C	-7.64	98.57	110.42
1	G	459	TYR	N-CA-C	-7.62	98.60	110.42
1	H	459	TYR	N-CA-C	-7.53	98.75	110.42
1	G	438	LYS	CG-CD-CE	-6.86	95.53	111.30
1	A	449	PRO	N-CA-C	6.83	121.54	111.03
1	H	37	ILE	CB-CA-C	-6.73	104.05	110.65
1	E	180	LYS	CG-CD-CE	-6.63	96.05	111.30
1	H	187	LEU	N-CA-C	6.61	119.09	111.02
1	E	187	LEU	N-CA-C	6.47	118.92	111.02
1	G	187	LEU	N-CA-C	6.47	118.92	111.02
1	A	495	SER	N-CA-C	6.37	118.54	110.33
1	H	449	PRO	N-CA-C	6.36	120.83	111.03
1	F	449	PRO	N-CA-C	6.33	120.78	111.03
1	A	187	LEU	N-CA-C	6.32	118.73	111.02
1	C	187	LEU	N-CA-C	6.29	118.70	111.02
1	C	359	LYS	CA-CB-CG	6.29	126.68	114.10
1	B	216	VAL	N-CA-CB	-6.29	103.67	112.35
1	E	449	PRO	N-CA-C	6.25	120.65	111.03
1	G	449	PRO	N-CA-C	6.22	120.61	111.03
1	D	449	PRO	N-CA-C	6.21	120.59	111.03
1	C	449	PRO	N-CA-C	6.15	120.50	111.03
1	C	123	ALA	N-CA-C	6.12	118.73	111.33
1	G	123	ALA	N-CA-C	6.10	118.72	111.33
1	C	269	ASP	N-CA-C	-6.06	100.43	109.15
1	H	123	ALA	N-CA-C	5.97	118.56	111.33
1	F	187	LEU	N-CA-C	5.92	118.24	111.02
1	D	123	ALA	N-CA-C	5.92	118.49	111.33
1	F	269	ASP	N-CA-C	-5.91	100.64	109.15
1	D	269	ASP	N-CA-C	-5.91	100.64	109.15
1	E	10	ARG	CB-CG-CD	5.90	124.86	111.30
1	C	359	LYS	N-CA-CB	-5.88	101.84	111.66
1	B	10	ARG	N-CA-C	5.88	117.53	109.18
1	C	364	ASP	N-CA-C	5.86	118.17	111.02
1	D	187	LEU	N-CA-C	5.84	118.14	111.02
1	A	43	GLY	N-CA-C	5.83	118.75	112.33
1	E	259	LYS	N-CA-C	-5.83	98.87	108.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	ASP	N-CA-C	-5.74	100.89	109.15
1	E	123	ALA	N-CA-C	5.72	118.26	111.33
1	B	123	ALA	N-CA-C	5.70	118.23	111.33
1	H	269	ASP	N-CA-C	-5.70	100.95	109.15
1	B	269	ASP	N-CA-C	-5.68	100.96	109.15
1	G	438	LYS	CD-CE-NZ	5.60	129.82	111.90
1	F	123	ALA	N-CA-C	5.59	118.09	111.33
1	H	259	LYS	N-CA-C	-5.58	99.28	108.49
1	E	269	ASP	N-CA-C	-5.55	101.16	109.15
1	C	339	LYS	CB-CG-CD	5.53	124.03	111.30
1	A	5	LYS	CB-CA-C	5.51	120.93	109.95
1	D	364	ASP	N-CA-C	5.51	117.75	111.02
1	B	449	PRO	N-CA-C	5.49	119.71	111.14
1	A	123	ALA	N-CA-C	5.47	117.95	111.33
1	G	2	GLU	N-CA-C	-5.42	106.74	113.19
1	H	364	ASP	N-CA-C	5.41	117.61	111.02
1	G	269	ASP	N-CA-C	-5.38	101.41	109.15
1	G	157	ASN	N-CA-C	5.37	116.94	111.14
1	B	187	LEU	N-CA-C	5.33	117.52	111.02
1	H	157	ASN	N-CA-C	5.33	116.89	111.14
1	G	364	ASP	N-CA-C	5.31	117.48	111.11
1	F	364	ASP	N-CA-C	5.29	117.47	111.02
1	H	37	ILE	N-CA-CB	-5.27	105.01	112.12
1	C	235	ILE	N-CA-C	5.25	115.46	110.42
1	H	43	GLY	N-CA-C	5.22	118.07	112.33
1	A	10	ARG	N-CA-C	5.22	117.21	109.69
1	E	43	GLY	N-CA-C	5.20	118.04	112.33
1	A	270	PHE	N-CA-C	5.19	117.34	111.11
1	C	10	ARG	N-CA-C	5.16	117.12	109.69
1	H	10	ARG	N-CA-C	5.16	117.12	109.69
1	A	388	GLN	N-CA-C	5.15	118.72	112.23
1	D	10	ARG	N-CA-C	5.15	117.69	109.81
1	B	43	GLY	N-CA-C	5.13	117.97	112.33
1	C	43	GLY	N-CA-C	5.12	117.96	112.33
1	E	364	ASP	N-CA-C	5.11	117.24	111.11
1	G	43	GLY	N-CA-C	5.09	117.93	112.33
1	A	157	ASN	N-CA-C	5.07	116.61	111.14
1	D	299	ASN	N-CA-C	5.07	117.19	111.11
1	E	448	HIS	N-CA-C	5.07	121.00	109.81
1	A	364	ASP	N-CA-C	5.06	117.18	111.11
1	E	10	ARG	N-CA-C	5.06	116.97	109.69
1	B	157	ASN	N-CA-C	5.05	116.60	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	270	PHE	N-CA-C	5.04	117.15	111.11

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	3899	0	3849	26	0
1	B	3881	0	3819	25	0
1	C	3847	0	3803	18	0
1	D	3856	0	3808	22	0
1	E	3847	0	3803	27	0
1	F	3847	0	3803	24	0
1	G	3847	0	3803	21	0
1	H	3847	0	3803	27	0
2	A	44	0	26	5	0
2	B	44	0	26	2	0
2	C	44	0	26	3	0
2	D	44	0	26	7	0
2	E	44	0	26	9	0
2	F	44	0	26	5	0
2	G	44	0	26	5	0
2	H	44	0	26	2	0
3	B	4	0	3	0	0
4	A	109	0	0	2	0
4	B	116	0	0	4	0
4	C	78	0	0	1	0
4	D	109	0	0	2	0
4	E	58	0	0	1	0
4	F	92	0	0	4	0
4	G	86	0	0	3	0
4	H	59	0	0	5	0
All	All	31934	0	30702	182	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 (182) 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:289:CYS:SG	2:A:501:NAD:C4N	2.37	1.12
1:D:289:CYS:SG	2:D:501:NAD:C4N	2.40	1.08
1:E:289:CYS:SG	2:E:501:NAD:C4N	2.45	1.04
1:G:289:CYS:SG	2:G:501:NAD:C4N	2.46	1.04
1:F:289:CYS:SG	2:F:501:NAD:C4N	2.48	1.00
1:H:30:ASN:HB2	1:H:37:ILE:CD1	1.92	0.99
1:B:289:CYS:SG	2:B:501:NAD:C4N	2.52	0.97
1:C:289:CYS:SG	2:C:501:NAD:C4N	2.56	0.93
1:H:289:CYS:SG	2:H:501:NAD:C4N	2.60	0.89
1:H:30:ASN:HB2	1:H:37:ILE:HD12	1.54	0.89
1:B:147:VAL:HG21	4:B:618:HOH:O	1.76	0.83
1:C:496:LYS:OXT	1:E:315:LYS:HD3	1.80	0.82
1:G:359:LYS:HD3	4:G:680:HOH:O	1.83	0.78
1:H:30:ASN:HB2	1:H:37:ILE:HD11	1.64	0.78
1:A:147:VAL:HG21	4:A:700:HOH:O	1.86	0.76
1:F:184:ILE:HG22	1:F:185:THR:HG23	1.70	0.73
1:D:144:LYS:NZ	1:D:474:GLU:OE2	2.23	0.72
1:A:289:CYS:SG	2:A:501:NAD:C5N	2.78	0.72
4:C:677:HOH:O	1:E:431:GLN:HG2	1.90	0.72
1:D:496:LYS:NZ	4:D:655:HOH:O	2.23	0.72
1:D:129:GLU:OE1	1:G:470:LYS:NZ	2.23	0.71
1:C:2:GLU:N	1:C:2:GLU:OE1	2.23	0.71
1:A:184:ILE:HG22	1:A:185:THR:HG23	1.73	0.70
1:A:215:GLU:HG3	4:A:643:HOH:O	1.91	0.69
1:G:311:ASP:HB3	4:G:636:HOH:O	1.92	0.68
1:A:2:GLU:OE1	1:A:5:LYS:NZ	2.26	0.68
1:A:289:CYS:SG	2:A:501:NAD:H4N	2.33	0.68
1:C:184:ILE:HG22	1:C:185:THR:HG23	1.75	0.67
1:H:403:THR:HB	4:H:624:HOH:O	1.95	0.67
1:B:289:CYS:SG	2:B:501:NAD:C3N	2.83	0.66
1:E:346:VAL:O	1:E:350:GLU:HG3	1.96	0.66
1:E:289:CYS:SG	2:E:501:NAD:C3N	2.84	0.66
1:B:346:VAL:O	1:B:350:GLU:HG3	1.95	0.66
1:C:346:VAL:O	1:C:350:GLU:HG3	1.97	0.65
1:A:346:VAL:O	1:A:350:GLU:HG3	1.97	0.65
1:H:346:VAL:O	1:H:350:GLU:HG3	1.97	0.65
1:D:346:VAL:O	1:D:350:GLU:HG3	1.96	0.64
1:H:147:VAL:HG21	4:H:651:HOH:O	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:346:VAL:O	1:G:350:GLU:HG3	1.98	0.63
1:F:346:VAL:O	1:F:350:GLU:HG3	1.97	0.63
1:F:364:ASP:HB3	4:F:668:HOH:O	1.99	0.63
1:C:492:ASN:OD1	1:C:496:LYS:NZ	2.32	0.62
1:D:471:GLU:HG3	4:D:613:HOH:O	2.01	0.61
1:D:289:CYS:SG	2:D:501:NAD:C5N	2.89	0.61
1:C:155:PRO:HB3	2:C:501:NAD:H52N	1.83	0.60
1:G:289:CYS:SG	2:G:501:NAD:H4N	2.41	0.60
1:C:289:CYS:SG	2:C:501:NAD:C3N	2.90	0.59
1:D:289:CYS:SG	2:D:501:NAD:C3N	2.90	0.58
1:F:289:CYS:SG	2:F:501:NAD:C3N	2.90	0.58
1:E:88:GLU:OE1	1:E:92:ARG:NH2	2.38	0.57
1:B:324:ALA:HB3	1:F:490:LEU:CD2	2.35	0.57
1:B:359:LYS:HD3	4:B:674:HOH:O	2.06	0.56
1:D:289:CYS:SG	2:D:501:NAD:H4N	2.40	0.55
1:G:289:CYS:SG	2:G:501:NAD:C3N	2.95	0.55
1:F:100:LYS:NZ	4:F:680:HOH:O	2.27	0.54
1:E:45:LYS:HB2	1:E:219:VAL:HG21	1.89	0.54
1:B:203:LYS:HE3	4:B:690:HOH:O	2.06	0.54
1:F:45:LYS:HB2	1:F:219:VAL:HG21	1.91	0.53
1:B:106:TYR:OH	1:F:496:LYS:OXT	2.24	0.53
1:H:45:LYS:HB2	1:H:219:VAL:HG21	1.91	0.53
1:H:203:LYS:HA	4:H:639:HOH:O	2.09	0.52
1:G:183:GLU:HB3	1:G:211:GLY:O	2.08	0.52
1:F:335:GLU:HG3	1:F:336:HIS:N	2.25	0.52
1:A:289:CYS:HB3	2:A:501:NAD:C6N	2.40	0.52
1:E:289:CYS:SG	2:E:501:NAD:C5N	2.97	0.51
1:C:45:LYS:HB2	1:C:219:VAL:HG21	1.92	0.51
1:G:289:CYS:SG	2:G:501:NAD:C5N	3.00	0.50
1:B:45:LYS:HB2	1:B:219:VAL:HG21	1.92	0.50
1:A:45:LYS:HB2	1:A:219:VAL:HG21	1.92	0.50
1:G:45:LYS:HB2	1:G:219:VAL:HG21	1.93	0.50
1:D:45:LYS:HB2	1:D:219:VAL:HG21	1.93	0.50
1:A:289:CYS:SG	2:A:501:NAD:C3N	3.00	0.50
1:E:257:GLY:HA2	2:E:501:NAD:O2D	2.12	0.49
1:C:447:PHE:O	1:C:448:HIS:HB2	2.13	0.49
1:D:60:GLU:OE2	1:G:438:LYS:NZ	2.46	0.49
1:A:447:PHE:O	1:A:448:HIS:HB2	2.13	0.48
1:B:447:PHE:O	1:B:448:HIS:HB2	2.12	0.48
1:H:4:LEU:CD1	1:H:37:ILE:CG2	2.92	0.48
1:E:447:PHE:O	1:E:448:HIS:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-2:PHE:O	1:A:2:GLU:HG2	2.14	0.48
1:B:77:ARG:NH1	1:C:77:ARG:HG2	2.28	0.48
1:F:289:CYS:SG	2:F:501:NAD:C5N	3.00	0.48
1:A:2:GLU:OE1	1:A:5:LYS:HE2	2.13	0.48
1:C:296:LEU:HD23	1:C:399:GLU:HB2	1.96	0.48
1:E:344:MET:CE	1:E:354:ILE:HD12	2.44	0.48
1:E:217:GLY:HA3	2:E:501:NAD:C2A	2.43	0.47
1:H:4:LEU:CD1	1:H:37:ILE:HG23	2.45	0.47
1:H:4:LEU:HD11	1:H:37:ILE:CG2	2.44	0.47
1:H:447:PHE:O	1:H:448:HIS:HB2	2.14	0.47
1:A:2:GLU:OE1	1:A:5:LYS:CE	2.62	0.47
1:A:344:MET:CE	1:A:354:ILE:HD12	2.44	0.47
1:C:344:MET:CE	1:C:354:ILE:HD12	2.44	0.47
1:H:289:CYS:SG	2:H:501:NAD:C3N	3.02	0.47
1:H:344:MET:CE	1:H:354:ILE:HD12	2.44	0.47
1:D:344:MET:CE	1:D:354:ILE:HD12	2.44	0.47
1:F:289:CYS:SG	2:F:501:NAD:H4N	2.51	0.47
1:F:344:MET:CE	1:F:354:ILE:HD12	2.45	0.47
1:G:447:PHE:O	1:G:448:HIS:HB2	2.15	0.47
1:A:296:LEU:HD23	1:A:399:GLU:HB2	1.97	0.47
1:G:344:MET:CE	1:G:354:ILE:HD12	2.45	0.47
1:H:37:ILE:CG2	1:H:37:ILE:O	2.59	0.46
1:F:150:VAL:HG11	1:F:476:TYR:HE1	1.80	0.46
1:H:296:LEU:HD23	1:H:399:GLU:HB2	1.97	0.46
1:D:447:PHE:O	1:D:448:HIS:HB2	2.16	0.46
1:B:344:MET:CE	1:B:354:ILE:HD12	2.45	0.46
1:B:104:GLU:HG3	1:H:493:TRP:CH2	2.49	0.46
1:D:296:LEU:HD23	1:D:399:GLU:HB2	1.98	0.46
1:B:296:LEU:HD23	1:B:399:GLU:HB2	1.98	0.46
1:D:150:VAL:HG11	1:D:476:TYR:HE1	1.80	0.46
1:H:74:LYS:HE2	4:H:621:HOH:O	2.16	0.46
1:A:150:VAL:HG11	1:A:476:TYR:HE1	1.81	0.45
1:F:447:PHE:O	1:F:448:HIS:HB2	2.14	0.45
1:E:289:CYS:SG	2:E:501:NAD:H4N	2.50	0.45
1:B:409:GLN:HB3	4:B:703:HOH:O	2.17	0.45
1:C:150:VAL:HG11	1:C:476:TYR:HE1	1.81	0.45
1:E:180:LYS:NZ	2:E:501:NAD:O2B	2.50	0.44
1:D:155:PRO:HB3	2:D:501:NAD:H52N	1.99	0.44
1:E:296:LEU:HD23	1:E:399:GLU:HB2	1.98	0.44
1:H:113:HIS:CE1	1:H:117:MET:CE	3.01	0.44
1:B:150:VAL:HG11	1:B:476:TYR:HE2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:296:LEU:HD23	1:F:399:GLU:HB2	1.99	0.44
1:G:296:LEU:HD23	1:G:399:GLU:HB2	1.98	0.44
1:E:177:LEU:C	1:E:177:LEU:HD12	2.42	0.44
1:B:177:LEU:C	1:B:177:LEU:HD12	2.43	0.44
1:F:74:LYS:HE2	4:F:651:HOH:O	2.17	0.43
1:F:183:GLU:HG2	1:F:184:ILE:HD13	2.00	0.43
1:G:113:HIS:CE1	1:G:117:MET:CE	3.01	0.43
1:A:177:LEU:C	1:A:177:LEU:HD12	2.42	0.43
1:E:20:SER:N	1:E:41:SER:OG	2.52	0.43
1:C:177:LEU:HD12	1:C:177:LEU:C	2.43	0.43
1:E:150:VAL:HG11	1:E:476:TYR:HE2	1.83	0.43
1:D:289:CYS:HB3	2:D:501:NAD:C6N	2.49	0.43
1:D:177:LEU:C	1:D:177:LEU:HD12	2.43	0.42
1:E:2:GLU:O	1:E:5:LYS:HG2	2.19	0.42
1:A:88:GLU:OE1	1:A:92:ARG:NH1	2.53	0.42
1:G:150:VAL:HG11	1:G:476:TYR:HE2	1.83	0.42
1:H:177:LEU:C	1:H:177:LEU:HD12	2.44	0.42
1:A:333:SER:HA	1:A:370:LEU:HD13	2.01	0.42
1:B:167:ILE:HD11	1:B:179:MET:HE3	2.01	0.42
1:G:177:LEU:HD12	1:G:177:LEU:C	2.44	0.42
1:F:448:HIS:N	1:F:449:PRO:CD	2.83	0.42
1:A:266:ASP:HA	1:A:301:ILE:HD13	2.02	0.42
1:B:113:HIS:CE1	1:B:117:MET:CE	3.03	0.42
1:B:333:SER:HA	1:B:370:LEU:HD13	2.02	0.42
1:F:98:THR:HG21	4:F:680:HOH:O	2.20	0.42
1:G:470:LYS:HE3	4:G:659:HOH:O	2.20	0.42
1:D:155:PRO:CB	2:D:501:NAD:H52N	2.50	0.41
1:E:471:GLU:HG3	4:E:627:HOH:O	2.20	0.41
1:F:113:HIS:CE1	1:F:117:MET:CE	3.03	0.41
1:H:471:GLU:HG3	4:H:637:HOH:O	2.19	0.41
1:B:480:LYS:NZ	1:H:437:LEU:O	2.53	0.41
1:E:260:ASN:HB2	1:E:291:ALA:O	2.20	0.41
1:G:448:HIS:N	1:G:449:PRO:CD	2.83	0.41
1:A:448:HIS:N	1:A:449:PRO:CD	2.83	0.41
1:G:392:PHE:CE1	2:G:501:NAD:H2D	2.55	0.41
1:D:266:ASP:HA	1:D:301:ILE:HD13	2.03	0.41
1:A:362:ASP:O	1:A:363:ARG:C	2.64	0.41
1:D:333:SER:HA	1:D:370:LEU:HD13	2.01	0.41
1:H:236:GLU:H	1:H:236:GLU:CD	2.28	0.41
1:E:333:SER:HA	1:E:370:LEU:HD13	2.02	0.41
1:E:448:HIS:N	1:E:449:PRO:CD	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:177:LEU:C	1:F:177:LEU:HD12	2.46	0.41
1:F:392:PHE:CE1	2:F:501:NAD:H2D	2.56	0.41
1:G:447:PHE:C	1:G:449:PRO:HD3	2.45	0.41
1:C:362:ASP:O	1:C:363:ARG:C	2.64	0.41
1:H:333:SER:HA	1:H:370:LEU:HD13	2.03	0.41
1:A:62:GLY:HA2	1:A:65:SER:OG	2.21	0.40
1:A:113:HIS:CE1	1:A:117:MET:CE	3.04	0.40
1:B:266:ASP:HA	1:B:301:ILE:HD13	2.03	0.40
1:B:362:ASP:O	1:B:363:ARG:C	2.64	0.40
1:C:260:ASN:HB2	1:C:291:ALA:O	2.21	0.40
1:E:266:ASP:HA	1:E:301:ILE:HD13	2.03	0.40
1:H:2:GLU:O	1:H:5:LYS:HG2	2.21	0.40
1:C:333:SER:HA	1:C:370:LEU:HD13	2.03	0.40
1:E:447:PHE:C	1:E:449:PRO:HD3	2.47	0.40
1:B:215:GLU:HG3	1:B:216:VAL:HG23	2.03	0.40
1:E:289:CYS:HB3	2:E:501:NAD:C6N	2.51	0.40
1:H:266:ASP:HA	1:H:301:ILE:HD13	2.03	0.40
1:B:447:PHE:C	1:B:449:PRO:HD3	2.47	0.40
1:D:62:GLY:HA2	1:D:65:SER:OG	2.22	0.40
1:E:217:GLY:HA3	2:E:501:NAD:N1A	2.36	0.40
1:F:266:ASP:HA	1:F:301:ILE:HD13	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	500/517 (97%)	485 (97%)	14 (3%)	1 (0%)	43	66
1	B	498/517 (96%)	486 (98%)	11 (2%)	1 (0%)	43	66
1	C	494/517 (96%)	480 (97%)	13 (3%)	1 (0%)	43	66
1	D	495/517 (96%)	483 (98%)	11 (2%)	1 (0%)	43	66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	494/517 (96%)	480 (97%)	13 (3%)	1 (0%)	43	66
1	F	494/517 (96%)	480 (97%)	13 (3%)	1 (0%)	43	66
1	G	494/517 (96%)	481 (97%)	12 (2%)	1 (0%)	43	66
1	H	494/517 (96%)	481 (97%)	12 (2%)	1 (0%)	43	66
All	All	3963/4136 (96%)	3856 (97%)	99 (2%)	8 (0%)	43	66

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	448	HIS
1	B	448	HIS
1	C	448	HIS
1	D	448	HIS
1	E	448	HIS
1	F	448	HIS
1	G	448	HIS
1	H	448	HIS

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	417/430 (97%)	410 (98%)	7 (2%)	53	78
1	B	416/430 (97%)	413 (99%)	3 (1%)	76	89
1	C	412/430 (96%)	406 (98%)	6 (2%)	57	80
1	D	413/430 (96%)	408 (99%)	5 (1%)	63	83
1	E	412/430 (96%)	405 (98%)	7 (2%)	53	78
1	F	412/430 (96%)	406 (98%)	6 (2%)	57	80
1	G	412/430 (96%)	408 (99%)	4 (1%)	68	86
1	H	412/430 (96%)	405 (98%)	7 (2%)	53	78
All	All	3306/3440 (96%)	3261 (99%)	45 (1%)	59	81

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	THR
1	A	215	GLU
1	A	236	GLU
1	A	252	ILE
1	A	288	VAL
1	A	367	LYS
1	A	436	LYS
1	B	252	ILE
1	B	288	VAL
1	B	436	LYS
1	C	2	GLU
1	C	95	THR
1	C	236	GLU
1	C	252	ILE
1	C	288	VAL
1	C	436	LYS
1	D	95	THR
1	D	236	GLU
1	D	252	ILE
1	D	288	VAL
1	D	436	LYS
1	E	10	ARG
1	E	95	THR
1	E	224	LYS
1	E	252	ILE
1	E	288	VAL
1	E	409	GLN
1	E	436	LYS
1	F	63	GLU
1	F	236	GLU
1	F	252	ILE
1	F	288	VAL
1	F	335	GLU
1	F	409	GLN
1	G	77	ARG
1	G	252	ILE
1	G	288	VAL
1	G	436	LYS
1	H	3	LEU
1	H	224	LYS
1	H	236	GLU
1	H	252	ILE

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Mol	Chain	Res	Type
1	H	288	VAL
1	H	311	ASP
1	H	436	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	492	ASN
1	B	9	GLN
1	B	34	GLN
1	C	34	GLN
1	C	113	HIS
1	C	260	ASN
1	C	431	GLN
1	D	34	GLN
1	D	405	GLN
1	D	409	GLN
1	E	431	GLN
1	E	492	ASN
1	F	34	GLN
1	G	34	GLN
1	H	34	GLN
1	H	409	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

9 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	G	501	-	46,48,48	1.15	6 (13%)	64,73,73	1.76	12 (18%)
2	NAD	C	501	-	46,48,48	1.38	5 (10%)	64,73,73	1.82	12 (18%)
2	NAD	H	501	-	46,48,48	1.33	6 (13%)	64,73,73	1.64	12 (18%)
2	NAD	B	501	-	46,48,48	1.39	4 (8%)	64,73,73	1.77	10 (15%)
3	ACT	B	502	-	3,3,3	0.84	0	3,3,3	0.76	0
2	NAD	F	501	-	46,48,48	1.09	2 (4%)	64,73,73	1.73	14 (21%)
2	NAD	D	501	-	46,48,48	1.30	6 (13%)	64,73,73	1.67	12 (18%)
2	NAD	A	501	-	46,48,48	1.30	3 (6%)	64,73,73	1.67	11 (17%)
2	NAD	E	501	-	46,48,48	1.52	5 (10%)	64,73,73	1.62	10 (15%)

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	G	501	-	-	6/30/62/62	0/5/5/5
2	NAD	C	501	-	-	7/30/62/62	0/5/5/5
2	NAD	H	501	-	-	6/30/62/62	0/5/5/5
2	NAD	B	501	-	-	8/30/62/62	0/5/5/5
2	NAD	F	501	-	-	5/30/62/62	0/5/5/5
2	NAD	D	501	-	-	8/30/62/62	0/5/5/5
2	NAD	A	501	-	-	8/30/62/62	0/5/5/5
2	NAD	E	501	-	-	8/30/62/62	0/5/5/5

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	501	NAD	PN-O3	6.27	1.66	1.59
2	H	501	NAD	O7N-C7N	5.52	1.34	1.24
2	A	501	NAD	O7N-C7N	5.46	1.34	1.24
2	C	501	NAD	O7N-C7N	4.91	1.33	1.24
2	E	501	NAD	O7N-C7N	4.89	1.33	1.24
2	B	501	NAD	O7N-C7N	4.88	1.33	1.24
2	C	501	NAD	PN-O3	4.71	1.64	1.59
2	B	501	NAD	PN-O3	4.69	1.64	1.59
2	G	501	NAD	O7N-C7N	4.30	1.32	1.24
2	A	501	NAD	PN-O3	3.88	1.63	1.59
2	D	501	NAD	O7N-C7N	3.74	1.31	1.24
2	F	501	NAD	O7N-C7N	3.69	1.31	1.24
2	D	501	NAD	PA-O3	3.52	1.63	1.59
2	D	501	NAD	PN-O3	3.23	1.63	1.59
2	E	501	NAD	C4A-N9A	-3.23	1.31	1.37
2	B	501	NAD	PA-O3	3.23	1.63	1.59
2	H	501	NAD	PN-O3	3.08	1.62	1.59
2	H	501	NAD	O4D-C1D	2.97	1.44	1.40
2	G	501	NAD	C8A-N9A	-2.61	1.33	1.37
2	D	501	NAD	C8A-N7A	2.56	1.36	1.31
2	C	501	NAD	PA-O3	2.55	1.62	1.59
2	C	501	NAD	C5A-N7A	-2.50	1.34	1.39
2	F	501	NAD	C8A-N7A	2.47	1.36	1.31
2	H	501	NAD	C4A-N9A	-2.45	1.32	1.37
2	H	501	NAD	C8A-N7A	2.45	1.36	1.31
2	H	501	NAD	PA-O3	2.45	1.62	1.59
2	C	501	NAD	C8A-N9A	-2.43	1.33	1.37
2	E	501	NAD	C5A-N7A	-2.42	1.34	1.39
2	A	501	NAD	C8A-N7A	2.41	1.36	1.31
2	G	501	NAD	O4D-C1D	2.32	1.43	1.40
2	G	501	NAD	C4A-N9A	-2.31	1.32	1.37
2	D	501	NAD	C4A-N9A	-2.31	1.32	1.37
2	D	501	NAD	C5A-N7A	-2.19	1.35	1.39
2	G	501	NAD	C8A-N7A	2.10	1.35	1.31
2	G	501	NAD	C5A-N7A	-2.08	1.35	1.39
2	B	501	NAD	C4A-N9A	-2.07	1.33	1.37
2	E	501	NAD	C8A-N7A	2.03	1.35	1.31

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	C5A-C4A-N3A	-6.21	118.17	126.72
2	C	501	NAD	C5A-C4A-N3A	-5.85	118.66	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	C5A-C4A-N3A	-5.44	119.22	126.72
2	E	501	NAD	C5A-C4A-N3A	-5.35	119.34	126.72
2	B	501	NAD	N3A-C2A-N1A	-5.28	120.59	128.58
2	H	501	NAD	N3A-C2A-N1A	-5.28	120.59	128.58
2	C	501	NAD	N3A-C2A-N1A	-5.19	120.73	128.58
2	F	501	NAD	C5A-C4A-N3A	-5.01	119.81	126.72
2	F	501	NAD	N3A-C2A-N1A	-5.00	121.02	128.58
2	G	501	NAD	C5A-C4A-N3A	-4.81	120.10	126.72
2	G	501	NAD	N3A-C2A-N1A	-4.80	121.31	128.58
2	A	501	NAD	N3A-C2A-N1A	-4.68	121.50	128.58
2	D	501	NAD	N3A-C2A-N1A	-4.65	121.54	128.58
2	C	501	NAD	N9A-C8A-N7A	-4.58	107.43	113.94
2	H	501	NAD	C5A-C4A-N3A	-4.46	120.58	126.72
2	H	501	NAD	N9A-C8A-N7A	-4.42	107.66	113.94
2	D	501	NAD	C5A-C4A-N3A	-4.42	120.63	126.72
2	B	501	NAD	C2A-N3A-C4A	4.30	122.33	111.83
2	G	501	NAD	N9A-C8A-N7A	-4.15	108.06	113.94
2	A	501	NAD	N9A-C8A-N7A	-4.12	108.10	113.94
2	D	501	NAD	C3N-C7N-N7N	4.08	122.77	117.74
2	F	501	NAD	N9A-C8A-N7A	-4.07	108.17	113.94
2	G	501	NAD	C4A-C5A-N7A	-4.02	105.98	110.58
2	E	501	NAD	N9A-C8A-N7A	-4.01	108.24	113.94
2	C	501	NAD	N3A-C4A-N9A	3.99	133.96	127.17
2	G	501	NAD	C3N-C7N-N7N	3.88	122.52	117.74
2	C	501	NAD	C2A-N3A-C4A	3.80	121.10	111.83
2	B	501	NAD	C4A-C5A-N7A	-3.79	106.25	110.58
2	B	501	NAD	O4B-C1B-N9A	3.78	115.36	108.09
2	F	501	NAD	C2A-N3A-C4A	3.73	120.94	111.83
2	E	501	NAD	N3A-C2A-N1A	-3.73	122.94	128.58
2	A	501	NAD	N3A-C4A-N9A	3.65	133.37	127.17
2	H	501	NAD	C2A-N3A-C4A	3.64	120.72	111.83
2	E	501	NAD	C4A-C5A-N7A	-3.59	106.48	110.58
2	A	501	NAD	C4A-N9A-C8A	3.55	109.47	105.74
2	A	501	NAD	C2A-N3A-C4A	3.53	120.46	111.83
2	B	501	NAD	N9A-C8A-N7A	-3.51	108.96	113.94
2	D	501	NAD	N9A-C8A-N7A	-3.49	108.99	113.94
2	C	501	NAD	C5A-N7A-C8A	3.40	108.79	103.45
2	B	501	NAD	N3A-C4A-N9A	3.39	132.93	127.17
2	H	501	NAD	C4A-N9A-C8A	3.39	109.29	105.74
2	G	501	NAD	C2A-N3A-C4A	3.37	120.06	111.83
2	D	501	NAD	C2A-N3A-C4A	3.36	120.04	111.83
2	H	501	NAD	C4A-C5A-N7A	-3.35	106.75	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	501	NAD	C5A-N7A-C8A	3.24	108.54	103.45
2	C	501	NAD	C4A-N9A-C8A	3.22	109.12	105.74
2	E	501	NAD	C2A-N3A-C4A	3.22	119.69	111.83
2	G	501	NAD	O7N-C7N-N7N	-3.17	118.04	122.62
2	G	501	NAD	C4A-N9A-C8A	3.16	109.05	105.74
2	C	501	NAD	O7N-C7N-N7N	-3.14	118.08	122.62
2	B	501	NAD	C5A-N7A-C8A	3.13	108.38	103.45
2	H	501	NAD	O4B-C1B-N9A	3.10	114.04	108.09
2	F	501	NAD	C4A-C5A-N7A	-3.01	107.14	110.58
2	C	501	NAD	O5D-PN-O1N	-3.00	97.06	108.94
2	D	501	NAD	O7N-C7N-N7N	-2.97	118.32	122.62
2	C	501	NAD	C4A-C5A-N7A	-2.90	107.27	110.58
2	F	501	NAD	O7N-C7N-N7N	-2.88	118.45	122.62
2	D	501	NAD	N3A-C4A-N9A	2.87	132.04	127.17
2	E	501	NAD	C5A-N7A-C8A	2.86	107.94	103.45
2	B	501	NAD	C5A-C4A-N9A	2.84	108.91	105.81
2	D	501	NAD	C4A-N9A-C8A	2.83	108.71	105.74
2	G	501	NAD	O2N-PN-O3	2.83	114.92	107.27
2	H	501	NAD	C5A-N7A-C8A	2.80	107.85	103.45
2	E	501	NAD	N3A-C4A-N9A	2.76	131.86	127.17
2	E	501	NAD	C5A-C4A-N9A	2.74	108.80	105.81
2	A	501	NAD	O5D-PN-O1N	-2.74	98.09	108.94
2	F	501	NAD	C5A-N7A-C8A	2.73	107.73	103.45
2	F	501	NAD	N3A-C4A-N9A	2.68	131.73	127.17
2	D	501	NAD	C4A-N9A-C1B	-2.68	120.37	126.63
2	E	501	NAD	C4A-N9A-C8A	2.67	108.54	105.74
2	F	501	NAD	C4A-N9A-C8A	2.62	108.49	105.74
2	G	501	NAD	N3A-C4A-N9A	2.56	131.53	127.17
2	A	501	NAD	C4A-C5A-N7A	-2.54	107.68	110.58
2	A	501	NAD	C5A-N7A-C8A	2.49	107.36	103.45
2	C	501	NAD	O7N-C7N-C3N	2.48	122.62	119.60
2	F	501	NAD	C5A-C4A-N9A	2.43	108.46	105.81
2	H	501	NAD	C5A-C4A-N9A	2.41	108.44	105.81
2	F	501	NAD	O7N-C7N-C3N	2.37	122.49	119.60
2	D	501	NAD	C5A-C6A-N6A	-2.36	117.43	123.29
2	G	501	NAD	C5A-C4A-N9A	2.34	108.36	105.81
2	F	501	NAD	O2N-PN-O3	2.34	113.58	107.27
2	D	501	NAD	O5D-PN-O1N	-2.25	100.01	108.94
2	E	501	NAD	O2N-PN-O3	2.25	113.36	107.27
2	H	501	NAD	N3A-C4A-N9A	2.23	130.97	127.17
2	H	501	NAD	O7N-C7N-N7N	-2.23	119.40	122.62
2	D	501	NAD	O2N-PN-O3	2.15	113.07	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	NAD	C4D-O4D-C1D	2.13	111.88	109.92
2	F	501	NAD	C4D-O4D-C1D	2.13	111.88	109.92
2	H	501	NAD	C4A-N9A-C1B	-2.13	121.65	126.63
2	A	501	NAD	O4D-C4D-C5D	-2.10	102.61	109.33
2	B	501	NAD	O7N-C7N-N7N	-2.08	119.61	122.62
2	F	501	NAD	C4A-N9A-C1B	-2.04	121.86	126.63
2	A	501	NAD	C5A-C6A-N6A	-2.04	118.24	123.29

There are no chirality outliers.

All (56) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAD	O4D-C1D-N1N-C6N
2	B	501	NAD	O4D-C1D-N1N-C6N
2	C	501	NAD	O4D-C1D-N1N-C6N
2	D	501	NAD	O4D-C1D-N1N-C6N
2	E	501	NAD	O4D-C1D-N1N-C6N
2	G	501	NAD	O4D-C1D-N1N-C6N
2	H	501	NAD	C5B-O5B-PA-O3
2	G	501	NAD	O4D-C4D-C5D-O5D
2	H	501	NAD	O4D-C4D-C5D-O5D
2	A	501	NAD	C3D-C4D-C5D-O5D
2	F	501	NAD	C3D-C4D-C5D-O5D
2	G	501	NAD	C3D-C4D-C5D-O5D
2	H	501	NAD	C3D-C4D-C5D-O5D
2	A	501	NAD	O4D-C4D-C5D-O5D
2	B	501	NAD	O4D-C4D-C5D-O5D
2	D	501	NAD	O4D-C4D-C5D-O5D
2	D	501	NAD	C3D-C4D-C5D-O5D
2	E	501	NAD	O4D-C4D-C5D-O5D
2	F	501	NAD	O4D-C4D-C5D-O5D
2	B	501	NAD	C3D-C4D-C5D-O5D
2	E	501	NAD	C3D-C4D-C5D-O5D
2	B	501	NAD	C2B-C1B-N9A-C8A
2	C	501	NAD	C4D-C5D-O5D-PN
2	A	501	NAD	C2B-C1B-N9A-C8A
2	C	501	NAD	C2B-C1B-N9A-C8A
2	D	501	NAD	C2B-C1B-N9A-C8A
2	E	501	NAD	C2B-C1B-N9A-C8A
2	F	501	NAD	C2B-C1B-N9A-C8A
2	G	501	NAD	C2B-C1B-N9A-C8A
2	G	501	NAD	C4D-C5D-O5D-PN

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Mol	Chain	Res	Type	Atoms
2	D	501	NAD	C4D-C5D-O5D-PN
2	E	501	NAD	C4D-C5D-O5D-PN
2	H	501	NAD	C4D-C5D-O5D-PN
2	A	501	NAD	C4D-C5D-O5D-PN
2	F	501	NAD	C4D-C5D-O5D-PN
2	D	501	NAD	C2B-C1B-N9A-C4A
2	H	501	NAD	C2B-C1B-N9A-C8A
2	B	501	NAD	C4D-C5D-O5D-PN
2	A	501	NAD	O4D-C1D-N1N-C2N
2	B	501	NAD	O4D-C1D-N1N-C2N
2	C	501	NAD	O4D-C1D-N1N-C2N
2	D	501	NAD	O4D-C1D-N1N-C2N
2	E	501	NAD	O4D-C1D-N1N-C2N
2	G	501	NAD	O4D-C1D-N1N-C2N
2	A	501	NAD	O4B-C1B-N9A-C8A
2	A	501	NAD	C2B-C1B-N9A-C4A
2	B	501	NAD	O4B-C1B-N9A-C8A
2	C	501	NAD	O4B-C1B-N9A-C8A
2	D	501	NAD	O4B-C1B-N9A-C8A
2	H	501	NAD	O4B-C1B-N9A-C8A
2	E	501	NAD	O4B-C1B-N9A-C8A
2	B	501	NAD	C2B-C1B-N9A-C4A
2	C	501	NAD	C2B-C1B-N9A-C4A
2	E	501	NAD	C2B-C1B-N9A-C4A
2	F	501	NAD	O4B-C1B-N9A-C8A
2	C	501	NAD	O4D-C4D-C5D-O5D

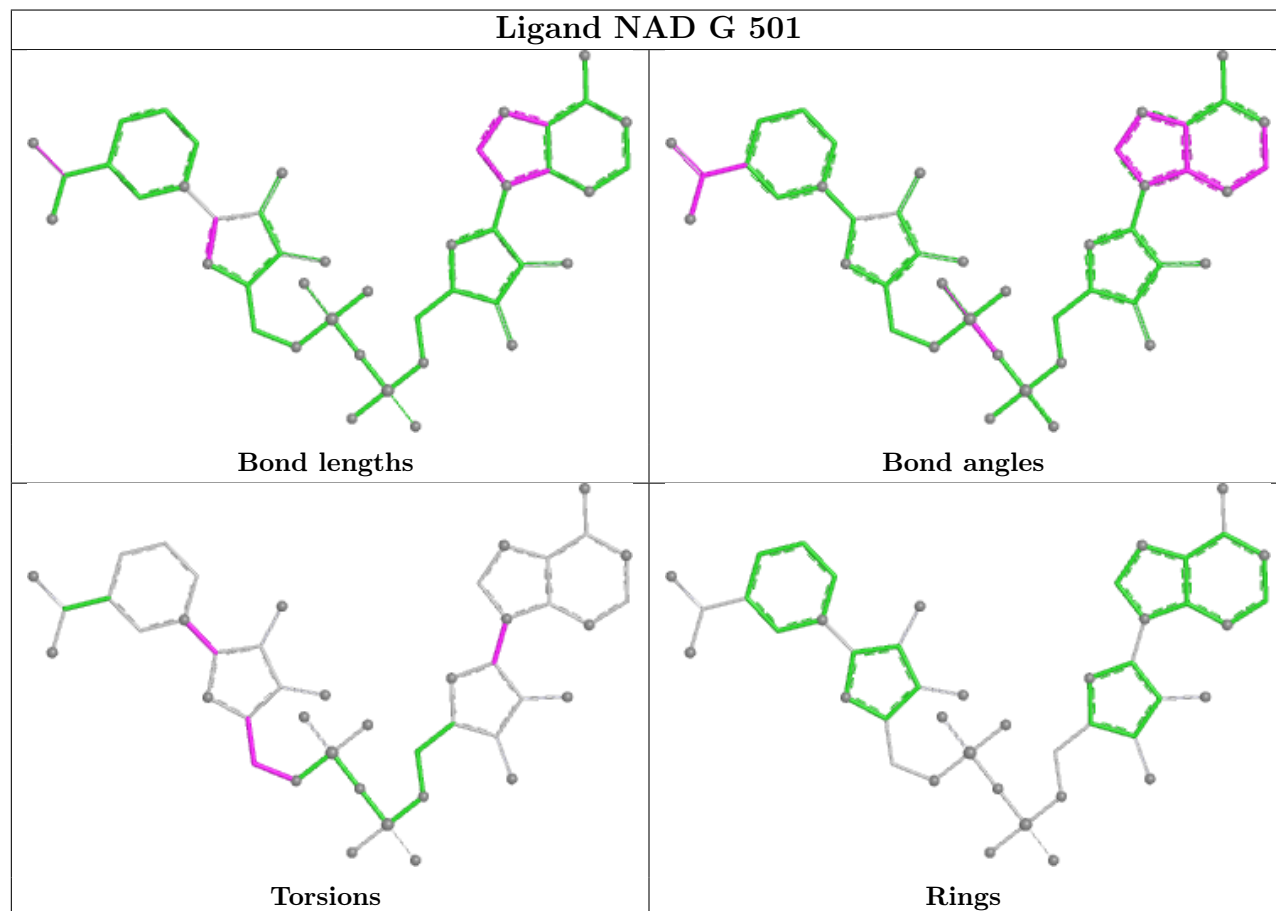
There are no ring outliers.

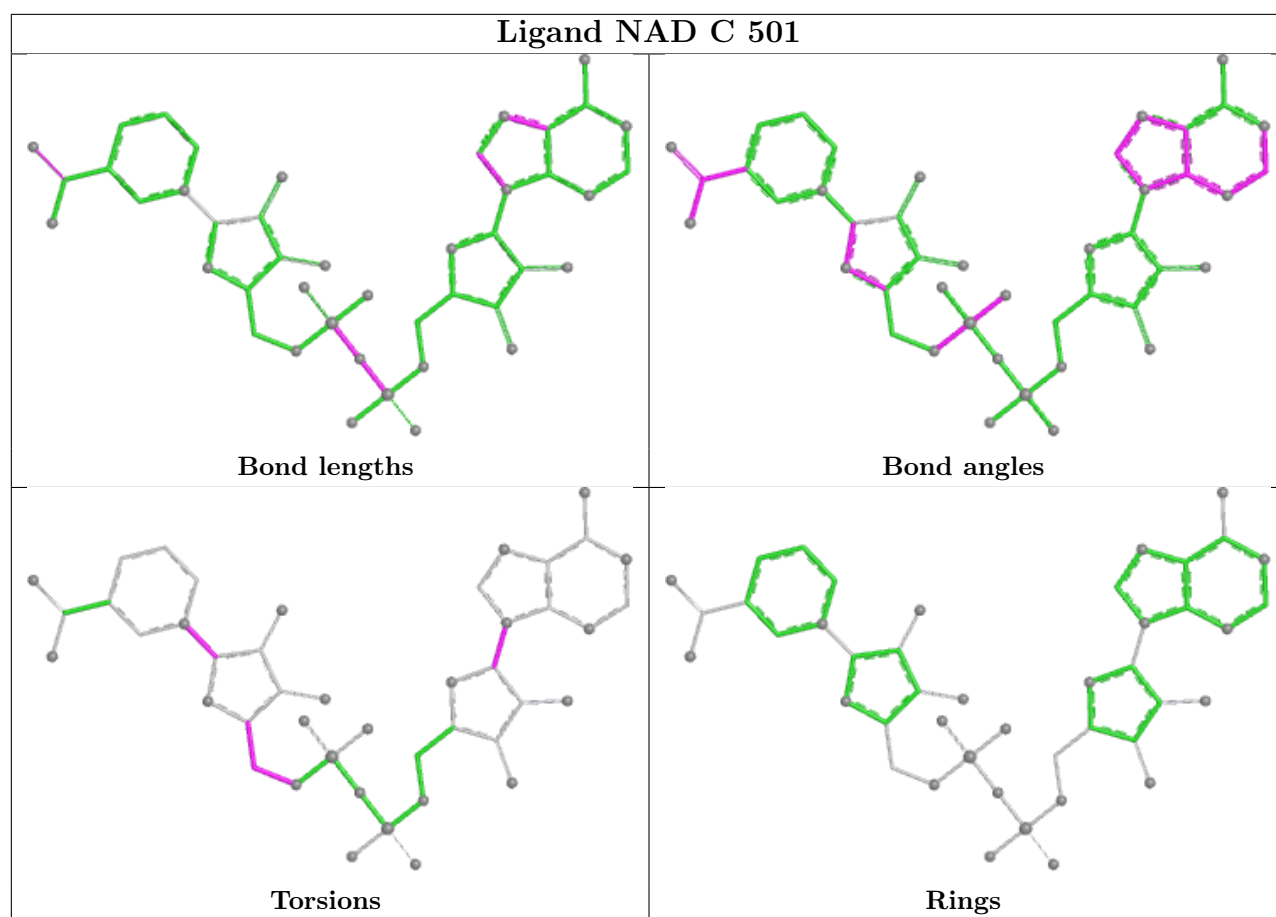
8 monomers are involved in 38 short contacts:

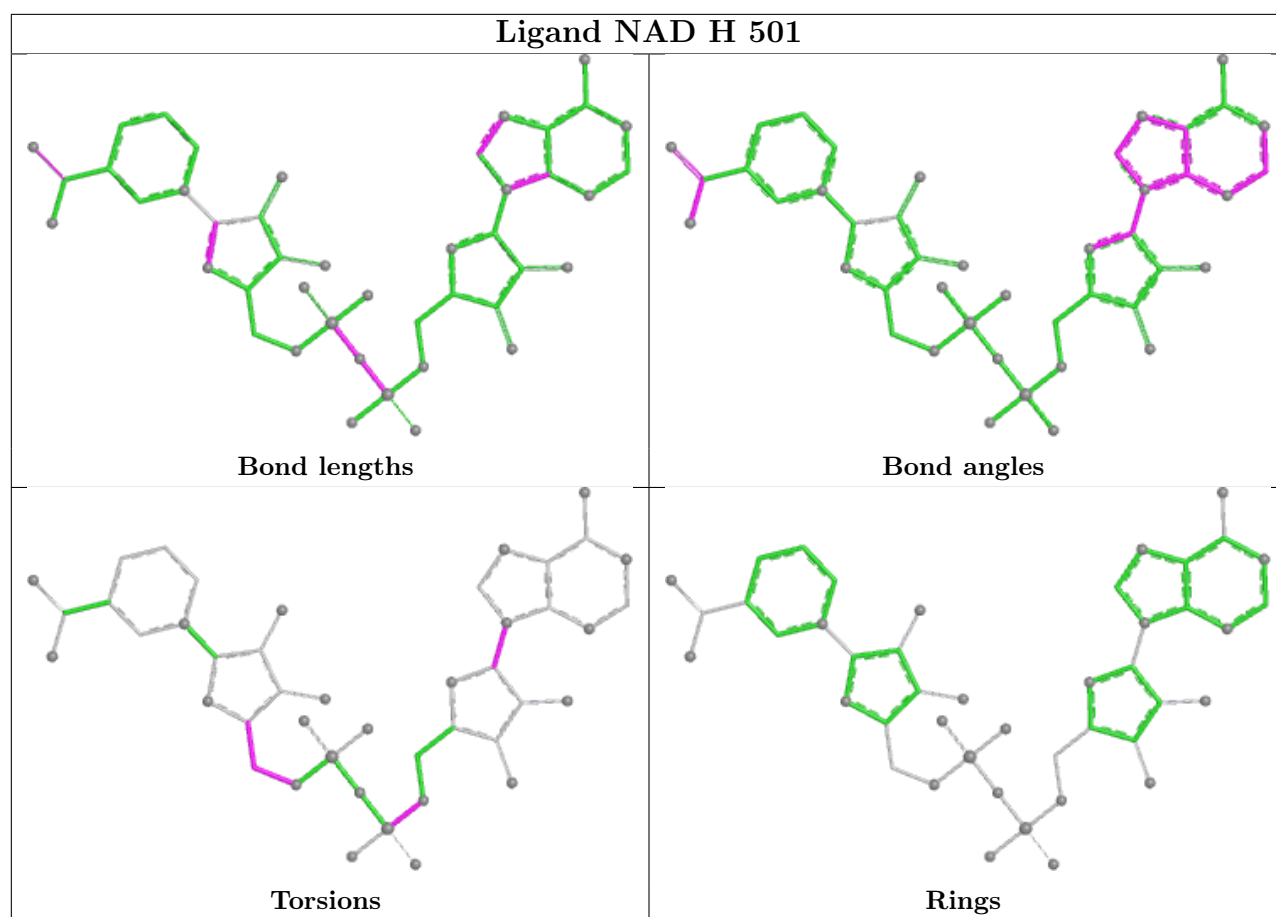
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	501	NAD	5	0
2	C	501	NAD	3	0
2	H	501	NAD	2	0
2	B	501	NAD	2	0
2	F	501	NAD	5	0
2	D	501	NAD	7	0
2	A	501	NAD	5	0
2	E	501	NAD	9	0

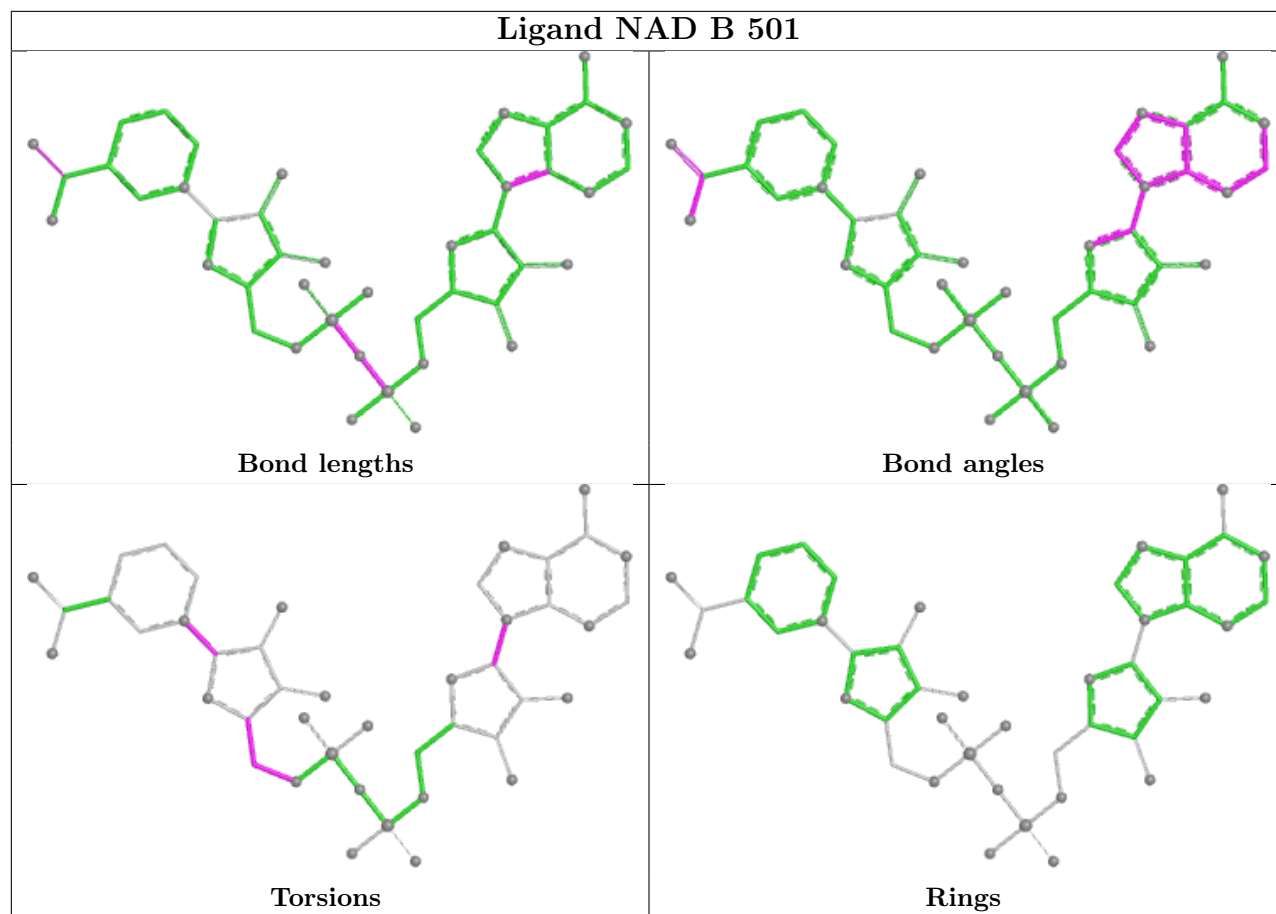
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

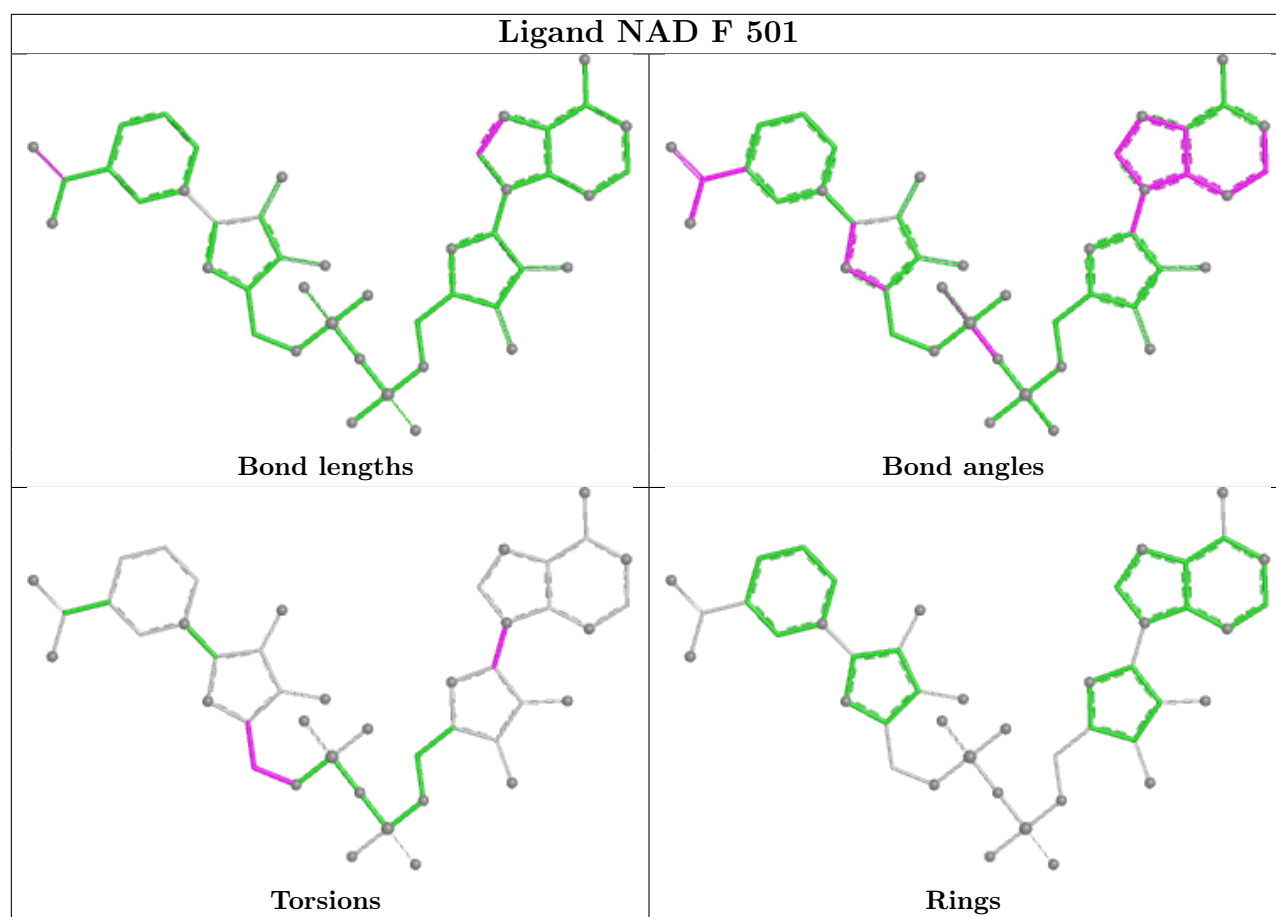
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. 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.

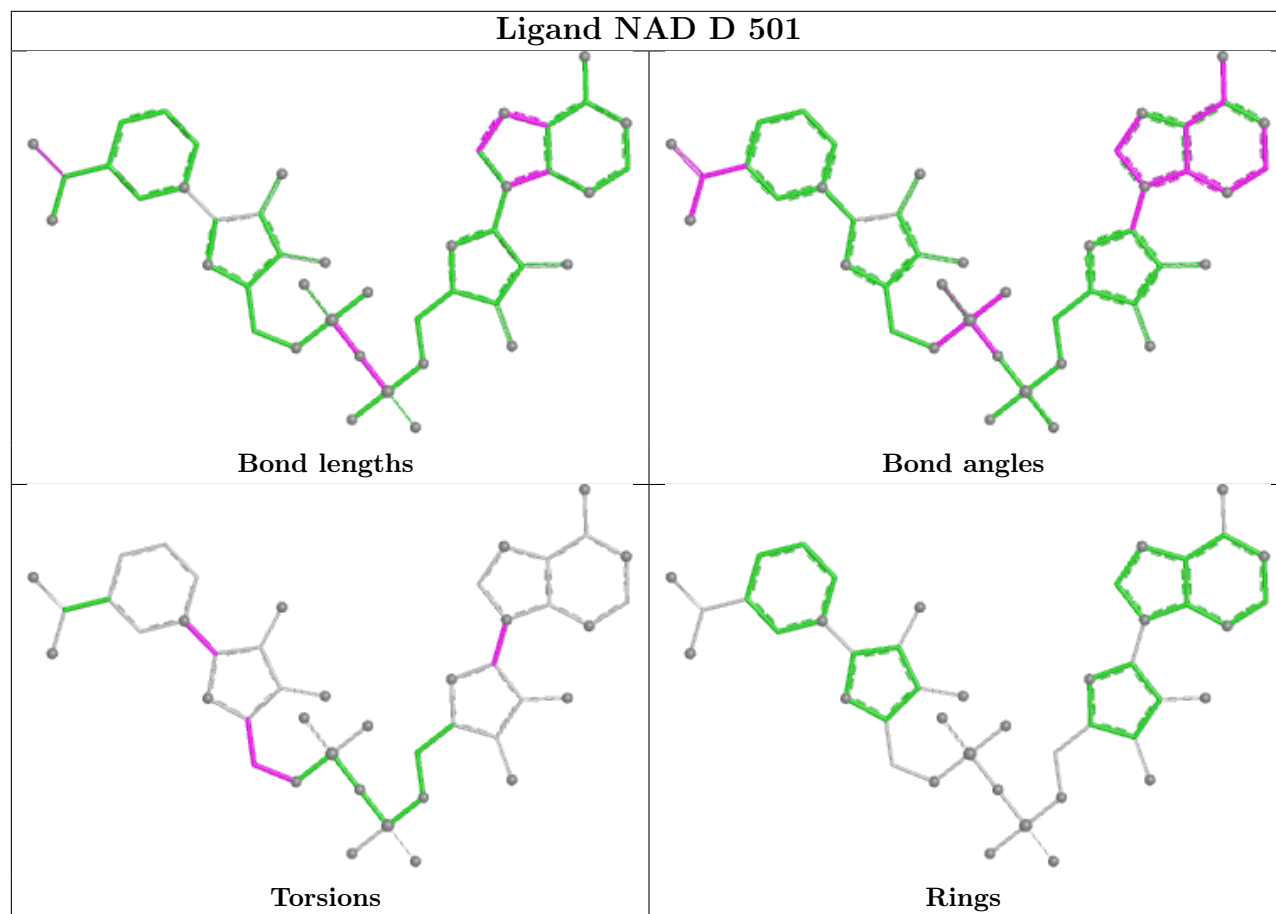


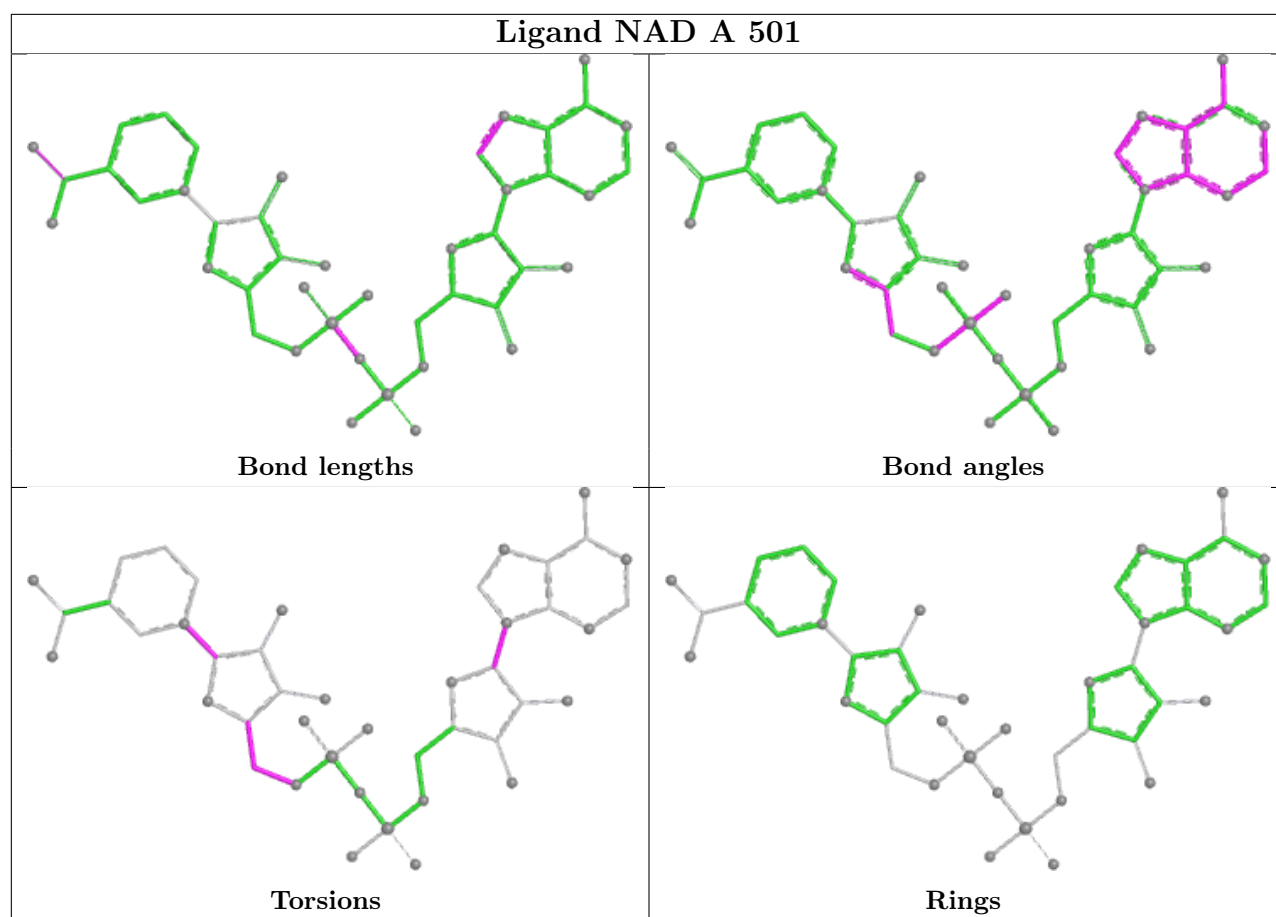


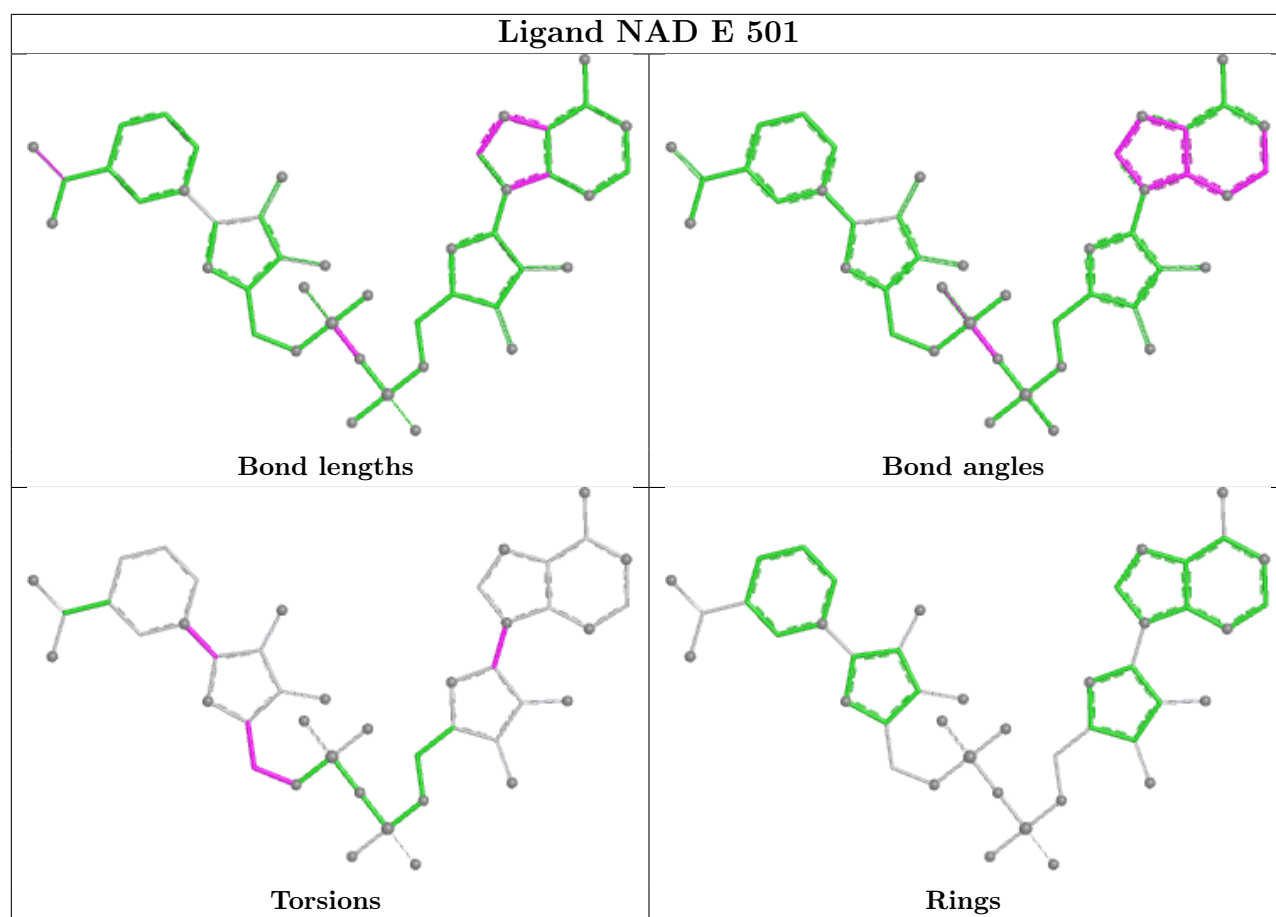












5.7 Other polymers [i](#)

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	502/517 (97%)	-0.62	0 100 100	21, 32, 53, 91	0
1	B	496/517 (95%)	-0.52	2 (0%) 88 86	17, 37, 59, 82	4 (0%)
1	C	496/517 (95%)	-0.47	1 (0%) 91 90	26, 40, 62, 116	0
1	D	496/517 (95%)	-0.53	0 100 100	16, 39, 63, 101	1 (0%)
1	E	496/517 (95%)	-0.31	1 (0%) 91 90	28, 49, 70, 96	0
1	F	496/517 (95%)	-0.50	1 (0%) 91 90	23, 39, 61, 94	0
1	G	496/517 (95%)	-0.46	2 (0%) 88 86	25, 41, 64, 104	0
1	H	496/517 (95%)	-0.42	0 100 100	26, 44, 65, 108	0
All	All	3974/4136 (96%)	-0.48	7 (0%) 91 90	16, 40, 63, 116	5 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	1	MET	3.4
1	B	1	MET	2.8
1	G	1	MET	2.5
1	C	1	MET	2.5
1	B	317	LYS	2.4
1	G	3	LEU	2.3
1	F	1	MET	2.2

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](#)

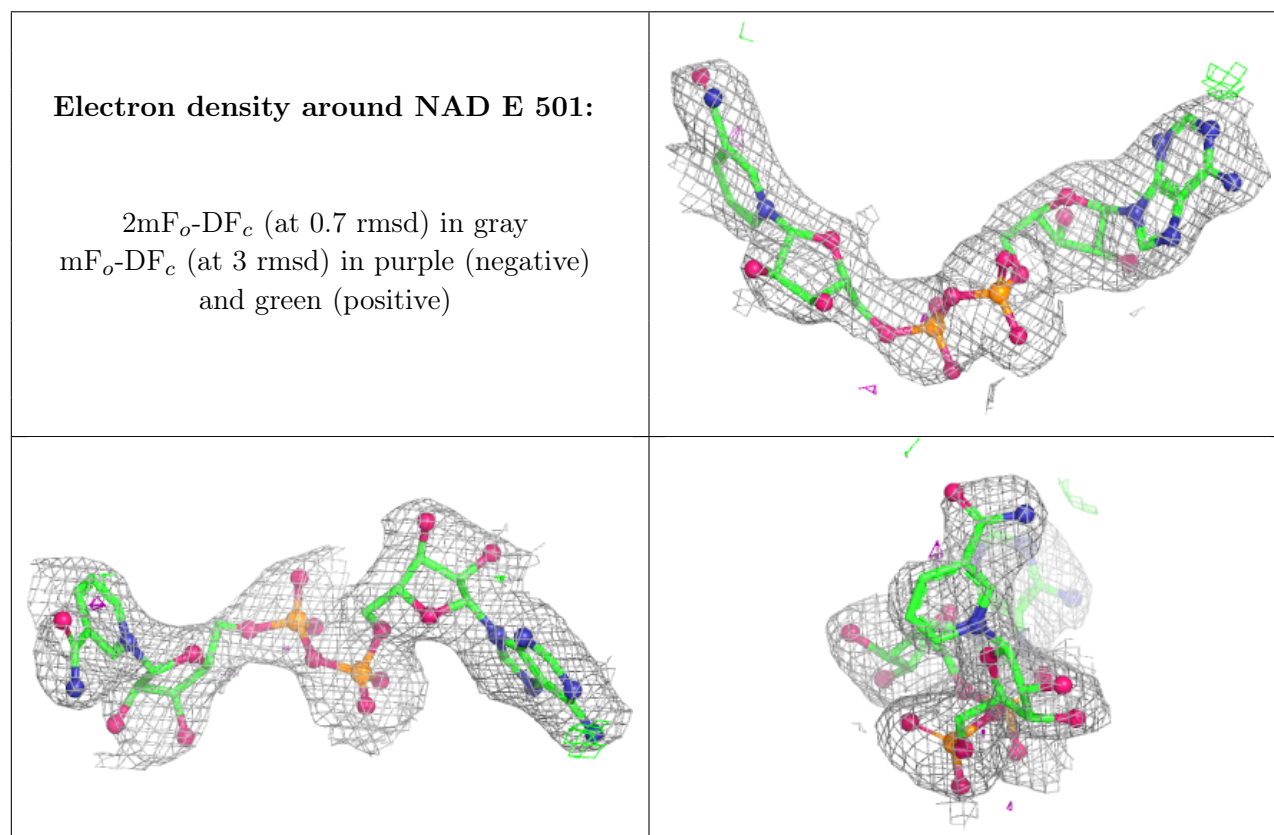
There are no oligosaccharides in this entry.

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.

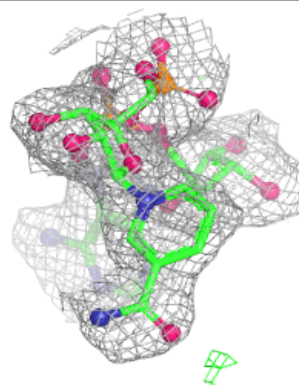
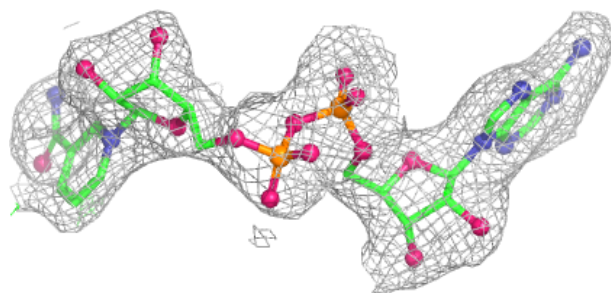
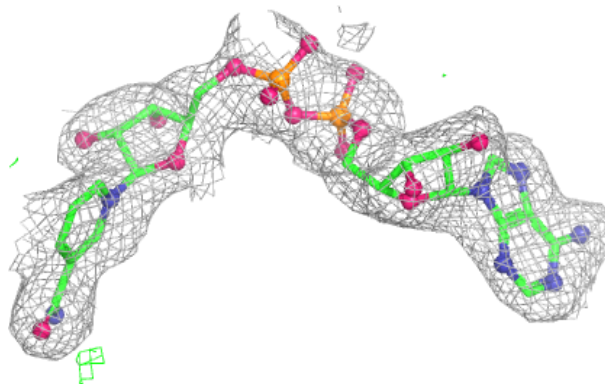
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ACT	B	502	4/4	0.74	0.18	43,45,50,53	0
2	NAD	E	501	44/44	0.95	0.08	34,47,59,60	0
2	NAD	G	501	44/44	0.96	0.06	29,36,51,65	0
2	NAD	D	501	44/44	0.96	0.06	23,32,41,54	0
2	NAD	B	501	44/44	0.97	0.06	23,35,50,54	0
2	NAD	F	501	44/44	0.97	0.06	24,33,39,44	0
2	NAD	C	501	44/44	0.97	0.06	35,41,49,54	0
2	NAD	H	501	44/44	0.97	0.06	33,41,48,58	0
2	NAD	A	501	44/44	0.97	0.05	23,30,39,46	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.

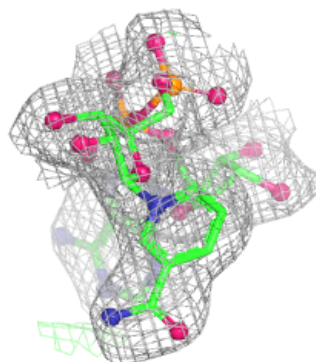
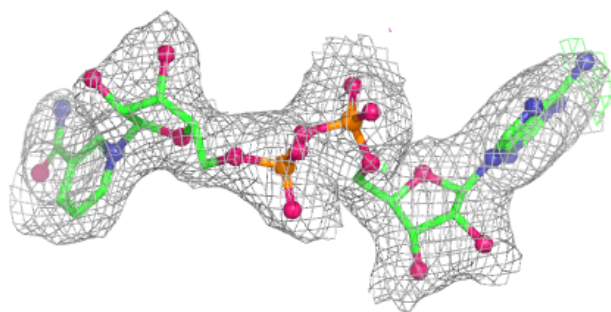
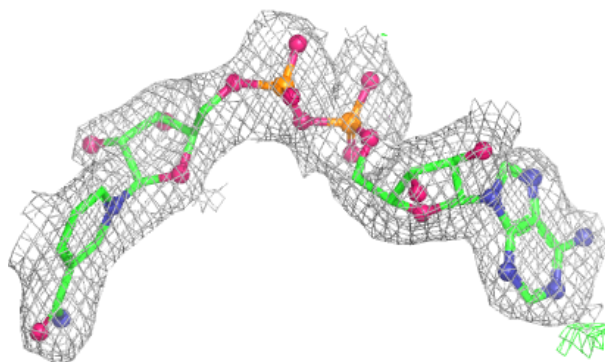


Electron density around NAD G 501:

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

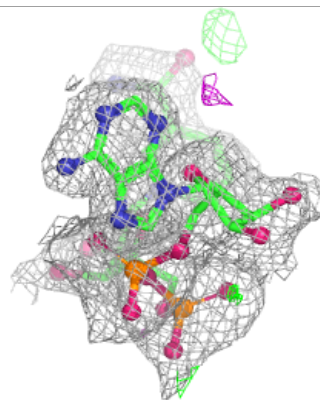
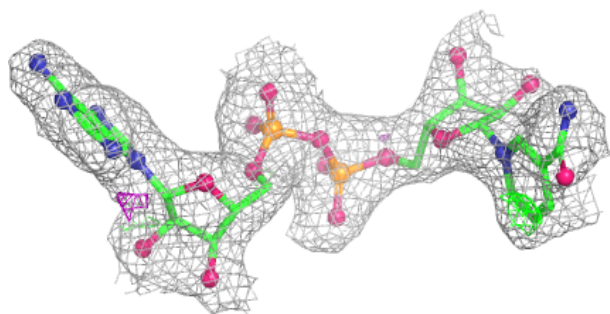
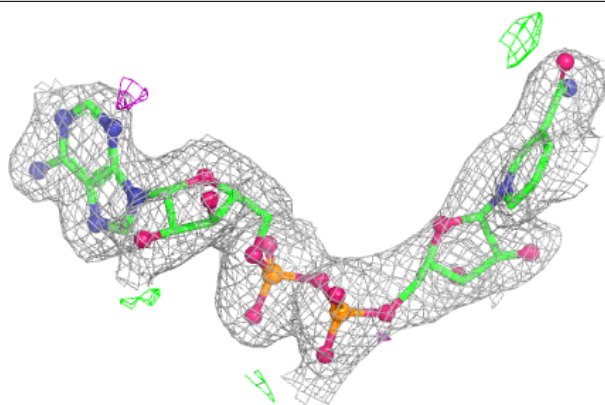
**Electron density around NAD D 501:**

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

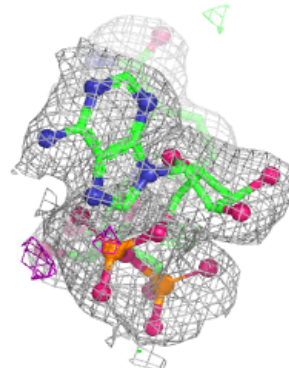
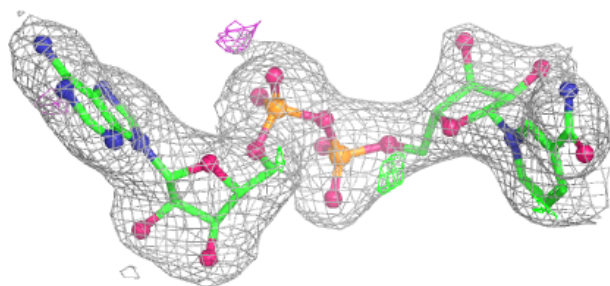
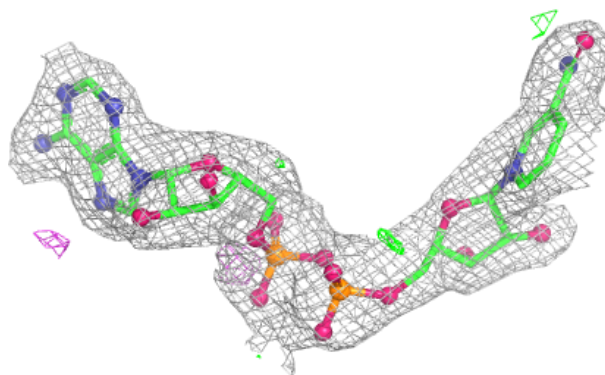


Electron density around NAD B 501:

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

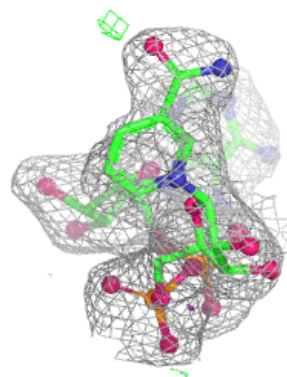
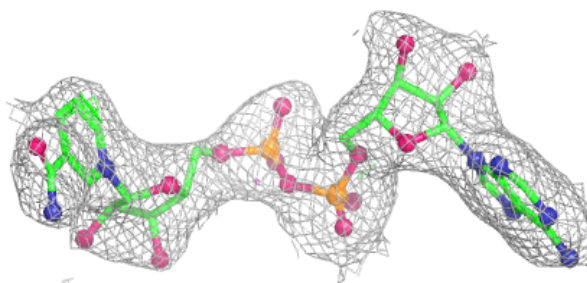
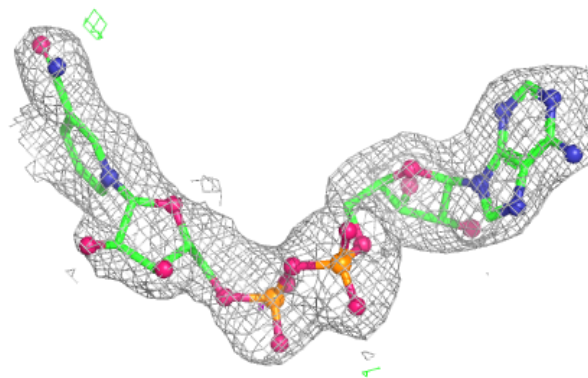
**Electron density around NAD F 501:**

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

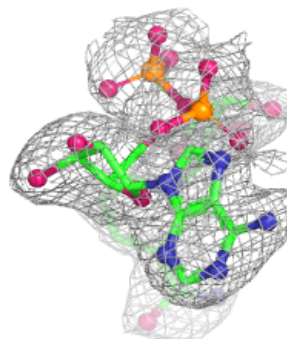
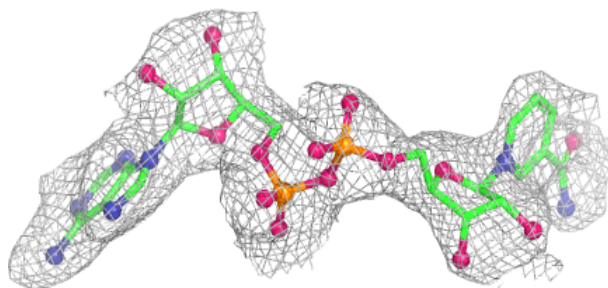
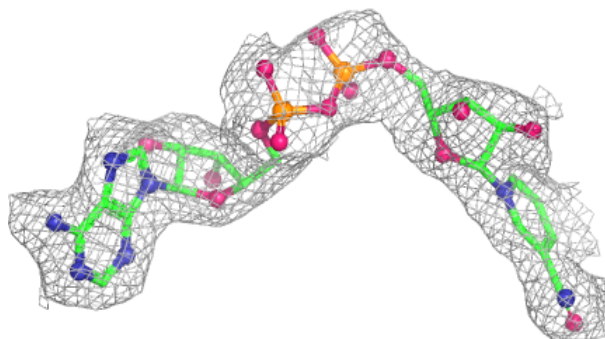


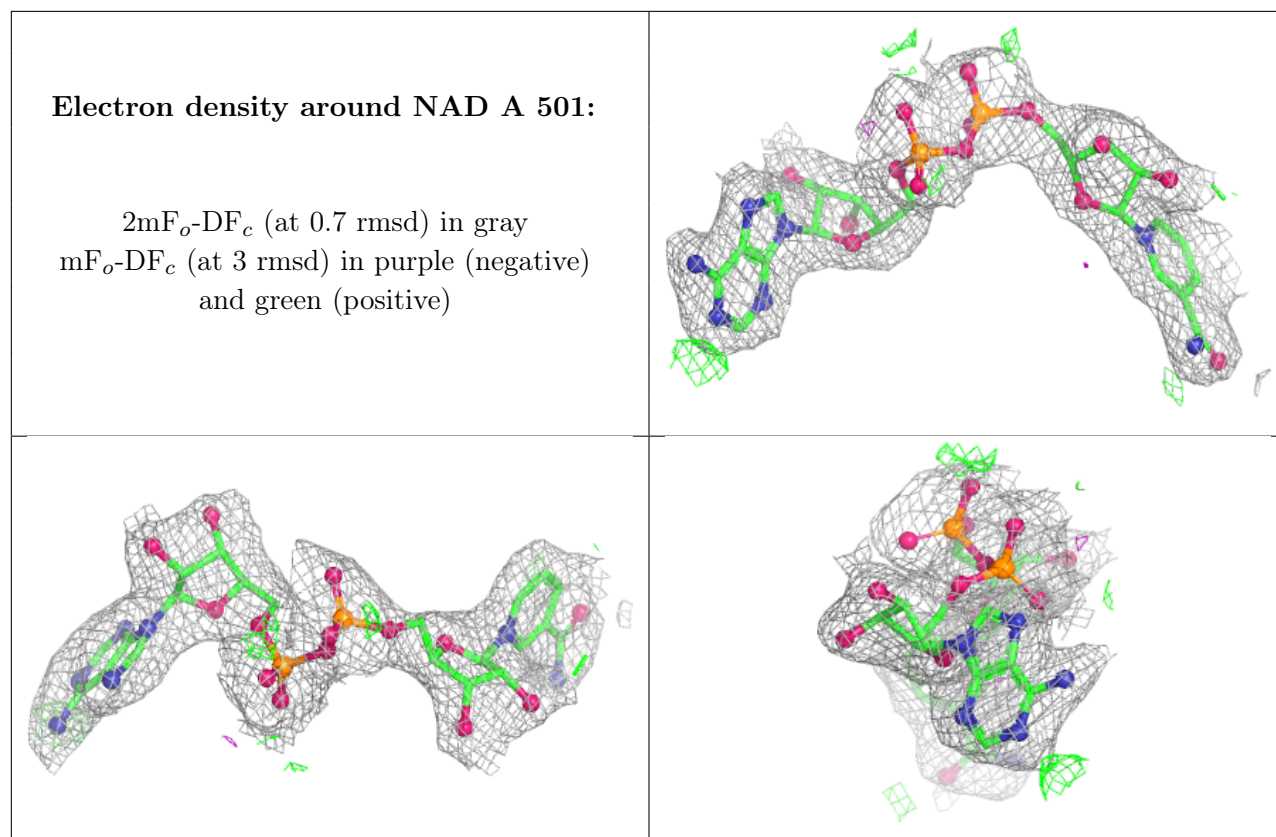
Electron density around NAD C 501:

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

**Electron density around NAD H 501:**

$2mF_o-DF_c$ (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.