



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 23, 2026 – 02:08 PM EDT

PDB ID : 9PXH / pdb_00009pxh
Title : Crystal structure of bovine ALDH3A1 in complex with Non-Canonical Redox Cofactors, NMN+
Authors : Hsu, N.; Saleh, S.; Kim, J.; Li, H.; Qiao, F.
Deposited on : 2025-08-05
Resolution : 1.77 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

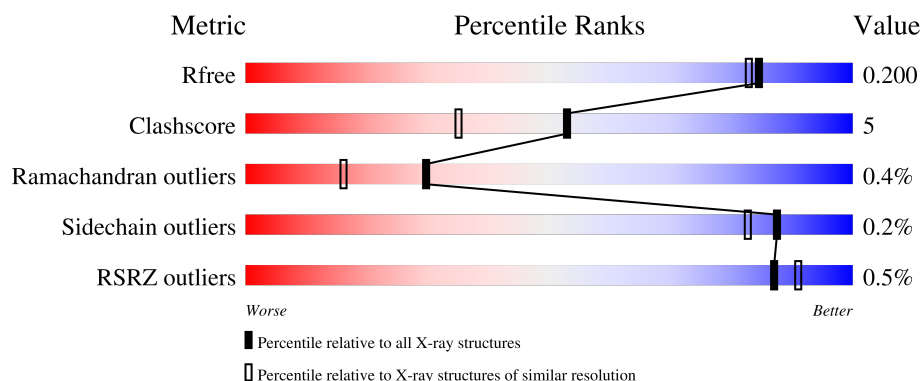
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	452	90% 9%
1	B	452	92% 8%
1	C	452	% 92% 7%
1	D	452	% 91% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	FMT	A	508	-	-	X	-
4	FMT	A	509	-	-	X	-
4	FMT	C	504	-	-	X	-
4	FMT	C	505	-	-	X	-
4	FMT	C	506	-	-	X	-
4	FMT	C	507	-	-	X	-
4	FMT	D	509	-	-	X	-
6	EDO	A	510	-	-	X	-

2 Entry composition [i](#)

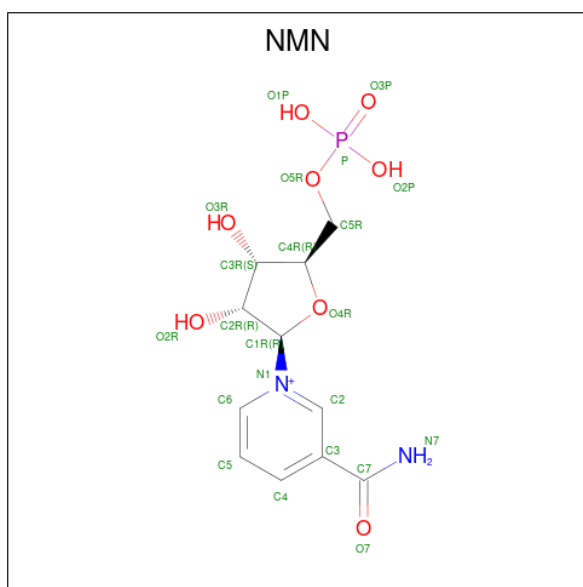
There are 8 unique types of molecules in this entry. The entry contains 30523 atoms, of which 14040 are hydrogens and 0 are deuteriums.

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 Aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	450	Total	C	H	N	O	S	0	1	0
			6981	2219	3483	602	659	18			
1	B	450	Total	C	H	N	O	S	0	1	0
			6976	2218	3484	600	656	18			
1	C	450	Total	C	H	N	O	S	0	2	0
			6950	2215	3458	599	660	18			
1	D	446	Total	C	H	N	O	S	0	0	0
			6876	2192	3422	591	654	17			

- Molecule 2 is BETA-NICOTINAMIDE RIBOSE MONOPHOSPHATE (CCD ID: NMN) (formula: C₁₁H₁₆N₂O₈P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			36	11	14	2	8	1		
2	B	1	Total	C	H	N	O	P	0	0
			36	11	14	2	8	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	H	N	O	P	
			36	11	14	2	8	1	0
2	D	1	Total	C	H	N	O	P	
			36	11	14	2	8	1	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



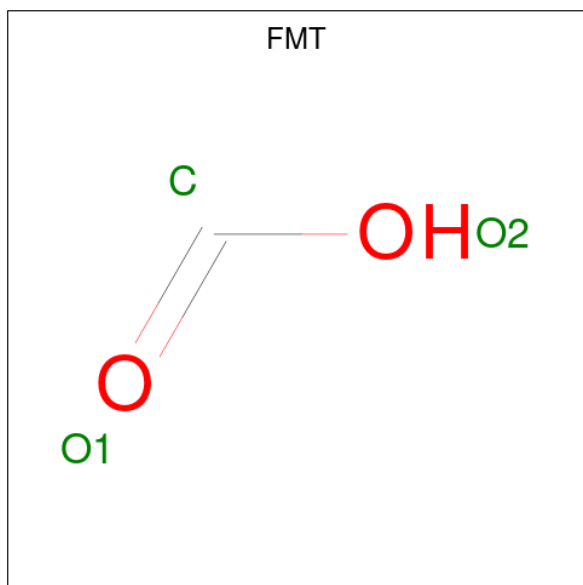
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	D	1	Total	C	H	O	0	0
			14	3	8	3		
3	D	1	Total	C	H	O	0	0
			14	3	8	3		
3	D	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 4 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



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

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

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			24	6	14	4		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 7 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total 1	Mg 1	0	0
7	D	1	Total 1	Mg 1	0	0

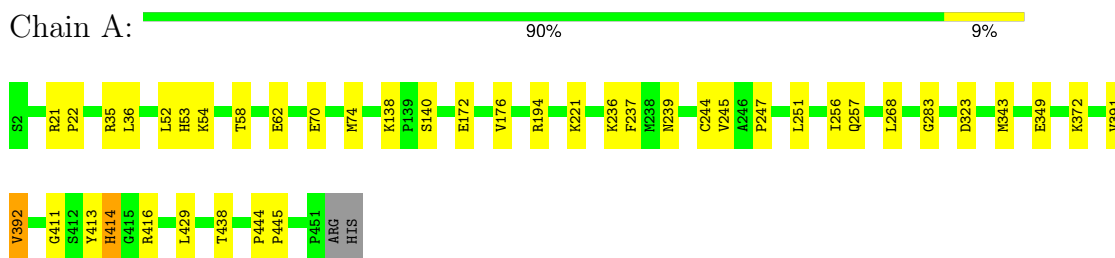
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	623	Total 623	O 623	0	0
8	B	616	Total 616	O 616	0	0
8	C	543	Total 543	O 543	0	0
8	D	526	Total 526	O 526	0	0

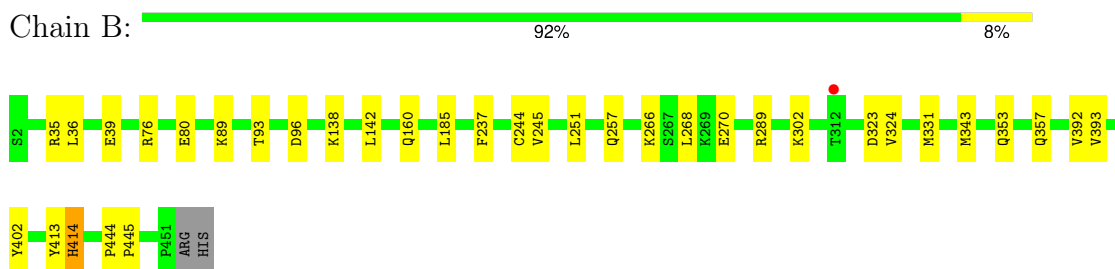
3 Residue-property plots

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.

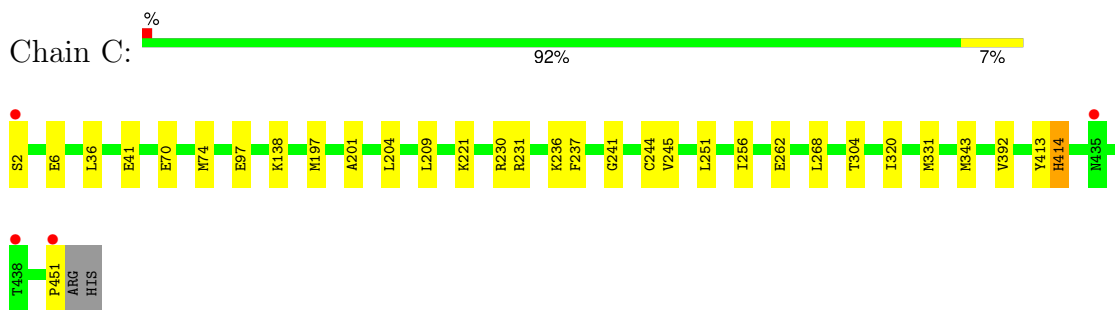
- Molecule 1: Aldehyde dehydrogenase



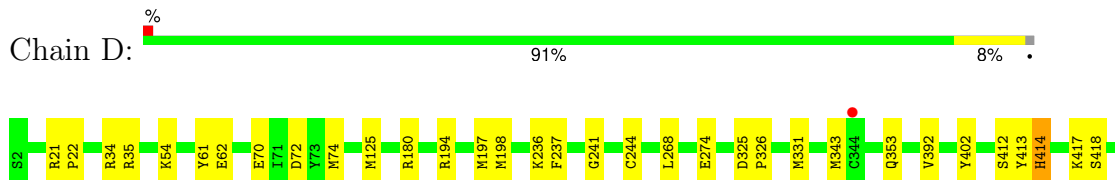
- Molecule 1: Aldehyde dehydrogenase



- Molecule 1: Aldehyde dehydrogenase



- Molecule 1: Aldehyde dehydrogenase



L434				
E435				
E436				
E437				
F447				
ALA				
LYS				
MET				
PRO				
ARG				
HIS				

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.17Å 59.27Å 158.64Å 90.00° 98.43° 90.00°	Depositor
Resolution (Å)	43.00 – 1.77 43.00 – 1.77	Depositor EDS
% Data completeness (in resolution range)	96.9 (43.00-1.77) 96.9 (43.00-1.77)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 1.77Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.153 , 0.200 0.153 , 0.200	Depositor DCC
R_{free} test set	2000 reflections (0.81%)	wwPDB-VP
Wilson B-factor (Å ²)	11.5	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	30523	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.02% 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: GOL, FMT, NMN, EDO, MG, PGE

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	0.48	0/3573	0.64	0/4854
1	B	0.48	0/3567	0.65	1/4845 (0.0%)
1	C	0.44	0/3567	0.62	0/4848
1	D	0.42	0/3528	0.62	2/4796 (0.0%)
All	All	0.45	0/14235	0.63	3/19343 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	125	MET	CB-CA-C	-5.48	102.05	110.81
1	D	331	MET	CA-CB-CG	-5.29	103.52	114.10
1	B	331	MET	CA-CB-CG	-5.14	103.82	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	35	ARG	Sidechain

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	3498	3483	3482	35	1
1	B	3492	3484	3483	25	0
1	C	3492	3458	3456	25	0
1	D	3454	3422	3422	35	0
2	A	22	14	13	4	0
2	B	22	14	12	5	0
2	C	22	14	13	4	0
2	D	22	14	12	5	0
3	A	18	24	24	1	0
3	B	24	32	32	2	0
3	C	12	16	16	0	0
3	D	18	24	24	1	0
4	A	18	6	6	11	0
4	B	6	2	2	0	0
4	C	24	8	8	7	0
4	D	15	5	5	5	0
5	A	10	14	14	0	0
6	A	4	6	6	9	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	A	623	0	0	12	6
8	B	616	0	0	12	2
8	C	543	0	0	11	4
8	D	526	0	0	15	4
All	All	16483	14040	14030	140	13

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 5.

The worst 5 of 140 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)
2:B:502:NMN:C4R	2:B:502:NMN:O4R	1.64	1.20
2:C:501:NMN:C4R	2:C:501:NMN:O4R	1.64	1.14
2:A:501:NMN:O4R	2:A:501:NMN:C4R	1.65	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:502:NMN:O1P	8:B:601:HOH:O	1.92	0.88
4:C:505:FMT:O1	8:C:601:HOH:O	1.92	0.86

The worst 5 of 13 symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:B:1169:HOH:O	8:C:919:HOH:O[2_555]	1.80	0.40
8:D:609:HOH:O	8:D:719:HOH:O[1_565]	1.99	0.21
8:C:1123:HOH:O	8:C:1129:HOH:O[2_546]	2.04	0.16
8:C:1046:HOH:O	8:D:950:HOH:O[1_565]	2.06	0.14
8:A:1165:HOH:O	8:D:1072:HOH:O[1_655]	2.07	0.13

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	449/452 (99%)	442 (98%)	5 (1%)	2 (0%)	30	16
1	B	449/452 (99%)	442 (98%)	5 (1%)	2 (0%)	30	16
1	C	450/452 (100%)	443 (98%)	5 (1%)	2 (0%)	30	16
1	D	444/452 (98%)	436 (98%)	6 (1%)	2 (0%)	24	11
All	All	1792/1808 (99%)	1763 (98%)	21 (1%)	8 (0%)	30	16

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	392	VAL
1	A	414	HIS
1	B	414	HIS
1	C	414	HIS
1	D	414	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	379/387 (98%)	379 (100%)	0	100	100
1	B	378/387 (98%)	376 (100%)	2 (0%)	81	73
1	C	376/387 (97%)	375 (100%)	1 (0%)	86	80
1	D	373/387 (96%)	373 (100%)	0	100	100
All	All	1506/1548 (97%)	1503 (100%)	3 (0%)	87	83

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	138	LYS
1	B	142	LEU
1	C	138	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	257	GLN
1	B	259	GLN
1	C	94	GLN
1	B	435	ASN
1	B	160	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 41 ligands modelled in this entry, 2 are monoatomic - leaving 39 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$
3	GOL	B	505	-	5,5,5	0.44	0	5,5,5	1.33	0
3	GOL	B	503	-	5,5,5	0.50	0	5,5,5	0.82	0
4	FMT	C	507	-	2,2,2	0.77	0	1,1,1	0.60	0
4	FMT	C	505	-	2,2,2	0.69	0	1,1,1	0.65	0
4	FMT	A	512	-	2,2,2	0.89	0	1,1,1	0.33	0
4	FMT	A	508	-	2,2,2	0.65	0	1,1,1	0.46	0
3	GOL	B	501	-	5,5,5	0.34	0	5,5,5	0.38	0
4	FMT	D	506	-	2,2,2	0.63	0	1,1,1	0.68	0
4	FMT	C	506	-	2,2,2	0.58	0	1,1,1	0.58	0
4	FMT	D	501	-	2,2,2	0.87	0	1,1,1	0.38	0
6	EDO	A	510	-	3,3,3	0.34	0	2,2,2	1.18	0
4	FMT	A	504	-	2,2,2	0.76	0	1,1,1	0.58	0
3	GOL	D	503	-	5,5,5	0.53	0	5,5,5	0.59	0
4	FMT	C	508	-	2,2,2	0.64	0	1,1,1	0.68	0
3	GOL	C	503	-	5,5,5	0.51	0	5,5,5	0.51	0
4	FMT	D	508	-	2,2,2	0.74	0	1,1,1	0.53	0
4	FMT	A	505	-	2,2,2	0.68	0	1,1,1	0.46	0
3	GOL	A	503	-	5,5,5	0.53	0	5,5,5	0.76	0
4	FMT	D	507	-	2,2,2	0.66	0	1,1,1	0.47	0
3	GOL	A	502	-	5,5,5	0.82	0	5,5,5	0.92	0
4	FMT	A	507	-	2,2,2	0.70	0	1,1,1	0.49	0
5	PGE	A	506	-	9,9,9	0.54	0	8,8,8	0.49	0
2	NMN	B	502	-	21,23,23	3.96	8 (38%)	27,34,34	2.76	9 (33%)
4	FMT	C	504	-	2,2,2	0.54	0	1,1,1	0.65	0
4	FMT	B	507	-	2,2,2	0.72	0	1,1,1	0.56	0
4	FMT	A	509	-	2,2,2	0.39	0	1,1,1	0.25	0
2	NMN	D	502	-	21,23,23	4.01	10 (47%)	27,34,34	1.70	2 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FMT	D	509	-	2,2,2	0.71	0	1,1,1	0.72	0
3	GOL	A	511	-	5,5,5	0.53	0	5,5,5	0.35	0
3	GOL	C	502	-	5,5,5	0.51	0	5,5,5	0.51	0
3	GOL	B	504	-	5,5,5	0.77	0	5,5,5	0.71	0
4	FMT	C	510	-	2,2,2	0.70	0	1,1,1	0.53	0
4	FMT	B	506	-	2,2,2	0.80	0	1,1,1	0.51	0
2	NMN	C	501	-	21,23,23	4.14	8 (38%)	27,34,34	1.99	3 (11%)
4	FMT	C	509	-	2,2,2	0.69	0	1,1,1	0.39	0
4	FMT	C	511	-	2,2,2	0.84	0	1,1,1	0.37	0
2	NMN	A	501	-	21,23,23	3.98	7 (33%)	27,34,34	2.08	5 (18%)
3	GOL	D	505	-	5,5,5	0.54	0	5,5,5	0.45	0
3	GOL	D	504	-	5,5,5	0.49	0	5,5,5	0.42	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	505	-	-	2/4/4/4	-
3	GOL	B	503	-	-	0/4/4/4	-
2	NMN	D	502	-	-	6/14/30/30	0/2/2/2
3	GOL	A	503	-	-	0/4/4/4	-
3	GOL	A	511	-	-	0/4/4/4	-
3	GOL	D	503	-	-	0/4/4/4	-
3	GOL	C	502	-	-	0/4/4/4	-
3	GOL	A	502	-	-	0/4/4/4	-
3	GOL	B	504	-	-	0/4/4/4	-
3	GOL	C	503	-	-	0/4/4/4	-
5	PGE	A	506	-	-	4/7/7/7	-
2	NMN	B	502	-	-	4/14/30/30	0/2/2/2
2	NMN	C	501	-	-	1/14/30/30	0/2/2/2
3	GOL	D	504	-	-	1/4/4/4	-
2	NMN	A	501	-	-	1/14/30/30	0/2/2/2
3	GOL	D	505	-	-	0/4/4/4	-
6	EDO	A	510	-	-	0/1/1/1	-
3	GOL	B	501	-	-	0/4/4/4	-

The worst 5 of 33 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	NMN	C3R-C4R	-11.31	1.24	1.53
2	D	502	NMN	C3R-C4R	-10.73	1.25	1.53
2	B	502	NMN	C3R-C4R	-10.72	1.25	1.53
2	A	501	NMN	C3R-C4R	-10.36	1.26	1.53
2	A	501	NMN	O4R-C4R	9.27	1.65	1.45

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	NMN	C4R-O4R-C1R	-10.43	100.38	109.92
2	A	501	NMN	C4R-O4R-C1R	-7.86	102.72	109.92
2	C	501	NMN	C4R-O4R-C1R	-7.80	102.78	109.92
2	D	502	NMN	C4R-O4R-C1R	-6.82	103.68	109.92
2	B	502	NMN	O7-C7-C3	5.59	126.44	119.60

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	502	NMN	C5R-O5R-P-O3P
2	D	502	NMN	C5R-O5R-P-O1P
2	D	502	NMN	C5R-O5R-P-O2P
3	B	505	GOL	C1-C2-C3-O3
5	A	506	PGE	O2-C3-C4-O3

There are no ring outliers.

21 monomers are involved in 51 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	505	GOL	1	0
4	C	507	FMT	2	0
4	C	505	FMT	2	0
4	A	508	FMT	2	0
3	B	501	GOL	1	0
4	D	506	FMT	1	0
4	C	506	FMT	2	0
6	A	510	EDO	9	0
4	A	504	FMT	1	0
3	D	503	GOL	1	0
4	C	508	FMT	1	0
4	A	505	FMT	1	0
3	A	503	GOL	1	0

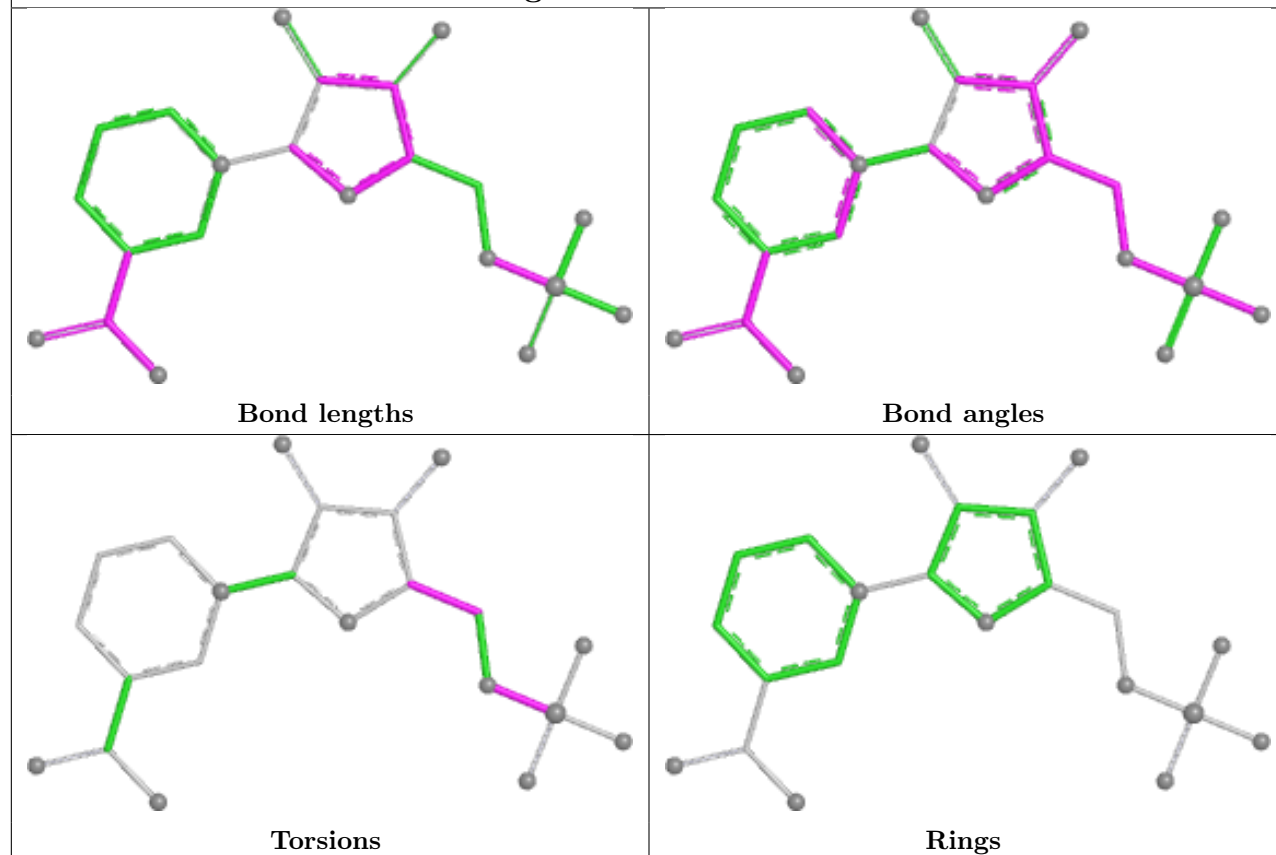
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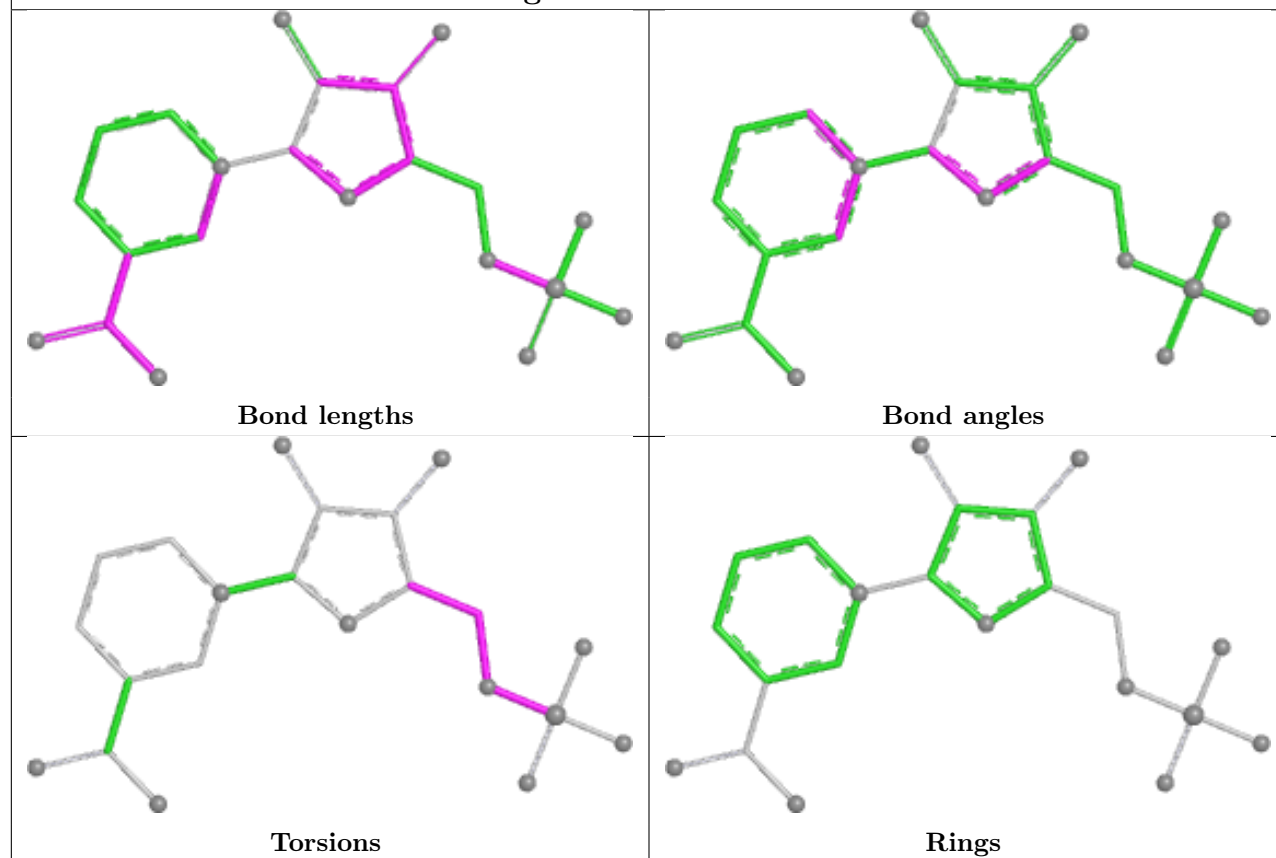
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	507	FMT	1	0
2	B	502	NMN	5	0
4	C	504	FMT	2	0
4	A	509	FMT	7	0
2	D	502	NMN	5	0
4	D	509	FMT	4	0
2	C	501	NMN	4	0
2	A	501	NMN	4	0

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.

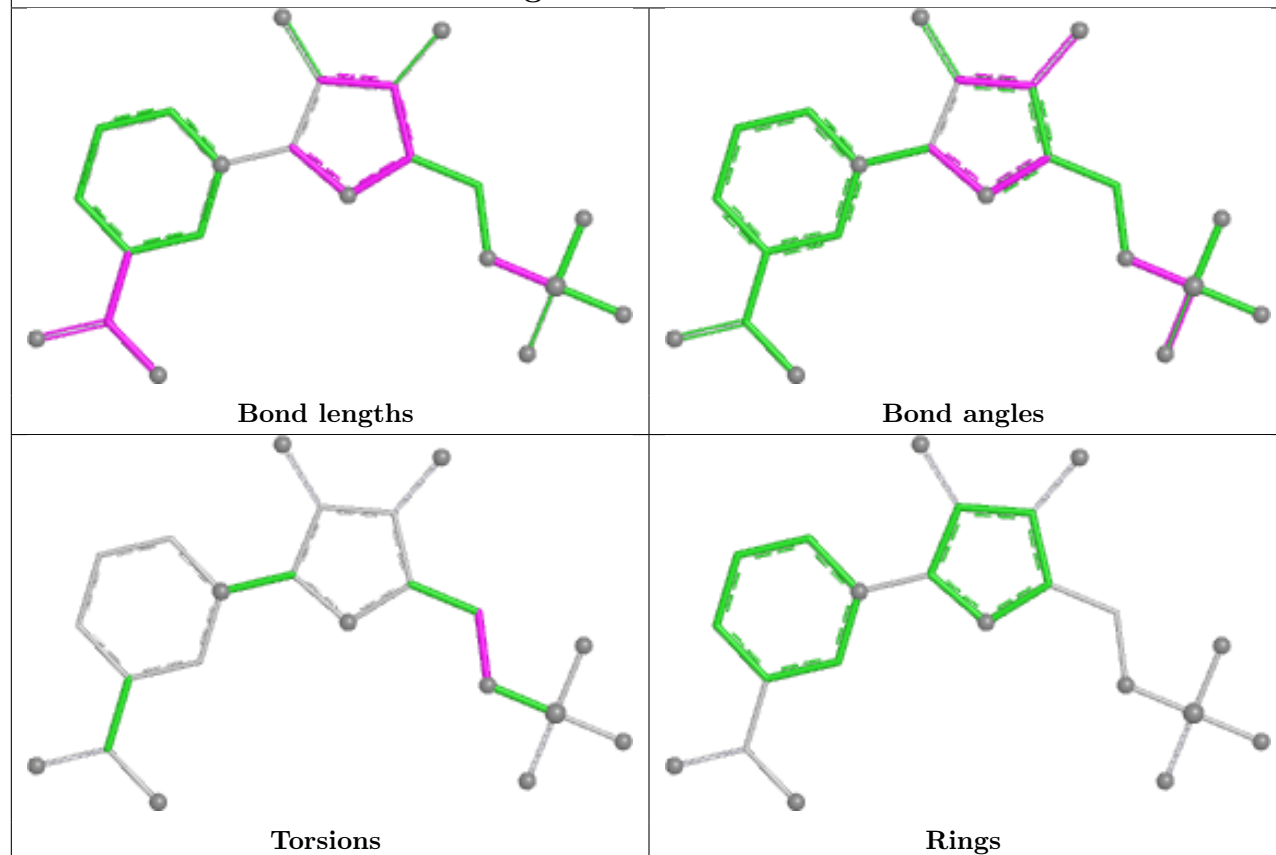
Ligand NMN B 502



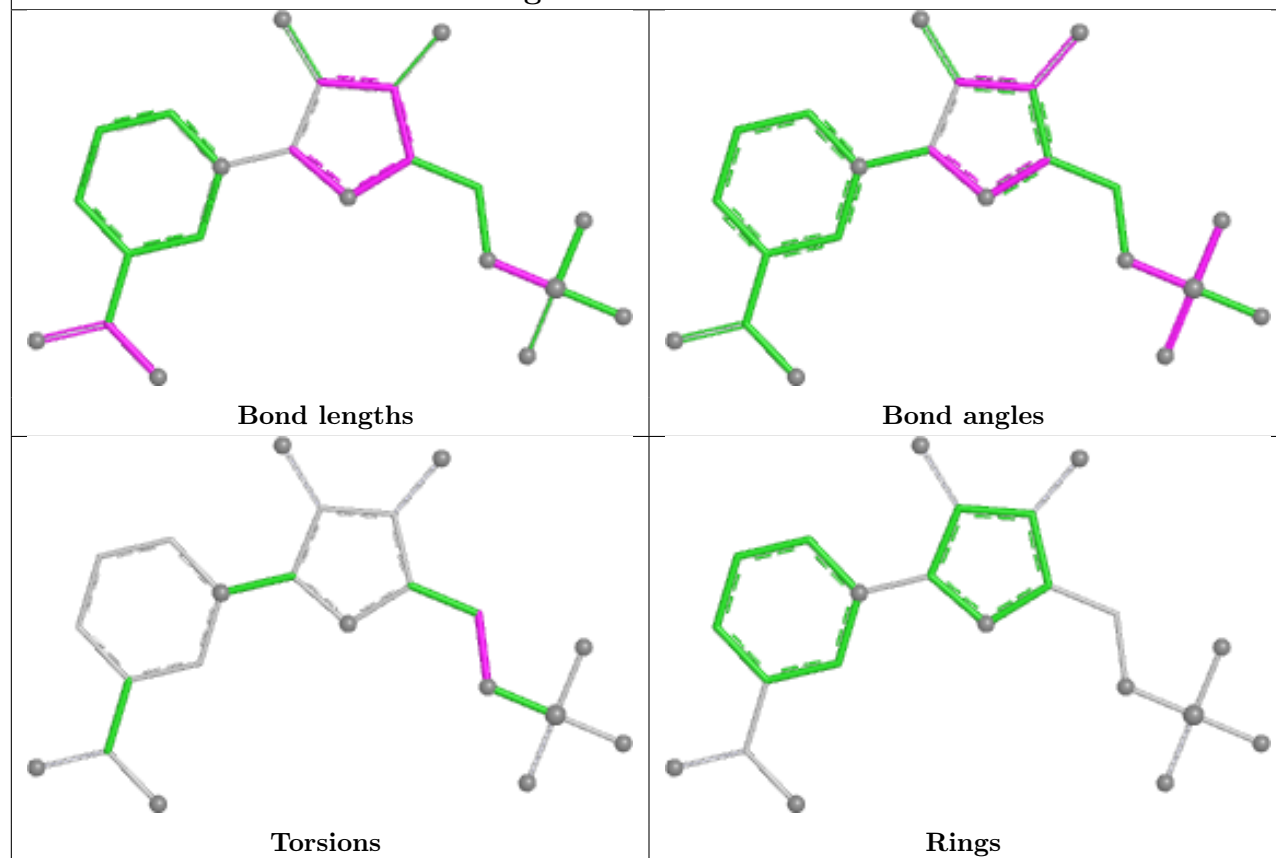
Ligand NMN D 502



Ligand NMN C 501



Ligand NMN A 501



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	450/452 (99%)	-0.69	0 100 100	3, 10, 24, 40	1 (0%)
1	B	450/452 (99%)	-0.57	1 (0%) 91 93	2, 12, 28, 37	1 (0%)
1	C	450/452 (99%)	-0.51	4 (0%) 81 86	5, 14, 29, 46	2 (0%)
1	D	446/452 (98%)	-0.35	4 (0%) 81 86	6, 16, 31, 55	0
All	All	1796/1808 (99%)	-0.53	9 (0%) 87 91	2, 13, 29, 55	4 (0%)

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	436	GLU	3.0
1	D	435	ASN	2.6
1	C	451	PRO	2.5
1	D	437	GLU	2.5
1	C	435	ASN	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

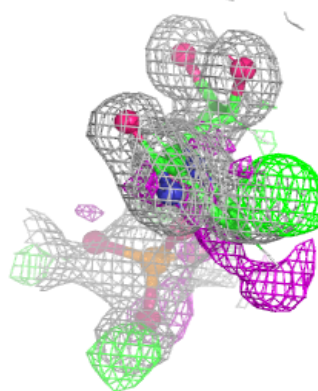
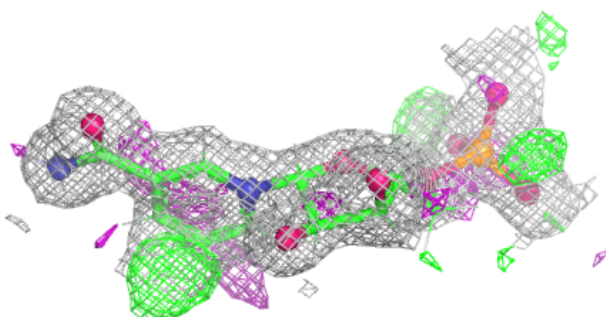
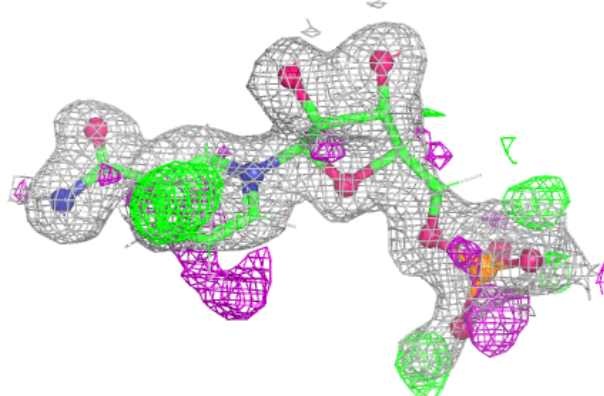
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FMT	C	508	3/3	0.59	0.27	35,40,41,48	0
4	FMT	B	506	3/3	0.61	0.20	35,36,42,50	0
4	FMT	C	506	3/3	0.75	0.17	32,38,41,42	0
5	PGE	A	506	10/10	0.77	0.15	29,38,51,57	0
4	FMT	A	505	3/3	0.78	0.19	29,35,35,39	0
4	FMT	D	506	3/3	0.80	0.14	44,44,49,53	0
4	FMT	D	501	3/3	0.81	0.19	24,31,34,37	0
2	NMN	B	502	22/22	0.82	0.13	6,23,42,45	0
4	FMT	A	507	3/3	0.83	0.12	31,35,41,49	0
4	FMT	C	510	3/3	0.83	0.15	38,39,45,47	0
4	FMT	C	511	3/3	0.83	0.13	22,23,37,45	0
4	FMT	D	507	3/3	0.84	0.13	34,36,41,43	0
4	FMT	C	504	3/3	0.84	0.16	21,21,26,42	0
4	FMT	D	508	3/3	0.85	0.20	43,43,45,55	0
4	FMT	A	512	3/3	0.86	0.14	20,22,37,45	0
4	FMT	B	507	3/3	0.86	0.12	38,39,42,50	0
4	FMT	D	509	3/3	0.87	0.12	26,32,36,39	0
4	FMT	C	505	3/3	0.87	0.10	33,33,36,40	0
6	EDO	A	510	4/4	0.87	0.18	16,20,30,30	0
2	NMN	A	501	22/22	0.88	0.11	9,17,38,48	0
2	NMN	D	502	22/22	0.89	0.10	10,23,45,58	0
3	GOL	B	501	6/6	0.90	0.10	13,18,26,26	0
3	GOL	B	505	6/6	0.90	0.12	11,20,29,29	0
3	GOL	D	503	6/6	0.90	0.10	14,19,26,26	0
4	FMT	A	504	3/3	0.90	0.14	17,21,33,35	0
4	FMT	C	507	3/3	0.90	0.12	22,22,27,41	0
4	FMT	A	508	3/3	0.91	0.19	27,28,29,35	0
3	GOL	B	503	6/6	0.92	0.09	14,17,21,23	0
3	GOL	D	504	6/6	0.92	0.09	15,22,32,32	0
4	FMT	C	509	3/3	0.92	0.12	25,26,32,32	0
2	NMN	C	501	22/22	0.92	0.09	8,18,40,48	0
3	GOL	D	505	6/6	0.94	0.06	12,15,20,24	0
4	FMT	A	509	3/3	0.94	0.25	14,17,20,24	0
3	GOL	C	503	6/6	0.94	0.10	11,19,24,29	0
3	GOL	A	511	6/6	0.94	0.09	19,25,43,43	0
3	GOL	A	502	6/6	0.94	0.08	10,15,17,18	0
7	MG	D	510	1/1	0.94	0.14	29,29,29,29	0
3	GOL	A	503	6/6	0.97	0.05	10,13,20,21	0
3	GOL	C	502	6/6	0.97	0.07	9,15,19,22	0
3	GOL	B	504	6/6	0.97	0.05	8,11,15,19	0
7	MG	C	512	1/1	0.98	0.04	24,24,24,24	1

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.

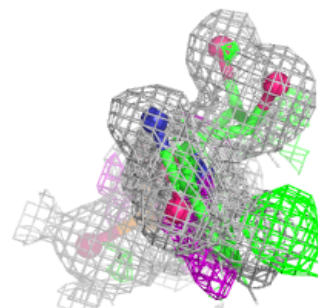
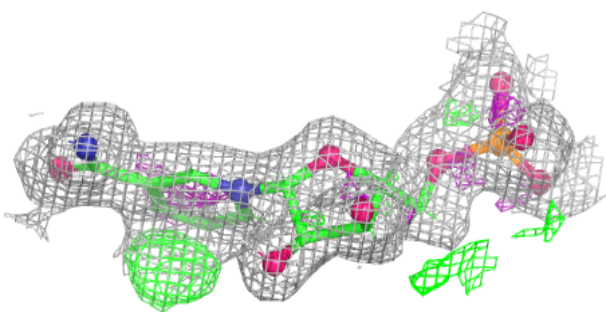
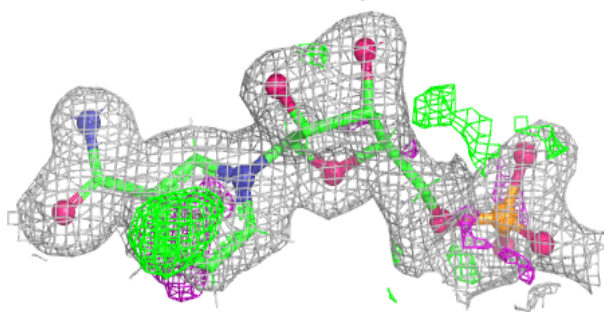
Electron density around NMN B 502:

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



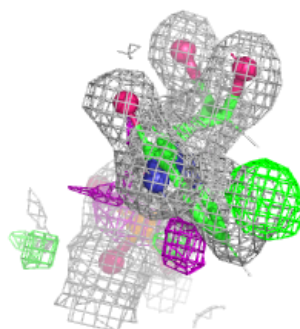
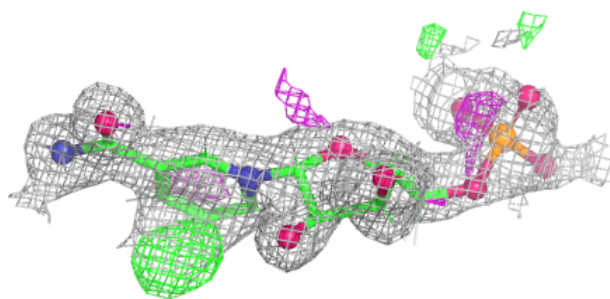
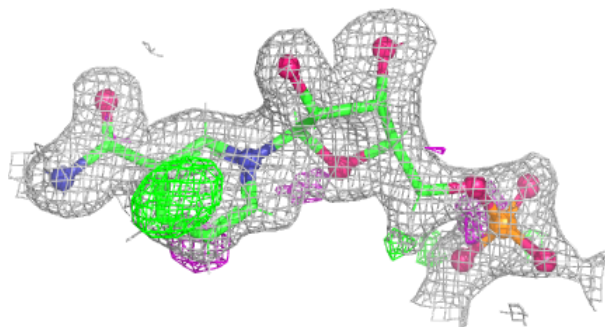
Electron density around NMN A 501:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

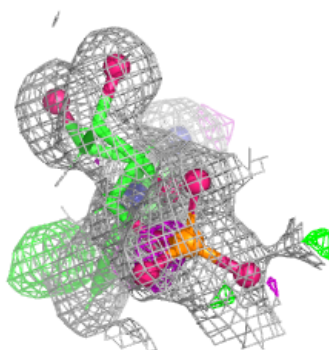
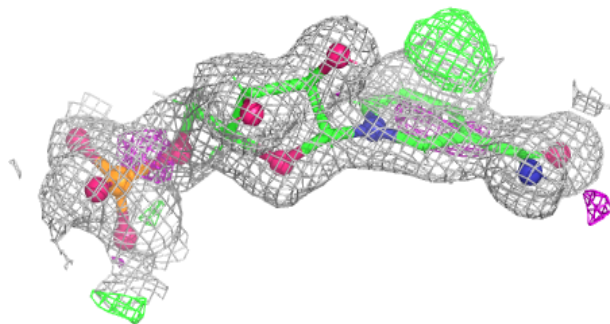
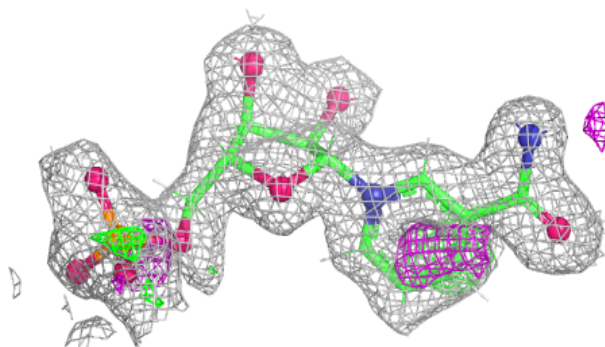


Electron density around NMN D 502:

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

**Electron density around NMN 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.