



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:51 AM UTC

PDB ID : 5EZ4 / pdb_00005ez4
Title : 2.11 Angstrom resolution crystal structure of betaine aldehyde dehydrogenase (betB) P449M/Y450L double mutant from *Staphylococcus aureus* in complex with NAD⁺ and BME-modified 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 : 2015-11-26
Resolution : 2.11 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

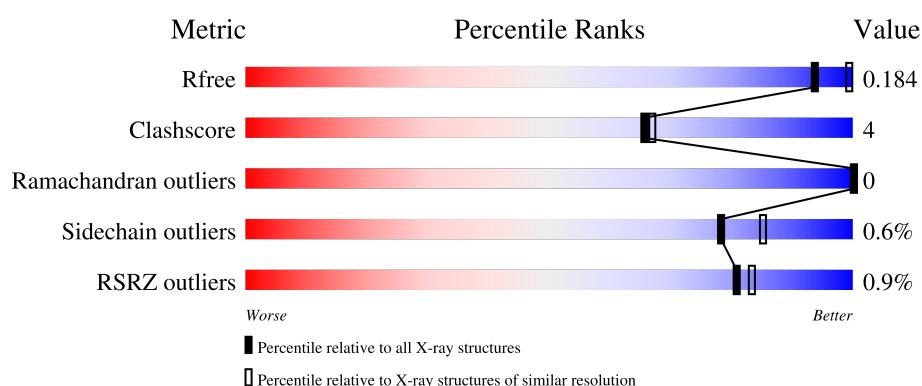
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	% 90% 6% .
1	B	517	% 89% 7% .
1	C	517	2% 89% 7% .
1	D	517	89% 5% 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 18119 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	496	Total	C	N	O	S	0	16	0
			3979	2498	678	787	16			
1	B	495	Total	C	N	O	S	0	15	0
			3952	2484	672	780	16			
1	C	498	Total	C	N	O	S	0	13	0
			3956	2485	672	781	18			
1	D	493	Total	C	N	O	S	0	12	0
			3910	2459	663	771	17			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q9L4P8
A	-19	GLY	-	expression tag	UNP Q9L4P8
A	-18	SER	-	expression tag	UNP Q9L4P8
A	-17	SER	-	expression tag	UNP Q9L4P8
A	-16	HIS	-	expression tag	UNP Q9L4P8
A	-15	HIS	-	expression tag	UNP Q9L4P8
A	-14	HIS	-	expression tag	UNP Q9L4P8
A	-13	HIS	-	expression tag	UNP Q9L4P8
A	-12	HIS	-	expression tag	UNP Q9L4P8
A	-11	HIS	-	expression tag	UNP Q9L4P8
A	-10	SER	-	expression tag	UNP Q9L4P8
A	-9	SER	-	expression tag	UNP Q9L4P8
A	-8	GLY	-	expression tag	UNP Q9L4P8
A	-7	ARG	-	expression tag	UNP Q9L4P8
A	-6	GLU	-	expression tag	UNP Q9L4P8
A	-5	ASN	-	expression tag	UNP Q9L4P8
A	-4	LEU	-	expression tag	UNP Q9L4P8
A	-3	TYR	-	expression tag	UNP Q9L4P8
A	-2	PHE	-	expression tag	UNP Q9L4P8
A	-1	GLN	-	expression tag	UNP Q9L4P8
A	0	GLY	-	expression tag	UNP Q9L4P8

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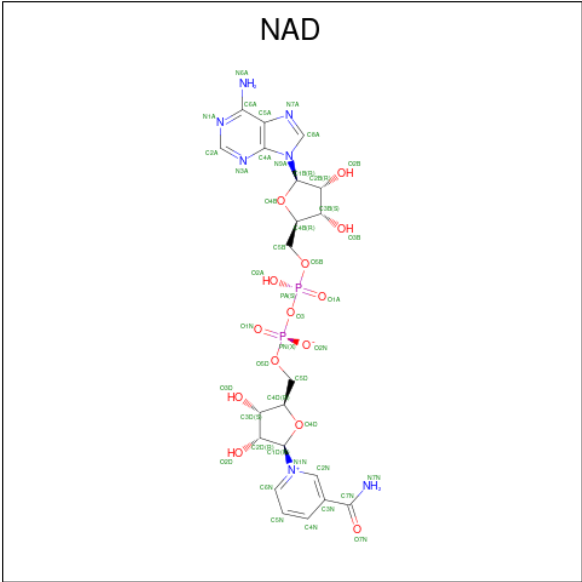
Chain	Residue	Modelled	Actual	Comment	Reference
A	449	MET	PRO	engineered mutation	UNP Q9L4P8
A	450	LEU	TYR	engineered mutation	UNP Q9L4P8
B	-20	MET	-	expression tag	UNP Q9L4P8
B	-19	GLY	-	expression tag	UNP Q9L4P8
B	-18	SER	-	expression tag	UNP Q9L4P8
B	-17	SER	-	expression tag	UNP Q9L4P8
B	-16	HIS	-	expression tag	UNP Q9L4P8
B	-15	HIS	-	expression tag	UNP Q9L4P8
B	-14	HIS	-	expression tag	UNP Q9L4P8
B	-13	HIS	-	expression tag	UNP Q9L4P8
B	-12	HIS	-	expression tag	UNP Q9L4P8
B	-11	HIS	-	expression tag	UNP Q9L4P8
B	-10	SER	-	expression tag	UNP Q9L4P8
B	-9	SER	-	expression tag	UNP Q9L4P8
B	-8	GLY	-	expression tag	UNP Q9L4P8
B	-7	ARG	-	expression tag	UNP Q9L4P8
B	-6	GLU	-	expression tag	UNP Q9L4P8
B	-5	ASN	-	expression tag	UNP Q9L4P8
B	-4	LEU	-	expression tag	UNP Q9L4P8
B	-3	TYR	-	expression tag	UNP Q9L4P8
B	-2	PHE	-	expression tag	UNP Q9L4P8
B	-1	GLN	-	expression tag	UNP Q9L4P8
B	0	GLY	-	expression tag	UNP Q9L4P8
B	449	MET	PRO	engineered mutation	UNP Q9L4P8
B	450	LEU	TYR	engineered mutation	UNP Q9L4P8
C	-20	MET	-	expression tag	UNP Q9L4P8
C	-19	GLY	-	expression tag	UNP Q9L4P8
C	-18	SER	-	expression tag	UNP Q9L4P8
C	-17	SER	-	expression tag	UNP Q9L4P8
C	-16	HIS	-	expression tag	UNP Q9L4P8
C	-15	HIS	-	expression tag	UNP Q9L4P8
C	-14	HIS	-	expression tag	UNP Q9L4P8
C	-13	HIS	-	expression tag	UNP Q9L4P8
C	-12	HIS	-	expression tag	UNP Q9L4P8
C	-11	HIS	-	expression tag	UNP Q9L4P8
C	-10	SER	-	expression tag	UNP Q9L4P8
C	-9	SER	-	expression tag	UNP Q9L4P8
C	-8	GLY	-	expression tag	UNP Q9L4P8
C	-7	ARG	-	expression tag	UNP Q9L4P8
C	-6	GLU	-	expression tag	UNP Q9L4P8
C	-5	ASN	-	expression tag	UNP Q9L4P8
C	-4	LEU	-	expression tag	UNP Q9L4P8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	TYR	-	expression tag	UNP Q9L4P8
C	-2	PHE	-	expression tag	UNP Q9L4P8
C	-1	GLN	-	expression tag	UNP Q9L4P8
C	0	GLY	-	expression tag	UNP Q9L4P8
C	449	MET	PRO	engineered mutation	UNP Q9L4P8
C	450	LEU	TYR	engineered mutation	UNP Q9L4P8
D	-20	MET	-	expression tag	UNP Q9L4P8
D	-19	GLY	-	expression tag	UNP Q9L4P8
D	-18	SER	-	expression tag	UNP Q9L4P8
D	-17	SER	-	expression tag	UNP Q9L4P8
D	-16	HIS	-	expression tag	UNP Q9L4P8
D	-15	HIS	-	expression tag	UNP Q9L4P8
D	-14	HIS	-	expression tag	UNP Q9L4P8
D	-13	HIS	-	expression tag	UNP Q9L4P8
D	-12	HIS	-	expression tag	UNP Q9L4P8
D	-11	HIS	-	expression tag	UNP Q9L4P8
D	-10	SER	-	expression tag	UNP Q9L4P8
D	-9	SER	-	expression tag	UNP Q9L4P8
D	-8	GLY	-	expression tag	UNP Q9L4P8
D	-7	ARG	-	expression tag	UNP Q9L4P8
D	-6	GLU	-	expression tag	UNP Q9L4P8
D	-5	ASN	-	expression tag	UNP Q9L4P8
D	-4	LEU	-	expression tag	UNP Q9L4P8
D	-3	TYR	-	expression tag	UNP Q9L4P8
D	-2	PHE	-	expression tag	UNP Q9L4P8
D	-1	GLN	-	expression tag	UNP Q9L4P8
D	0	GLY	-	expression tag	UNP Q9L4P8
D	449	MET	PRO	engineered mutation	UNP Q9L4P8
D	450	LEU	TYR	engineered mutation	UNP Q9L4P8

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

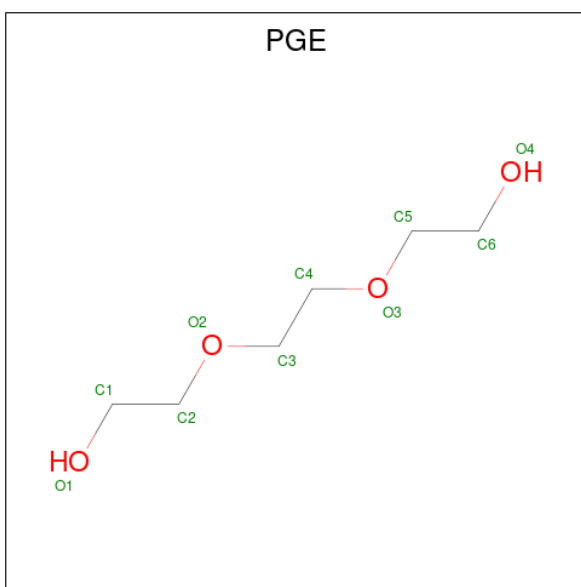


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

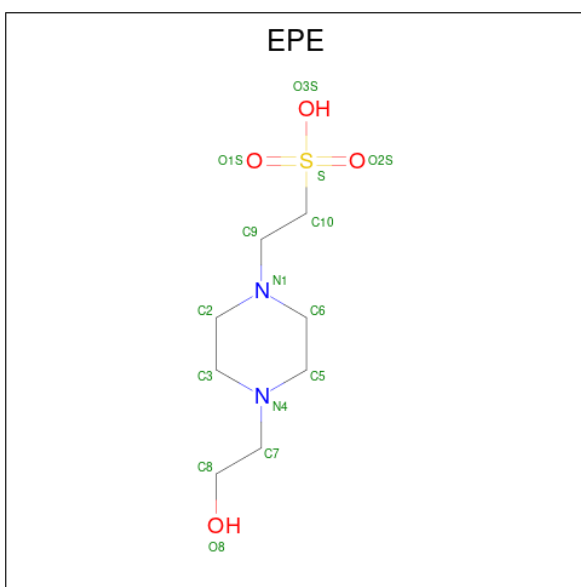
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		
3	C	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		

- Molecule 4 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	C	1	Total	C	O	0	0
			10	6	4		
4	C	1	Total	C	O	0	0
			10	6	4		
4	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



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

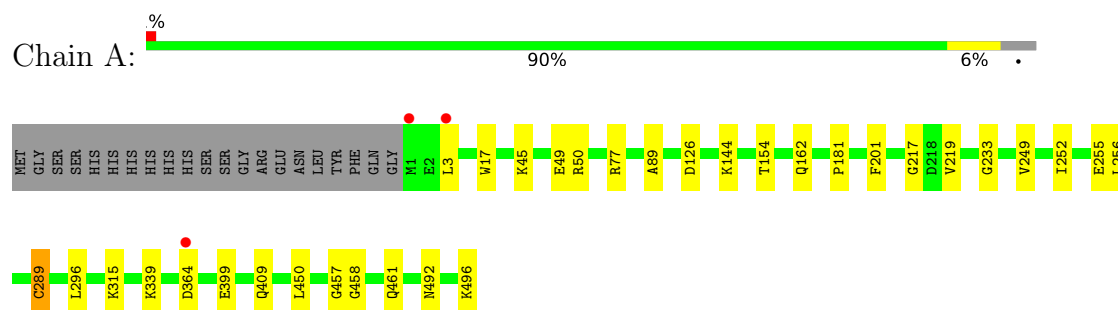
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	498	Total	O	0	10
			505	505		
6	B	483	Total	O	0	9
			487	487		
6	C	515	Total	O	0	6
			519	519		
6	D	505	Total	O	0	3
			506	506		

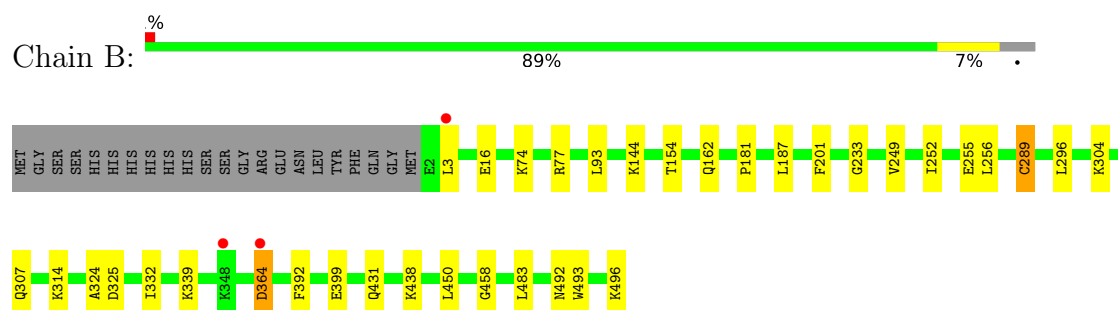
3 Residue-property plots [i](#)

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.

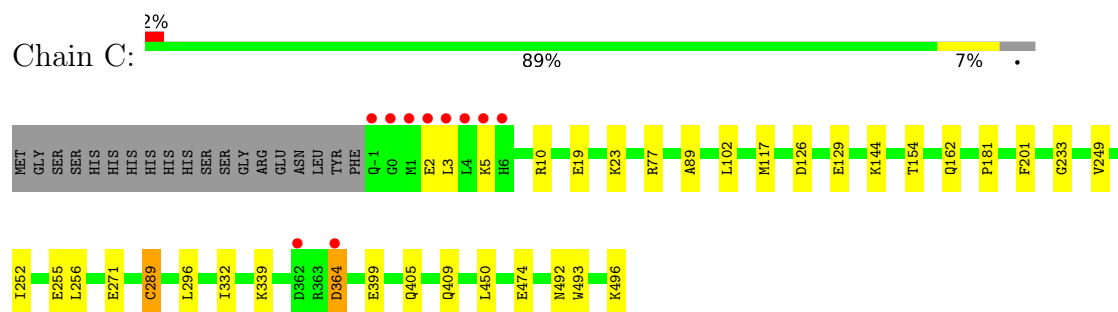
- Molecule 1: Betaine aldehyde dehydrogenase



- Molecule 1: Betaine aldehyde dehydrogenase

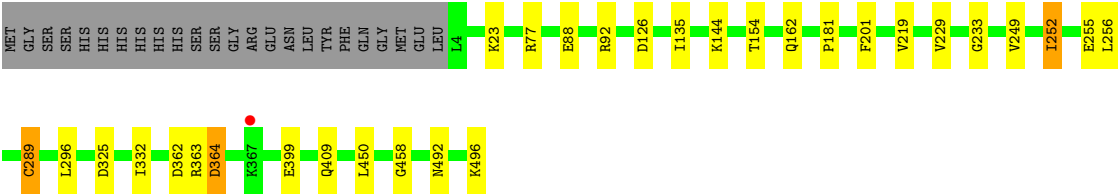


- Molecule 1: Betaine aldehyde dehydrogenase



- Molecule 1: Betaine aldehyde dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	225.03Å 102.16Å 118.33Å 90.00° 104.85° 90.00°	Depositor
Resolution (Å)	29.69 – 2.11 29.69 – 2.11	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.69-2.11) 99.9 (29.69-2.11)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 2.12Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.141 , 0.175 0.152 , 0.184	Depositor DCC
R_{free} test set	7478 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18119	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% 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: NA, CME, PGE, 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	0.77	0/4037	0.88	6/5453 (0.1%)
1	B	0.78	0/4013	0.89	7/5424 (0.1%)
1	C	0.77	0/4014	0.89	4/5423 (0.1%)
1	D	0.78	0/3968	0.89	5/5362 (0.1%)
All	All	0.78	0/16032	0.89	22/21662 (0.1%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	187	LEU	N-CA-C	6.78	118.32	111.07
1	A	458	GLY	N-CA-C	6.38	122.36	111.19
1	B	364	ASP	N-CA-C	6.09	117.59	111.07
1	C	364	ASP	N-CA-C	5.85	117.33	111.07
1	B	201	PHE	CA-C-N	-5.80	113.81	119.78
1	B	201	PHE	C-N-CA	-5.80	113.81	119.78
1	B	458	GLY	N-CA-C	5.80	121.30	111.04
1	C	201	PHE	CA-C-N	-5.77	114.00	119.89
1	C	201	PHE	C-N-CA	-5.77	114.00	119.89
1	D	332	ILE	N-CA-C	5.63	118.13	111.09
1	A	457	GLY	CA-C-N	-5.59	118.33	122.18
1	A	457	GLY	C-N-CA	-5.59	118.33	122.18
1	C	332	ILE	N-CA-C	5.56	118.89	111.44
1	B	332	ILE	N-CA-C	5.49	118.80	111.44
1	A	364	ASP	N-CA-C	5.44	117.63	111.11
1	B	304	LYS	CB-CG-CD	5.38	123.66	111.30
1	D	364	ASP	N-CA-C	5.35	116.80	111.07
1	D	458	GLY	N-CA-C	5.23	120.02	110.71
1	D	201	PHE	CA-C-N	-5.05	114.58	119.78
1	D	201	PHE	C-N-CA	-5.05	114.58	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	PHE	CA-C-N	-5.02	114.61	119.78
1	A	201	PHE	C-N-CA	-5.02	114.61	119.78

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	3979	0	3910	23	0
1	B	3952	0	3892	35	0
1	C	3956	0	3890	40	0
1	D	3910	0	3845	22	0
2	A	44	0	26	3	0
2	B	44	0	26	4	0
2	C	44	0	26	4	0
2	D	44	0	26	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	10	0	14	1	0
4	B	10	0	14	0	0
4	C	20	0	28	0	0
4	D	10	0	14	0	0
5	A	30	0	36	2	0
5	B	30	0	36	1	0
5	C	15	0	18	2	0
6	A	505	0	0	13	0
6	B	487	0	0	9	0
6	C	519	0	0	21	0
6	D	506	0	0	11	0
All	All	18119	0	15801	123	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 (123) 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[A]:ASP:HB3	6:D:601[A]:HOH:O	1.61	0.97
1:B:3:LEU:HD23	1:B:93:LEU:HG	1.53	0.90
1:D:409[A]:GLN:HG2	6:D:1027:HOH:O	1.72	0.88
1:B:77[B]:ARG:NH1	1:C:117[B]:MET:CE	2.39	0.85
1:C:409[A]:GLN:HG2	6:C:1062:HOH:O	1.78	0.83
1:B:77[B]:ARG:HD3	1:C:77:ARG:HD3	1.60	0.82
1:A:126[A]:ASP:HB3	6:A:606[A]:HOH:O	1.78	0.82
1:C:3:LEU:HD23	6:C:1054:HOH:O	1.83	0.79
1:A:77[A]:ARG:HD3	1:D:77:ARG:HD3	1.63	0.77
1:D:126[A]:ASP:CB	6:D:601[A]:HOH:O	2.23	0.76
1:B:249:VAL:HG12	1:B:249:VAL:O	1.85	0.76
1:D:126[A]:ASP:CG	6:D:601[A]:HOH:O	2.30	0.75
1:C:249:VAL:HG12	1:C:249:VAL:O	1.86	0.75
1:A:249:VAL:O	1:A:249:VAL:HG12	1.87	0.74
2:C:501:NAD:C5N	6:C:969:HOH:O	2.37	0.72
1:D:144:LYS:HD2	6:D:780:HOH:O	1.88	0.72
1:B:77[B]:ARG:NH1	1:C:117[B]:MET:HE2	2.05	0.72
1:C:126[B]:ASP:CG	6:C:606[B]:HOH:O	2.32	0.71
1:C:126[B]:ASP:HB3	6:C:606[B]:HOH:O	1.92	0.69
1:D:77:ARG:NH2	6:D:602:HOH:O	2.27	0.66
1:B:77[B]:ARG:CZ	1:C:117[B]:MET:CE	2.74	0.64
1:B:77[B]:ARG:NH2	6:B:605:HOH:O	2.30	0.63
1:B:77[B]:ARG:CZ	1:C:117[B]:MET:HE1	2.30	0.62
1:D:162[A]:GLN:NE2	1:D:289:CME:HE2	2.14	0.62
6:A:983:HOH:O	1:B:144:LYS:HE2	2.01	0.61
1:C:126[B]:ASP:CB	6:C:606[B]:HOH:O	2.48	0.61
6:A:701:HOH:O	1:B:438:LYS:HE3	2.01	0.59
1:A:162[A]:GLN:NE2	1:A:289:CME:HE2	2.18	0.59
2:B:501:NAD:H5N	6:B:911:HOH:O	2.02	0.59
1:A:409[A]:GLN:HG3	6:A:781:HOH:O	2.03	0.58
1:B:307:GLN:CD	6:B:610:HOH:O	2.48	0.57
1:B:162:GLN:NE2	1:B:289:CME:HE2	2.19	0.56
1:C:77:ARG:NH2	6:C:613:HOH:O	2.38	0.56
1:C:19:GLU:OE1	1:C:23:LYS:HE3	2.05	0.56
1:C:409[A]:GLN:HG3	6:C:861:HOH:O	2.05	0.56
2:C:501:NAD:C4N	6:C:969:HOH:O	2.53	0.56
1:B:77[B]:ARG:NH1	1:C:117[B]:MET:HE1	2.20	0.55
2:A:501:NAD:H5N	6:A:957:HOH:O	2.07	0.55
1:C:162:GLN:NE2	1:C:289:CME:HE2	2.22	0.54
2:A:501:NAD:C5N	6:A:957:HOH:O	2.55	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:492:ASN:OD1	1:B:496:LYS:HE3	2.07	0.54
1:A:492:ASN:OD1	1:A:496:LYS:HE3	2.08	0.54
1:B:364:ASP:O	1:B:364:ASP:OD1	2.26	0.54
1:B:16:GLU:HG3	6:B:1077:HOH:O	2.08	0.53
1:C:339:LYS:HD3	6:C:841:HOH:O	2.08	0.53
2:D:501:NAD:H5N	6:D:951:HOH:O	2.08	0.53
1:C:493:TRP:CH2	5:C:505:EPE:H71	2.44	0.53
1:A:450:LEU:HD12	1:A:450:LEU:C	2.34	0.53
1:C:409[A]:GLN:CG	6:C:1062:HOH:O	2.47	0.53
1:C:3:LEU:HD22	1:C:89:ALA:HB1	1.89	0.52
2:C:501:NAD:H5N	6:C:969:HOH:O	2.06	0.52
1:C:450:LEU:C	1:C:450:LEU:HD12	2.33	0.52
1:B:450:LEU:C	1:B:450:LEU:HD12	2.35	0.52
5:A:505:EPE:H81	1:B:324:ALA:O	2.10	0.51
1:C:2:GLU:HA	1:C:5:LYS:CE	2.40	0.51
1:C:405[A]:GLN:HG3	6:C:1018:HOH:O	2.10	0.51
6:A:1014:HOH:O	1:B:431:GLN:HG2	2.09	0.51
1:C:364:ASP:OD1	1:C:364:ASP:O	2.28	0.51
2:B:501:NAD:C5N	6:B:911:HOH:O	2.58	0.51
1:A:126[A]:ASP:CB	6:A:606[A]:HOH:O	2.49	0.51
1:D:450:LEU:C	1:D:450:LEU:HD12	2.35	0.50
2:D:501:NAD:C5N	6:D:951:HOH:O	2.59	0.50
1:A:50[B]:ARG:NH2	6:A:611:HOH:O	2.42	0.49
1:B:493:TRP:HA	5:B:505:EPE:H21	1.94	0.49
1:B:364:ASP:OD1	1:B:364:ASP:C	2.54	0.49
1:A:3:LEU:HD22	1:A:89:ALA:HB1	1.94	0.49
1:D:296:LEU:HD23	1:D:399:GLU:HB2	1.95	0.49
1:D:325[B]:ASP:OD1	1:D:325[B]:ASP:N	2.45	0.49
1:D:364:ASP:O	1:D:364:ASP:OD1	2.30	0.48
1:D:233:GLY:O	1:D:256:LEU:HA	2.13	0.48
1:B:154:THR:OG1	1:B:181:PRO:HA	2.13	0.48
1:A:233:GLY:O	1:A:256:LEU:HA	2.14	0.48
4:A:503:PGE:H42	6:A:944:HOH:O	2.15	0.47
1:B:74:LYS:HE2	6:B:799:HOH:O	2.13	0.47
1:D:88:GLU:OE2	1:D:92:ARG:NH1	2.47	0.47
2:D:501:NAD:H5N	6:D:623:HOH:O	2.14	0.47
1:C:3:LEU:CD2	1:C:89:ALA:HB1	2.45	0.47
1:C:233:GLY:O	1:C:256:LEU:HA	2.14	0.47
1:C:249:VAL:O	1:C:249:VAL:CG1	2.60	0.47
1:A:296:LEU:HD23	1:A:399:GLU:HB2	1.97	0.47
1:C:296:LEU:HD23	1:C:399:GLU:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:364:ASP:OD1	1:C:364:ASP:C	2.53	0.46
1:C:271:GLU:HG2	6:C:1014:HOH:O	2.16	0.46
1:B:296:LEU:HD23	1:B:399:GLU:HB2	1.96	0.46
1:B:74:LYS:CE	6:B:799:HOH:O	2.64	0.46
1:D:23:LYS:HE2	6:D:642:HOH:O	2.16	0.46
1:C:154:THR:OG1	1:C:181:PRO:HA	2.15	0.46
1:B:144:LYS:HD2	6:B:754:HOH:O	2.15	0.45
1:B:392:PHE:CG	2:B:501:NAD:H2D	2.51	0.45
1:C:474:GLU:HG2	6:C:1050:HOH:O	2.16	0.45
1:B:483:LEU:C	1:B:483:LEU:HD13	2.41	0.45
1:D:154:THR:OG1	1:D:181:PRO:HA	2.16	0.45
1:A:154:THR:OG1	1:A:181:PRO:HA	2.17	0.45
1:B:492:ASN:OD1	1:B:496:LYS:CE	2.65	0.45
1:D:364:ASP:OD1	1:D:364:ASP:C	2.55	0.45
1:B:339:LYS:HE3	2:B:501:NAD:O3D	2.17	0.45
1:B:233:GLY:O	1:B:256:LEU:HA	2.17	0.43
2:C:501:NAD:O5D	2:C:501:NAD:H6N	2.17	0.43
1:B:314:LYS:NZ	6:B:616:HOH:O	2.46	0.43
1:D:409[A]:GLN:HG3	6:D:708:HOH:O	2.19	0.43
1:A:17:TRP:CD1	5:A:504:EPE:H92	2.54	0.43
1:C:102:LEU:HD23	1:C:102:LEU:HA	1.86	0.43
1:A:144:LYS:HD2	6:A:811:HOH:O	2.18	0.42
1:A:461:GLN:HA	1:B:249:VAL:HB	2.01	0.42
1:D:362:ASP:OD1	1:D:363:ARG:N	2.52	0.42
1:D:229[B]:VAL:CG1	1:D:252:ILE:HG22	2.50	0.42
1:A:49[A]:GLU:HG3	1:A:219:VAL:HG11	2.02	0.42
1:A:315:LYS:HE3	1:B:496:LYS:O	2.19	0.42
1:C:496:LYS:HB3	1:C:496:LYS:HE2	1.85	0.42
1:A:339:LYS:CD	6:A:755:HOH:O	2.67	0.42
1:B:325[B]:ASP:OD1	1:B:325[B]:ASP:N	2.53	0.42
1:C:129:GLU:HG3	6:C:842:HOH:O	2.18	0.42
1:A:126[A]:ASP:CG	6:A:606[A]:HOH:O	2.61	0.41
1:A:217:GLY:HA3	2:A:501:NAD:C2A	2.51	0.41
1:C:162:GLN:HG3	6:C:933:HOH:O	2.21	0.41
1:D:492:ASN:OD1	1:D:496:LYS:HE3	2.21	0.41
1:C:144:LYS:HD2	6:C:867:HOH:O	2.21	0.41
1:C:492:ASN:O	5:C:505:EPE:H92	2.20	0.41
1:A:45:LYS:HB2	1:A:219:VAL:HG21	2.03	0.41
1:C:339:LYS:CE	6:C:802[A]:HOH:O	2.69	0.41
1:A:249:VAL:O	1:A:249:VAL:CG1	2.62	0.41
1:C:10:ARG:HD2	6:C:983:HOH:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:995:HOH:O	1:D:135:ILE:HD11	2.20	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	509/517 (98%)	500 (98%)	9 (2%)	0	100	100
1	B	507/517 (98%)	497 (98%)	10 (2%)	0	100	100
1	C	508/517 (98%)	495 (97%)	13 (3%)	0	100	100
1	D	502/517 (97%)	494 (98%)	8 (2%)	0	100	100
All	All	2026/2068 (98%)	1986 (98%)	40 (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	425/428 (99%)	423 (100%)	2 (0%)	81	87
1	B	423/428 (99%)	421 (100%)	2 (0%)	81	87
1	C	423/428 (99%)	421 (100%)	2 (0%)	81	87
1	D	418/428 (98%)	414 (99%)	4 (1%)	68	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1689/1712 (99%)	1679 (99%)	10 (1%)	78	85

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	252	ILE
1	A	255	GLU
1	B	252	ILE
1	B	255	GLU
1	C	252	ILE
1	C	255	GLU
1	D	219	VAL
1	D	249	VAL
1	D	252	ILE
1	D	255	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	251	ASN
1	A	435	ASN
1	B	162	GLN
1	B	251	ASN
1	C	6	HIS
1	C	251	ASN
1	D	251	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CME	B	289	1	8,9,10	0.67	0	6,9,11	1.57	2 (33%)
1	CME	D	289	1	8,9,10	0.56	0	6,9,11	1.45	2 (33%)
1	CME	C	289	1	8,9,10	0.63	0	6,9,11	1.31	2 (33%)
1	CME	A	289	1	8,9,10	0.84	0	6,9,11	1.41	2 (33%)

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
1	CME	B	289	1	-	3/5/8/10	-
1	CME	D	289	1	-	2/5/8/10	-
1	CME	C	289	1	-	2/5/8/10	-
1	CME	A	289	1	-	3/5/8/10	-

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	289	CME	CB-SG-SD	-2.96	96.19	103.86
1	D	289	CME	CB-SG-SD	-2.62	97.09	103.86
1	A	289	CME	CB-SG-SD	-2.39	97.68	103.86
1	B	289	CME	CE-SD-SG	-2.22	93.71	103.46
1	A	289	CME	CE-SD-SG	-2.20	93.80	103.46
1	C	289	CME	CB-SG-SD	-2.20	98.18	103.86
1	D	289	CME	CE-SD-SG	-2.19	93.85	103.46
1	C	289	CME	CE-SD-SG	-2.16	93.99	103.46

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	289	CME	CE-SD-SG-CB
1	A	289	CME	SD-CE-CZ-OH

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Mol	Chain	Res	Type	Atoms
1	B	289	CME	CE-SD-SG-CB
1	B	289	CME	SD-CE-CZ-OH
1	C	289	CME	CE-SD-SG-CB
1	C	289	CME	SD-CE-CZ-OH
1	D	289	CME	CE-SD-SG-CB
1	D	289	CME	SD-CE-CZ-OH
1	A	289	CME	CA-CB-SG-SD
1	B	289	CME	CA-CB-SG-SD

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	289	CME	1	0
1	D	289	CME	1	0
1	C	289	CME	1	0
1	A	289	CME	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 for Mogul analysis.

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	$\# Z > 2$
5	EPE	B	505	-	15,15,15	1.84	1 (6%)	19,20,20	1.42	4 (21%)
2	NAD	D	501	-	46,48,48	1.46	7 (15%)	64,73,73	1.65	13 (20%)
4	PGE	A	503	-	9,9,9	0.61	0	8,8,8	0.63	0
4	PGE	B	503	-	9,9,9	0.66	0	8,8,8	0.77	0
5	EPE	A	505	-	15,15,15	1.87	1 (6%)	19,20,20	1.19	2 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAD	A	501	-	46,48,48	1.53	9 (19%)	64,73,73	1.87	14 (21%)
4	PGE	C	503	-	9,9,9	0.67	0	8,8,8	0.55	0
5	EPE	B	504	-	15,15,15	1.58	1 (6%)	19,20,20	1.76	4 (21%)
2	NAD	B	501	-	46,48,48	1.32	6 (13%)	64,73,73	1.68	10 (15%)
5	EPE	A	504	-	15,15,15	1.69	1 (6%)	19,20,20	1.57	4 (21%)
2	NAD	C	501	-	46,48,48	1.43	6 (13%)	64,73,73	1.79	12 (18%)
4	PGE	D	503	-	9,9,9	0.62	0	8,8,8	0.74	0
4	PGE	C	504	-	9,9,9	0.52	0	8,8,8	0.52	0
5	EPE	C	505	-	15,15,15	1.87	1 (6%)	19,20,20	1.55	3 (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
5	EPE	B	505	-	-	7/9/19/19	0/1/1/1
2	NAD	D	501	-	-	7/30/62/62	0/5/5/5
4	PGE	A	503	-	-	5/7/7/7	-
4	PGE	B	503	-	-	4/7/7/7	-
5	EPE	A	505	-	-	3/9/19/19	0/1/1/1
2	NAD	A	501	-	-	7/30/62/62	0/5/5/5
4	PGE	C	503	-	-	3/7/7/7	-
5	EPE	B	504	-	-	1/9/19/19	0/1/1/1
2	NAD	B	501	-	-	7/30/62/62	0/5/5/5
5	EPE	A	504	-	-	5/9/19/19	0/1/1/1
2	NAD	C	501	-	-	9/30/62/62	0/5/5/5
4	PGE	D	503	-	-	4/7/7/7	-
4	PGE	C	504	-	-	3/7/7/7	-
5	EPE	C	505	-	-	4/9/19/19	0/1/1/1

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	505	EPE	C10-S	-6.79	1.68	1.77
5	C	505	EPE	C10-S	-6.77	1.68	1.77
5	B	505	EPE	C10-S	-6.64	1.68	1.77
2	A	501	NAD	O7N-C7N	6.19	1.35	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	504	EPE	C10-S	-6.16	1.69	1.77
2	D	501	NAD	O7N-C7N	6.07	1.35	1.24
2	B	501	NAD	O7N-C7N	5.89	1.35	1.24
5	B	504	EPE	C10-S	-5.50	1.69	1.77
2	C	501	NAD	O7N-C7N	5.08	1.33	1.24
2	C	501	NAD	PA-O3	4.29	1.64	1.59
2	D	501	NAD	C2N-N1N	3.59	1.38	1.35
2	A	501	NAD	C2N-N1N	3.52	1.38	1.35
2	C	501	NAD	C8A-N7A	3.13	1.37	1.31
2	A	501	NAD	C8A-N7A	2.81	1.37	1.31
2	C	501	NAD	PN-O3	2.74	1.62	1.59
2	A	501	NAD	PN-O3	2.64	1.62	1.59
2	A	501	NAD	O4D-C1D	2.62	1.44	1.40
2	A	501	NAD	C3N-C7N	2.55	1.54	1.50
2	D	501	NAD	C8A-N7A	2.55	1.36	1.31
2	B	501	NAD	C5A-N7A	-2.51	1.34	1.39
2	B	501	NAD	C8A-N7A	2.43	1.36	1.31
2	C	501	NAD	O4D-C1D	2.42	1.44	1.40
2	B	501	NAD	C4A-N9A	-2.40	1.32	1.37
2	D	501	NAD	C2N-C3N	2.31	1.42	1.39
2	D	501	NAD	PN-O3	2.26	1.61	1.59
2	A	501	NAD	C2N-C3N	2.24	1.42	1.39
2	B	501	NAD	O4D-C1D	2.21	1.43	1.40
2	D	501	NAD	O4D-C1D	2.17	1.43	1.40
2	A	501	NAD	PA-O3	2.10	1.61	1.59
2	D	501	NAD	C3N-C7N	2.08	1.53	1.50
2	B	501	NAD	C8A-N9A	-2.07	1.34	1.37
2	C	501	NAD	C4A-N9A	-2.05	1.33	1.37
2	A	501	NAD	C5A-N7A	-2.04	1.35	1.39

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	C5A-C4A-N3A	-5.69	118.88	126.72
2	C	501	NAD	C5A-C4A-N3A	-5.35	119.34	126.72
2	D	501	NAD	C5A-C4A-N3A	-5.01	119.82	126.72
5	B	504	EPE	O1S-S-C10	4.80	113.98	106.73
2	B	501	NAD	C5A-C4A-N3A	-4.60	120.38	126.72
2	A	501	NAD	C4A-C5A-N7A	-4.47	105.47	110.58
2	B	501	NAD	N3A-C2A-N1A	-4.47	121.82	128.58
2	C	501	NAD	N3A-C2A-N1A	-4.27	122.12	128.58
2	B	501	NAD	N9A-C8A-N7A	-4.26	107.90	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	C4A-C5A-N7A	-4.17	105.81	110.58
2	A	501	NAD	N3A-C2A-N1A	-4.12	122.34	128.58
2	C	501	NAD	C3N-C7N-N7N	4.10	122.78	117.74
2	D	501	NAD	N9A-C8A-N7A	-3.98	108.28	113.94
2	D	501	NAD	N3A-C2A-N1A	-3.94	122.62	128.58
2	A	501	NAD	C4D-O4D-C1D	3.92	113.52	109.92
2	C	501	NAD	C4A-C5A-N7A	-3.85	106.18	110.58
2	C	501	NAD	O7N-C7N-N7N	-3.84	117.06	122.62
5	A	504	EPE	O3S-S-C10	3.66	113.17	106.00
2	A	501	NAD	N9A-C8A-N7A	-3.61	108.81	113.94
2	C	501	NAD	N9A-C8A-N7A	-3.58	108.86	113.94
2	A	501	NAD	C2A-N3A-C4A	3.58	120.56	111.83
2	A	501	NAD	O7N-C7N-N7N	-3.54	117.49	122.62
2	C	501	NAD	C2A-N3A-C4A	3.53	120.45	111.83
2	A	501	NAD	C5A-N7A-C8A	3.52	108.98	103.45
2	C	501	NAD	C5A-C4A-N9A	3.44	109.56	105.81
2	A	501	NAD	C5A-C4A-N9A	3.43	109.55	105.81
2	B	501	NAD	C5A-N7A-C8A	3.43	108.83	103.45
2	D	501	NAD	N3A-C4A-N9A	3.29	132.77	127.17
2	D	501	NAD	C2A-N3A-C4A	3.13	119.48	111.83
5	C	505	EPE	C6-N1-C2	3.05	115.41	108.84
5	B	504	EPE	O2S-S-C10	-3.04	102.13	106.73
2	A	501	NAD	O4B-C1B-N9A	3.03	113.92	108.09
5	C	505	EPE	O3S-S-C10	3.00	111.88	106.00
2	B	501	NAD	C2A-N3A-C4A	2.97	119.09	111.83
2	D	501	NAD	C4A-N9A-C8A	2.96	108.85	105.74
2	B	501	NAD	C5A-C4A-N9A	2.91	108.98	105.81
2	D	501	NAD	O7N-C7N-N7N	-2.89	118.45	122.62
5	A	505	EPE	O1S-S-C10	2.88	111.08	106.73
5	B	504	EPE	C9-N1-C6	-2.82	103.72	111.24
2	C	501	NAD	C5A-N7A-C8A	2.80	107.85	103.45
2	D	501	NAD	C5A-N7A-C8A	2.75	107.78	103.45
5	C	505	EPE	O1S-S-C10	2.73	110.86	106.73
2	D	501	NAD	C4D-O4D-C1D	2.69	112.39	109.92
5	B	504	EPE	C7-N4-C5	-2.65	104.19	111.24
5	B	505	EPE	C7-N4-C3	2.59	118.14	111.24
2	B	501	NAD	C4A-N9A-C8A	2.58	108.45	105.74
2	A	501	NAD	N3A-C4A-N9A	2.58	131.55	127.17
2	D	501	NAD	C4A-C5A-N7A	-2.54	107.68	110.58
5	B	505	EPE	O3S-S-C10	2.51	110.91	106.00
2	A	501	NAD	O7N-C7N-C3N	2.31	122.42	119.60
2	A	501	NAD	C2N-C3N-C7N	2.31	126.12	119.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	O4B-C1B-C2B	-2.30	101.69	106.62
5	A	504	EPE	O2S-S-C10	2.30	110.20	106.73
5	B	505	EPE	C6-C5-N4	-2.27	106.07	110.65
2	C	501	NAD	O2A-PA-O1A	2.27	122.99	112.44
2	B	501	NAD	C5B-C4B-C3B	-2.26	107.07	115.21
2	C	501	NAD	N3A-C4A-N9A	2.25	130.99	127.17
5	A	505	EPE	O2S-S-C10	2.24	110.11	106.73
2	D	501	NAD	C5B-C4B-C3B	-2.19	107.31	115.21
5	B	505	EPE	O2S-S-C10	2.16	110.00	106.73
2	D	501	NAD	C3N-C7N-N7N	2.14	120.38	117.74
2	D	501	NAD	C2N-C3N-C7N	2.09	125.49	119.46
5	A	504	EPE	C6-N1-C2	2.05	113.26	108.84
2	B	501	NAD	N3A-C4A-N9A	2.04	130.63	127.17
2	C	501	NAD	C5B-C4B-C3B	-2.03	107.91	115.21
5	A	504	EPE	C7-N4-C5	2.01	116.59	111.24

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAD	C5D-O5D-PN-O3
2	A	501	NAD	C5D-O5D-PN-O1N
2	A	501	NAD	C5D-O5D-PN-O2N
2	B	501	NAD	C5D-O5D-PN-O3
2	B	501	NAD	C5D-O5D-PN-O1N
2	B	501	NAD	C5D-O5D-PN-O2N
2	C	501	NAD	C5D-O5D-PN-O3
2	C	501	NAD	C5D-O5D-PN-O1N
2	C	501	NAD	C5D-O5D-PN-O2N
2	D	501	NAD	C5D-O5D-PN-O3
5	A	504	EPE	C10-C9-N1-C2
5	A	504	EPE	C10-C9-N1-C6
5	A	505	EPE	C9-C10-S-O2S
5	A	505	EPE	C9-C10-S-O3S
5	B	505	EPE	C10-C9-N1-C2
5	B	505	EPE	C10-C9-N1-C6
5	B	505	EPE	C8-C7-N4-C3
5	B	505	EPE	C9-C10-S-O1S
5	B	505	EPE	C9-C10-S-O2S
5	C	505	EPE	C10-C9-N1-C6
4	D	503	PGE	C6-C5-O3-C4
4	B	503	PGE	C6-C5-O3-C4

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Mol	Chain	Res	Type	Atoms
5	A	504	EPE	C8-C7-N4-C5
5	C	505	EPE	C8-C7-N4-C5
4	A	503	PGE	O2-C3-C4-O3
4	D	503	PGE	O2-C3-C4-O3
2	A	501	NAD	O4B-C4B-C5B-O5B
2	A	501	NAD	C3B-C4B-C5B-O5B
4	C	503	PGE	O1-C1-C2-O2
4	A	503	PGE	O1-C1-C2-O2
4	C	504	PGE	O3-C5-C6-O4
5	A	504	EPE	N4-C7-C8-O8
5	C	505	EPE	N4-C7-C8-O8
4	D	503	PGE	O1-C1-C2-O2
4	C	503	PGE	O2-C3-C4-O3
4	C	504	PGE	O2-C3-C4-O3
5	B	505	EPE	C9-C10-S-O3S
4	A	503	PGE	O3-C5-C6-O4
5	C	505	EPE	C8-C7-N4-C3
2	D	501	NAD	O4D-C4D-C5D-O5D
4	B	503	PGE	O2-C3-C4-O3
4	A	503	PGE	C6-C5-O3-C4
5	A	505	EPE	C9-C10-S-O1S
5	B	504	EPE	S-C10-C9-N1
4	C	504	PGE	C3-C4-O3-C5
2	C	501	NAD	C4N-C3N-C7N-N7N
4	B	503	PGE	C4-C3-O2-C2
4	C	503	PGE	O3-C5-C6-O4
2	C	501	NAD	C2N-C3N-C7N-N7N
4	A	503	PGE	C4-C3-O2-C2
2	B	501	NAD	C3B-C4B-C5B-O5B
2	A	501	NAD	C5B-O5B-PA-O1A
2	B	501	NAD	C5B-O5B-PA-O1A
2	D	501	NAD	C5B-O5B-PA-O1A
2	D	501	NAD	C5D-O5D-PN-O1N
4	D	503	PGE	C4-C3-O2-C2
5	B	505	EPE	C8-C7-N4-C5
4	B	503	PGE	C1-C2-O2-C3
2	C	501	NAD	C4N-C3N-C7N-O7N
2	C	501	NAD	PA-O3-PN-O2N
5	A	504	EPE	C9-C10-S-O2S
2	C	501	NAD	C2N-C3N-C7N-O7N
2	B	501	NAD	O4B-C4B-C5B-O5B
2	A	501	NAD	C2B-C1B-N9A-C8A

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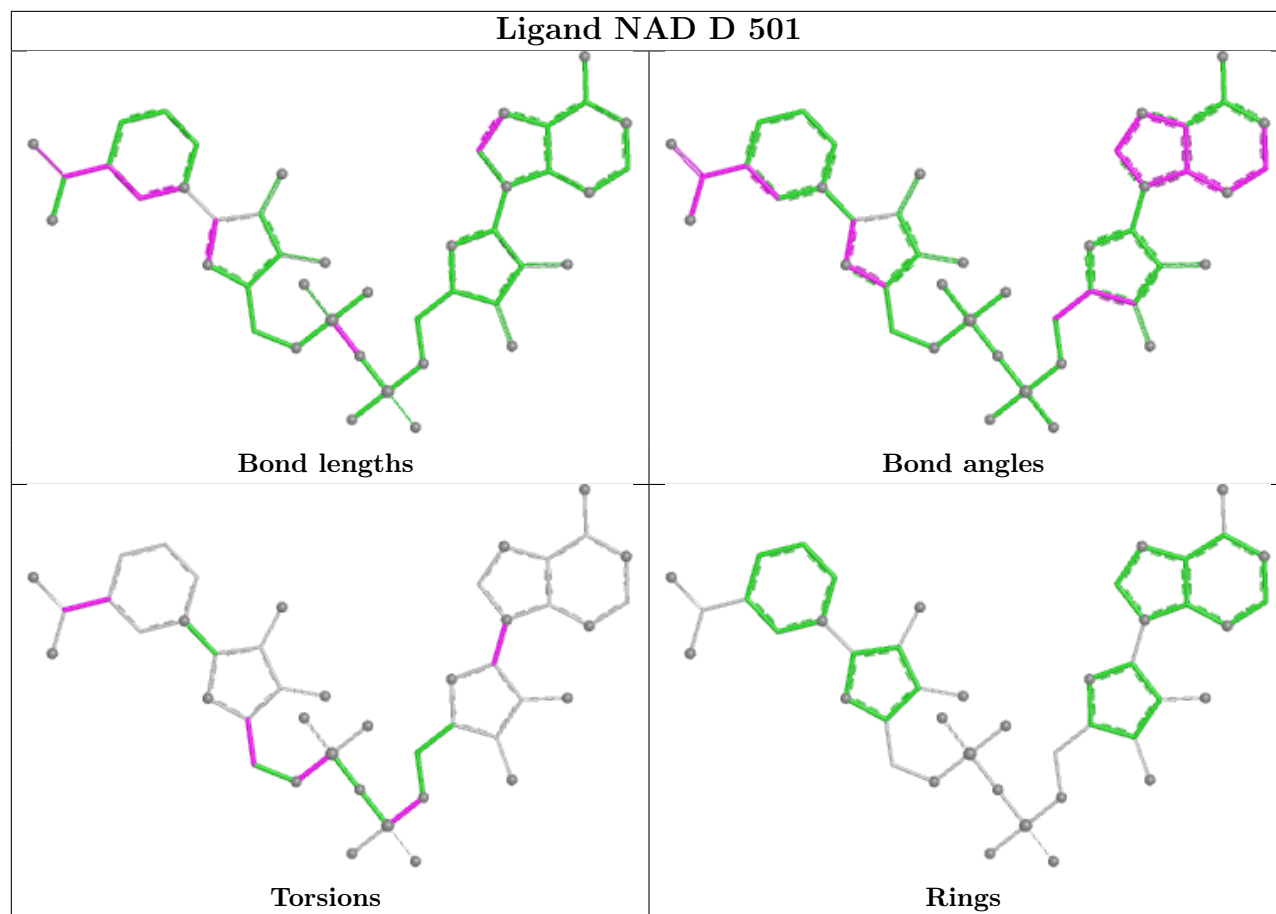
Mol	Chain	Res	Type	Atoms
2	D	501	NAD	C4N-C3N-C7N-N7N
2	C	501	NAD	C2B-C1B-N9A-C8A
2	B	501	NAD	C2B-C1B-N9A-C8A
2	D	501	NAD	C2B-C1B-N9A-C8A
2	D	501	NAD	C3D-C4D-C5D-O5D

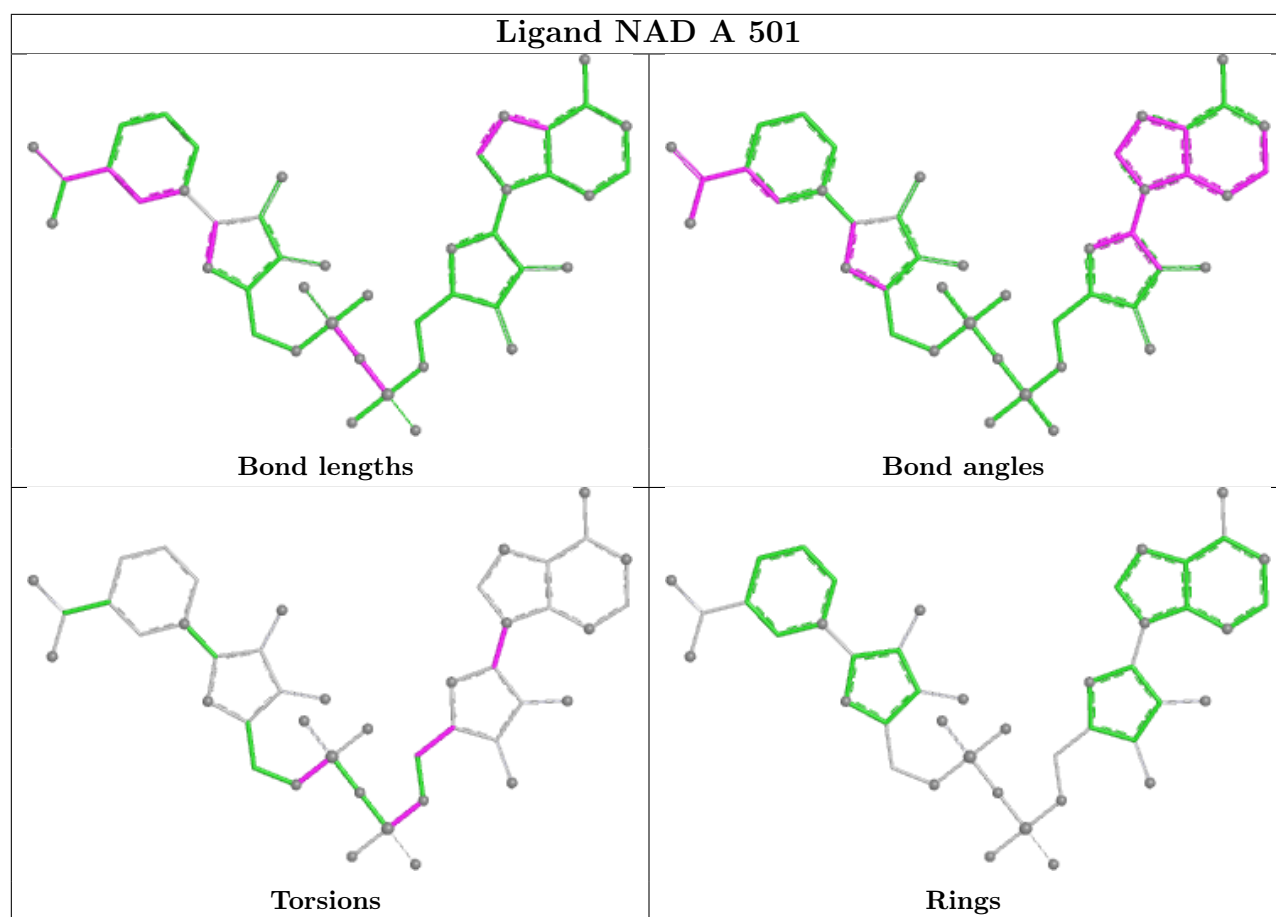
There are no ring outliers.

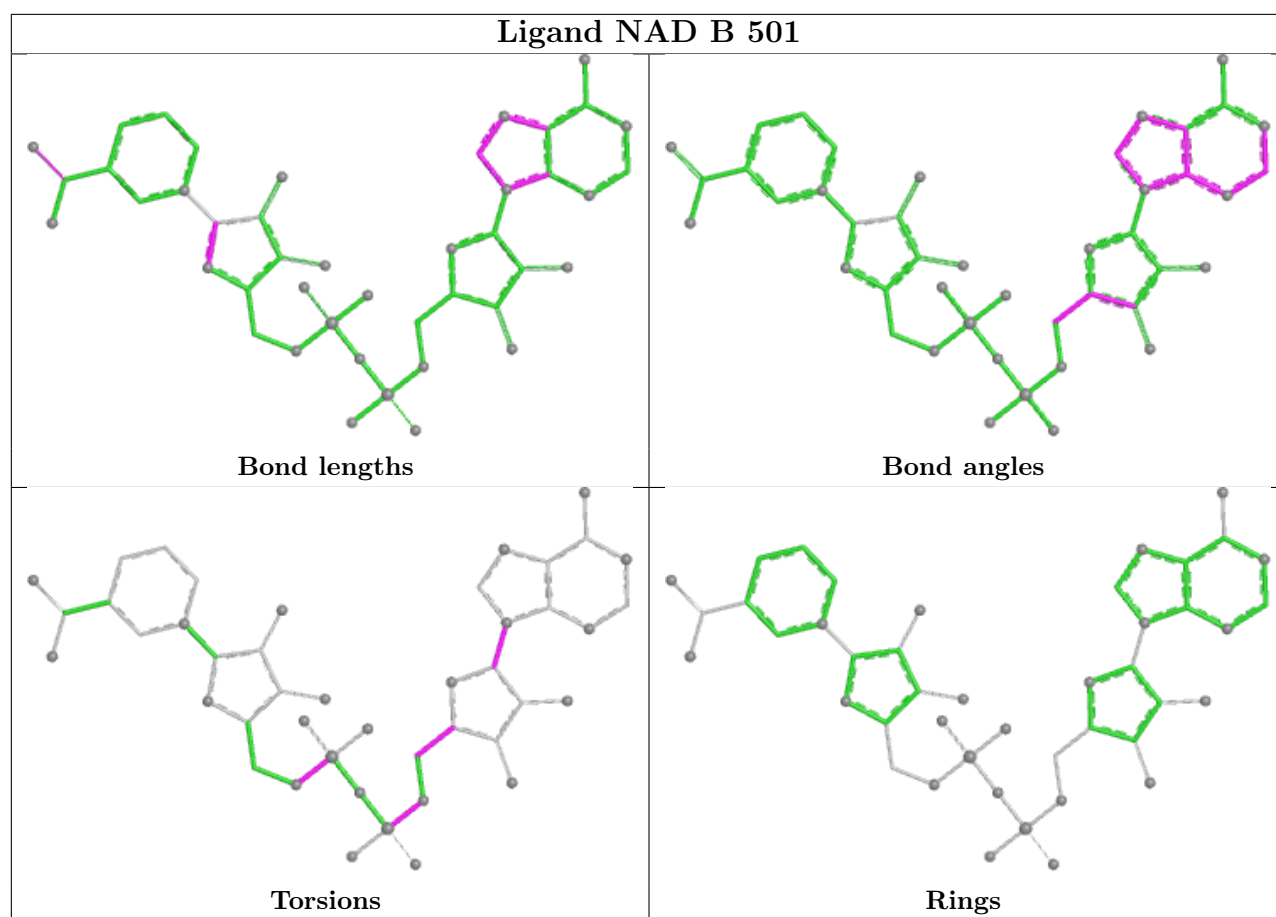
9 monomers are involved in 20 short contacts:

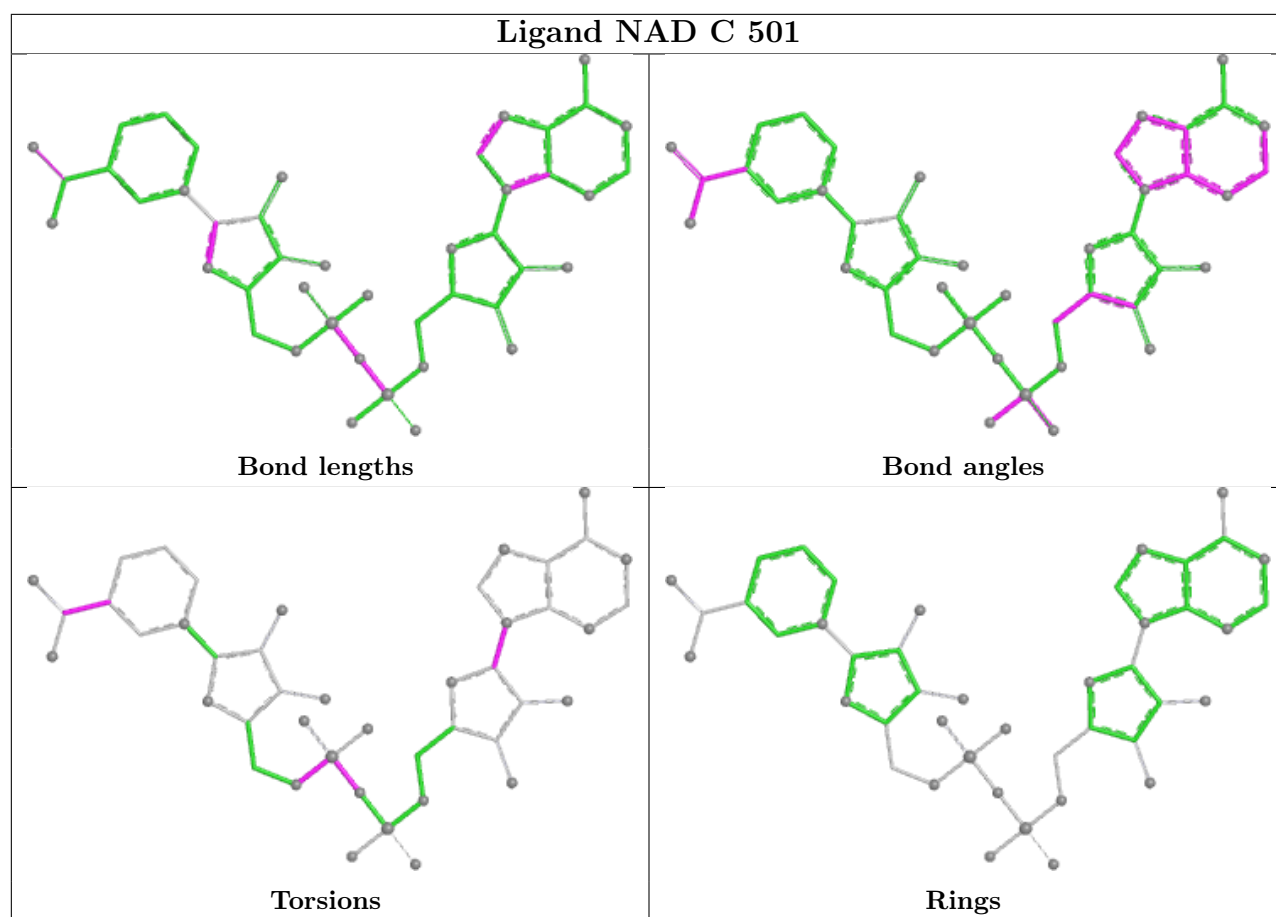
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	505	EPE	1	0
2	D	501	NAD	3	0
4	A	503	PGE	1	0
5	A	505	EPE	1	0
2	A	501	NAD	3	0
2	B	501	NAD	4	0
5	A	504	EPE	1	0
2	C	501	NAD	4	0
5	C	505	EPE	2	0

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	495/517 (95%)	-0.57	3 (0%) 85 87	7, 18, 36, 88	16 (3%)
1	B	494/517 (95%)	-0.52	3 (0%) 85 87	7, 19, 39, 85	15 (3%)
1	C	497/517 (96%)	-0.55	10 (2%) 65 68	7, 20, 35, 97	13 (2%)
1	D	492/517 (95%)	-0.60	1 (0%) 91 92	7, 19, 34, 72	12 (2%)
All	All	1978/2068 (95%)	-0.56	17 (0%) 81 83	7, 19, 36, 97	56 (2%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	LEU	5.6
1	C	3	LEU	4.9
1	B	3	LEU	4.6
1	C	4	LEU	4.1
1	A	1	MET	4.0
1	C	6	HIS	3.8
1	C	0	GLY	3.5
1	C	1	MET	3.1
1	C	-1	GLN	2.7
1	D	367	LYS	2.6
1	B	364	ASP	2.4
1	C	5	LYS	2.4
1	A	364	ASP	2.2
1	C	362	ASP	2.1
1	C	364	ASP	2.1
1	C	2	GLU	2.1
1	B	348	LYS	2.1

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

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
1	CME	C	289	10/11	0.90	0.14	18,24,49,63	0
1	CME	A	289	10/11	0.94	0.11	18,25,52,63	0
1	CME	D	289	10/11	0.94	0.12	19,24,49,63	0
1	CME	B	289	10/11	0.95	0.10	20,24,52,60	0

6.3 Carbohydrates ⓘ

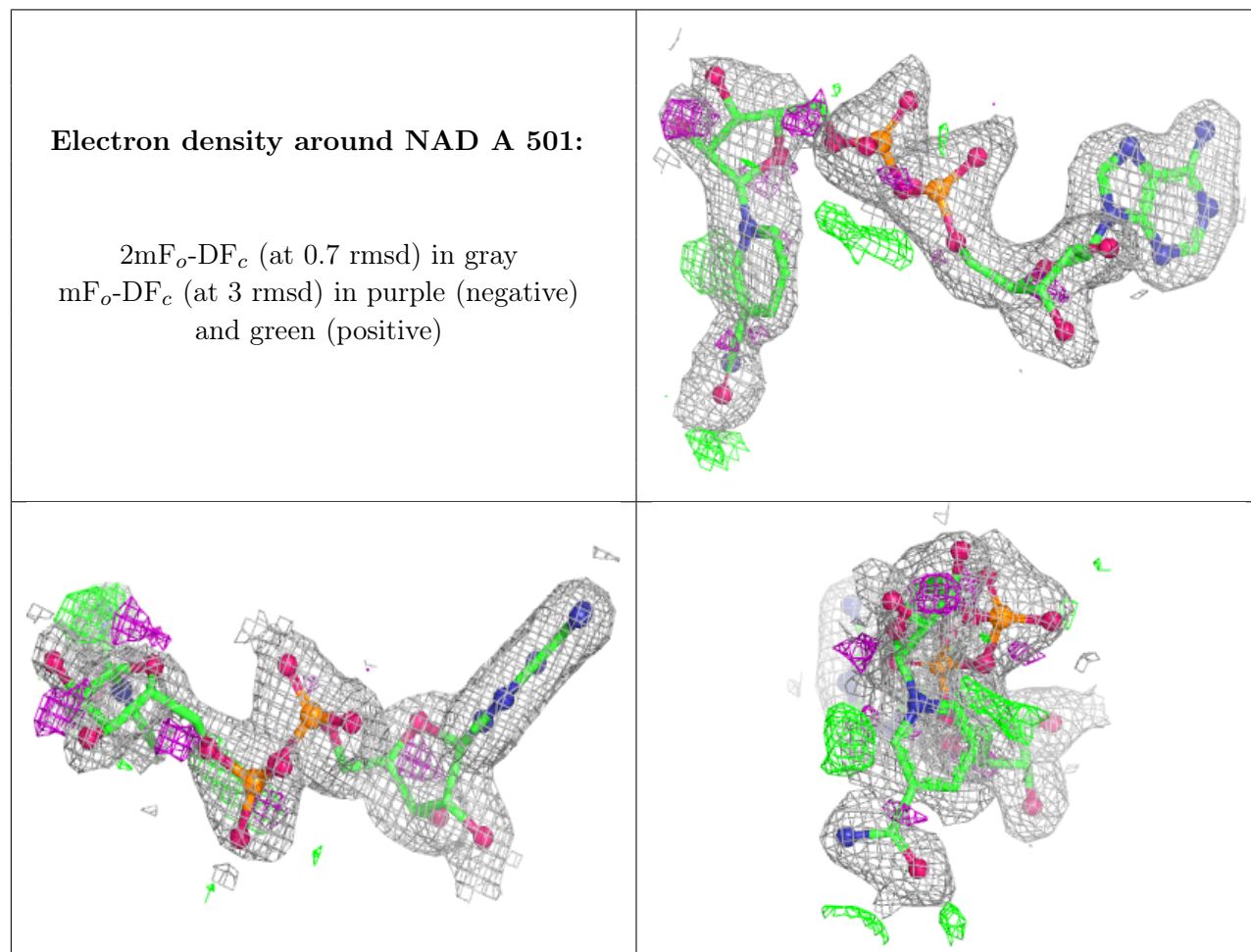
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

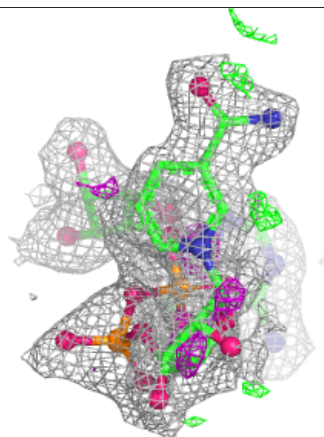
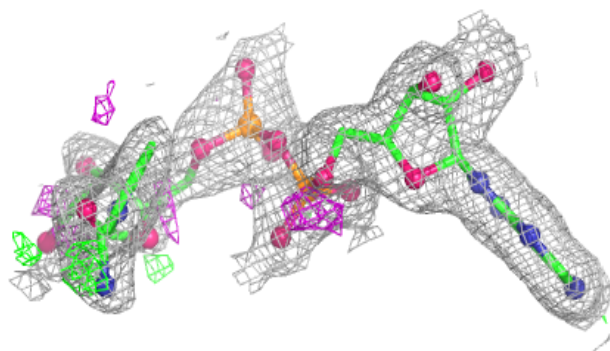
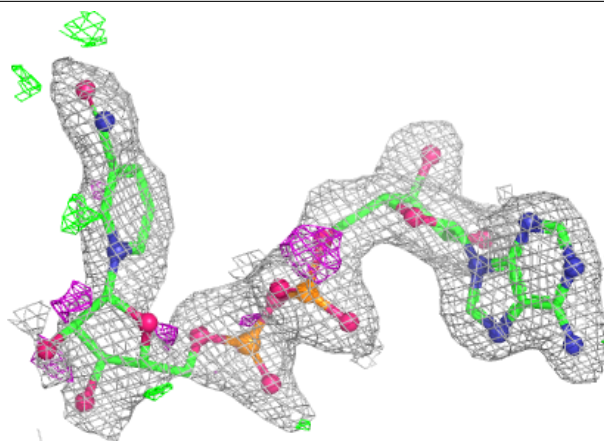
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	EPE	A	505	15/15	0.68	0.20	55,69,110,120	0
5	EPE	B	505	15/15	0.73	0.19	51,65,124,130	0
5	EPE	C	505	15/15	0.74	0.19	52,65,114,128	0
4	PGE	C	504	10/10	0.82	0.14	43,47,58,59	0
4	PGE	C	503	10/10	0.84	0.15	46,53,61,63	0
4	PGE	D	503	10/10	0.87	0.13	40,49,52,54	0
4	PGE	B	503	10/10	0.87	0.13	39,47,51,52	0
4	PGE	A	503	10/10	0.88	0.12	29,42,46,51	0
5	EPE	B	504	15/15	0.90	0.11	28,32,52,57	0
2	NAD	A	501	44/44	0.93	0.10	22,35,56,66	0
5	EPE	A	504	15/15	0.93	0.10	32,39,53,59	0
2	NAD	B	501	44/44	0.94	0.09	18,35,58,70	0
2	NAD	C	501	44/44	0.95	0.08	18,35,52,57	0
2	NAD	D	501	44/44	0.95	0.08	18,34,55,64	0
3	NA	B	502	1/1	0.96	0.07	15,15,15,15	0
3	NA	A	502	1/1	0.98	0.04	15,15,15,15	0
3	NA	C	502	1/1	0.99	0.03	15,15,15,15	0
3	NA	D	502	1/1	0.99	0.07	13,13,13,13	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.



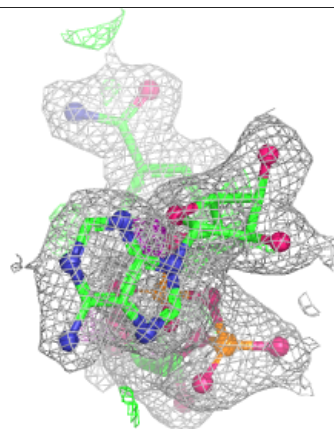
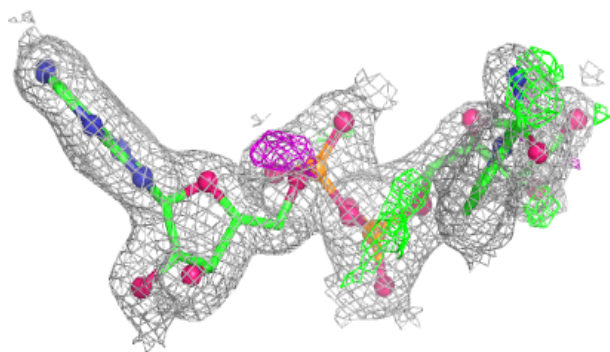
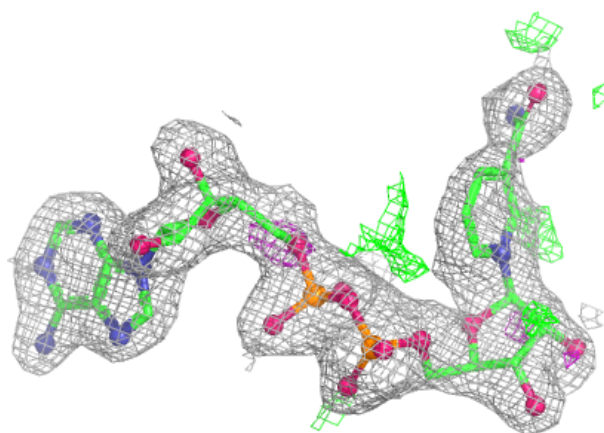
Electron density around NAD B 501:

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



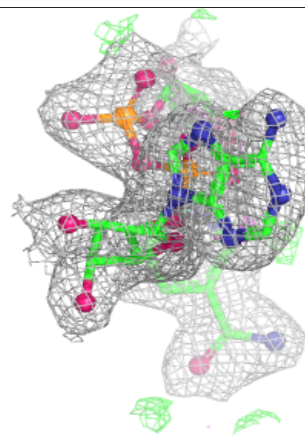
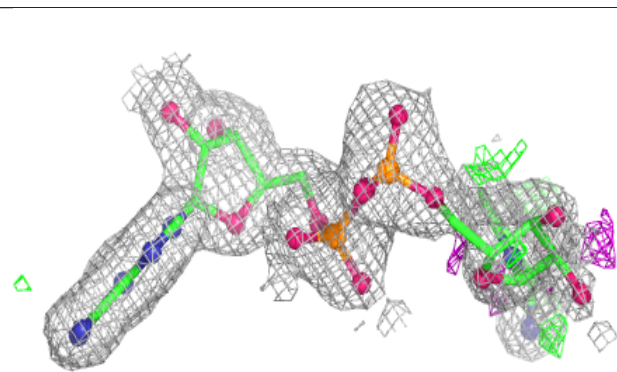
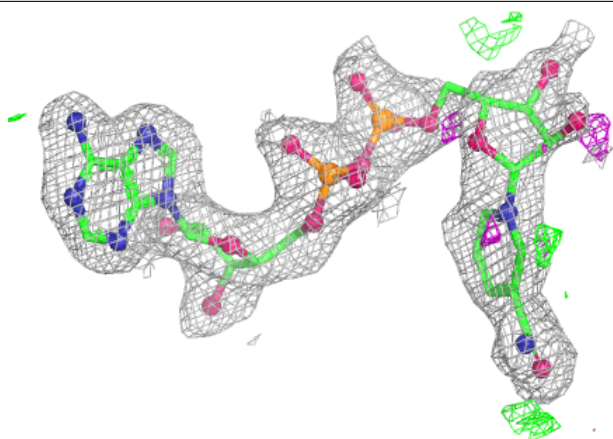
Electron density around NAD C 501:

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



Electron density around NAD 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.