



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 10, 2024 – 02:56 PM EST

PDB ID : 1M2W
Title : Pseudomonas fluorescens mannitol 2-dehydrogenase ternary complex with NAD and D-mannitol
Authors : Kavanagh, K.L.; Klimacek, M.; Nidetzky, B.; Wilson, D.K.
Deposited on : 2002-06-25
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

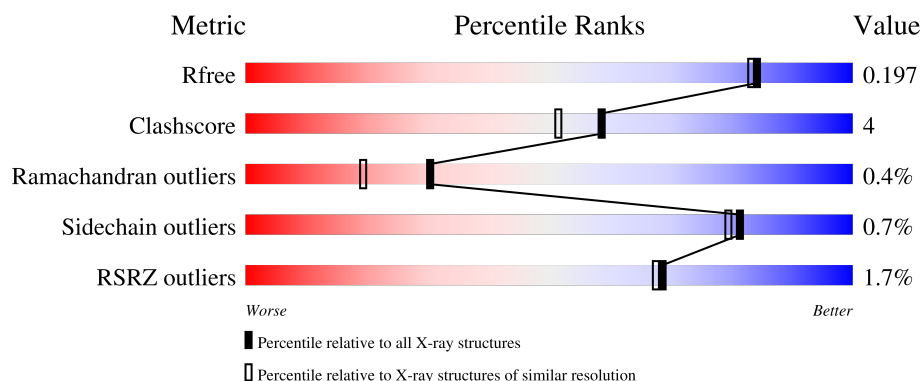
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

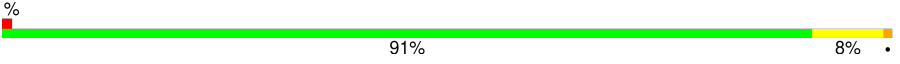

The reported resolution of this entry is 1.80 Å.

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}	164625	7108 (1.80-1.80)
Clashscore	180529	8162 (1.80-1.80)
Ramachandran outliers	177936	8077 (1.80-1.80)
Sidechain outliers	177891	8076 (1.80-1.80)
RSRZ outliers	164620	7108 (1.80-1.80)

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	493	 91% 8%
1	B	493	 90% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8400 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.

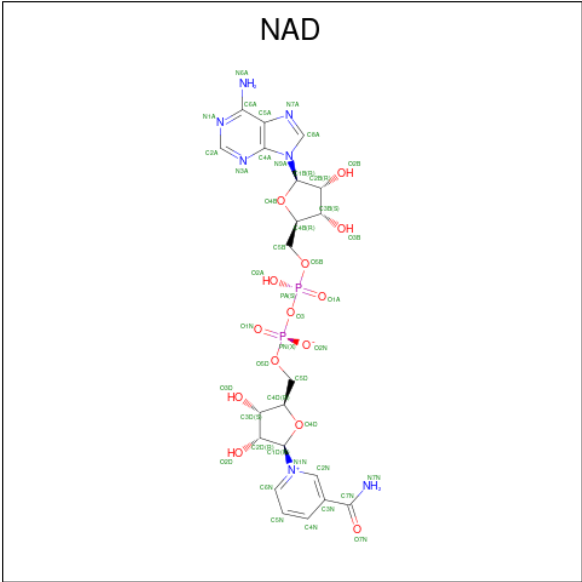
- Molecule 1 is a protein called mannitol dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	Se	0	0	0
			3829	2429	655	726	8	11			
1	B	492	Total	C	N	O	S	Se	0	0	0
			3829	2429	655	726	8	11			

There are 22 discrepancies between the modelled and reference sequences:

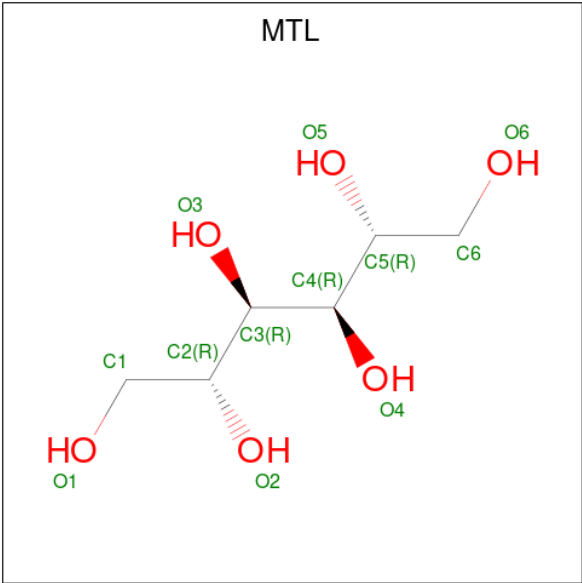
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	cloning artifact	UNP O08355
A	50	MSE	MET	cloning artifact	UNP O08355
A	105	MSE	MET	cloning artifact	UNP O08355
A	146	MSE	MET	cloning artifact	UNP O08355
A	187	MSE	MET	cloning artifact	UNP O08355
A	228	MSE	MET	cloning artifact	UNP O08355
A	235	MSE	MET	cloning artifact	UNP O08355
A	294	MSE	MET	cloning artifact	UNP O08355
A	322	MSE	MET	cloning artifact	UNP O08355
A	331	MSE	MET	cloning artifact	UNP O08355
A	335	MSE	MET	cloning artifact	UNP O08355
B	1	MSE	MET	cloning artifact	UNP O08355
B	50	MSE	MET	cloning artifact	UNP O08355
B	105	MSE	MET	cloning artifact	UNP O08355
B	146	MSE	MET	cloning artifact	UNP O08355
B	187	MSE	MET	cloning artifact	UNP O08355
B	228	MSE	MET	cloning artifact	UNP O08355
B	235	MSE	MET	cloning artifact	UNP O08355
B	294	MSE	MET	cloning artifact	UNP O08355
B	322	MSE	MET	cloning artifact	UNP O08355
B	331	MSE	MET	cloning artifact	UNP O08355
B	335	MSE	MET	cloning artifact	UNP O08355

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



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

- Molecule 3 is D-MANNITOL (three-letter code: MTL) (formula: C₆H₁₄O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		

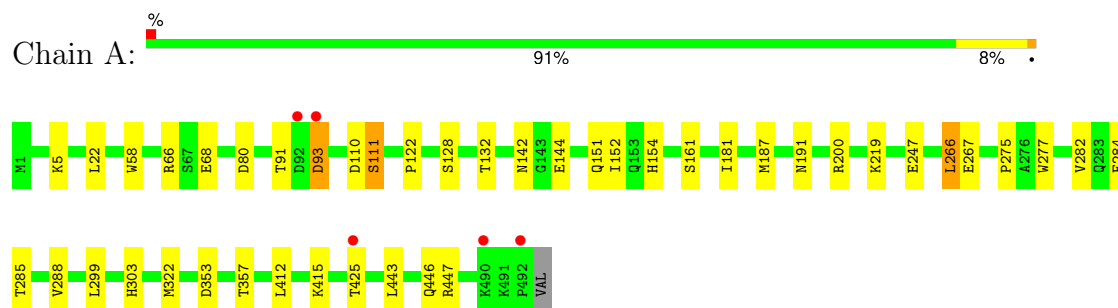
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	350	Total 350	O 350	0	0
4	B	280	Total 280	O 280	0	0

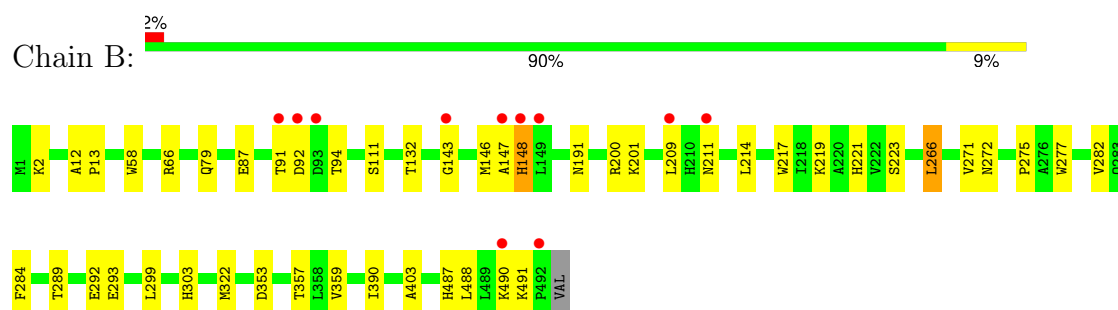
3 Residue-property plots

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: mannitol dehydrogenase



- Molecule 1: mannitol dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.95Å 104.53Å 101.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.80 30.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.00-1.80) 98.7 (30.00-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 1.80Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.176 , 0.202 0.171 , 0.197	Depositor DCC
R_{free} test set	5257 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for k,h,-l 0.008 for -l,-k,-h 0.011 for -h,l,k 0.000 for l,h,k 0.000 for k,l,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8400	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.05% 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: MTL, NAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.74	1/3907 (0.0%)	0.79	1/5301 (0.0%)
1	B	0.65	0/3907	0.77	1/5301 (0.0%)
All	All	0.70	1/7814 (0.0%)	0.78	2/10602 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	322	MSE	SE-CE	-8.76	1.43	1.95

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	LEU	CA-CB-CG	8.43	134.69	115.30
1	B	266	LEU	CA-CB-CG	6.80	130.95	115.30

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	3829	0	3757	29	0
1	B	3829	0	3757	33	0
2	A	44	0	26	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	44	0	26	5	0
3	A	12	0	14	2	0
3	B	12	0	14	2	0
4	A	350	0	0	3	0
4	B	280	0	0	0	0
All	All	8400	0	7594	66	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 (66) 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:6500:NAD:C4N	3:B:6600:MTL:H2	1.98	0.93
2:A:5500:NAD:C4N	3:A:5600:MTL:H2	2.01	0.90
2:A:5500:NAD:H4N	3:A:5600:MTL:H2	1.62	0.81
2:B:6500:NAD:H4N	3:B:6600:MTL:H2	1.62	0.81
1:A:200:ARG:NH2	1:A:219:LYS:HE3	2.01	0.75
1:A:267:GLU:HG2	1:A:288:VAL:HG13	1.80	0.64
1:B:87:GLU:OE1	1:B:94:THR:HA	2.01	0.60
1:A:412:LEU:HD12	1:A:415:LYS:HD3	1.81	0.60
1:B:66:ARG:HD3	2:B:6500:NAD:O2B	2.02	0.58
1:B:200:ARG:HH21	1:B:219:LYS:HZ2	1.52	0.57
1:B:200:ARG:NH2	1:B:219:LYS:NZ	2.53	0.57
1:A:22:LEU:HD22	4:A:5944:HOH:O	2.03	0.57
1:A:142:ASN:OD1	1:A:144:GLU:HG2	2.05	0.57
1:B:211:ASN:HD22	1:B:214:LEU:HB2	1.69	0.57
1:B:146:MSE:HB3	1:B:148:HIS:CE1	2.40	0.56
1:B:211:ASN:ND2	1:B:214:LEU:H	2.03	0.56
1:A:58:TRP:CG	1:A:275:PRO:HG3	2.40	0.56
1:B:58:TRP:CG	1:B:275:PRO:HG3	2.40	0.56
1:B:487:HIS:O	1:B:490:LYS:HG2	2.06	0.55
1:B:322:MSE:CE	1:B:359:VAL:HG13	2.38	0.54
1:A:200:ARG:HH22	1:A:219:LYS:HE3	1.72	0.52
1:B:292:GLU:HG3	1:B:293:GLU:N	2.24	0.52
1:A:277:TRP:HB3	1:A:282:VAL:HG21	1.93	0.51
1:A:5:LYS:HE2	1:A:22:LEU:HD13	1.92	0.50
1:B:322:MSE:HE1	1:B:359:VAL:HG13	1.93	0.50
1:B:132:THR:HA	2:B:6500:NAD:O3D	2.11	0.49
1:A:267:GLU:CG	1:A:288:VAL:HG13	2.42	0.48
1:A:299:LEU:C	1:A:299:LEU:HD23	2.34	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:488:LEU:HD12	1:B:491:LYS:HD2	1.95	0.48
1:B:277:TRP:HB3	1:B:282:VAL:HG21	1.96	0.48
1:B:299:LEU:C	1:B:299:LEU:HD23	2.35	0.47
1:B:2:LYS:HE3	1:B:79:GLN:HG2	1.96	0.47
1:B:200:ARG:NH2	1:B:219:LYS:HZ2	2.13	0.47
1:B:143:GLY:O	1:B:201:LYS:HE3	2.14	0.47
1:B:289:THR:HA	1:B:292:GLU:HG2	1.97	0.47
1:A:154:HIS:HE1	1:A:161:SER:O	1.99	0.46
1:B:223:SER:HB3	1:B:271:VAL:HG13	1.97	0.46
1:A:5:LYS:HE3	1:A:80:ASP:HB3	1.98	0.45
1:A:5:LYS:HG3	1:A:80:ASP:CB	2.47	0.45
1:B:266:LEU:O	1:B:284:PHE:HA	2.17	0.45
1:B:390:ILE:HG12	1:B:403:ALA:HB1	1.98	0.44
1:A:191:ASN:HD21	2:A:5500:NAD:C6N	2.30	0.44
1:B:200:ARG:NH2	1:B:219:LYS:HZ3	2.16	0.44
1:A:122:PRO:HG3	1:A:181:ILE:HD13	1.99	0.44
1:B:277:TRP:HB3	1:B:282:VAL:CG2	2.48	0.43
1:B:191:ASN:HD21	2:B:6500:NAD:C6N	2.32	0.43
1:A:266:LEU:O	1:A:284:PHE:HA	2.19	0.42
1:A:446:GLN:HG3	4:A:5881:HOH:O	2.19	0.42
1:A:353:ASP:O	1:A:357:THR:HG23	2.20	0.42
1:A:128:SER:HB3	1:A:187:MSE:HE2	2.00	0.41
1:B:211:ASN:HD22	1:B:214:LEU:CB	2.34	0.41
1:B:353:ASP:O	1:B:357:THR:HG23	2.20	0.41
1:B:92:ASP:C	1:B:94:THR:H	2.23	0.41
1:A:68:GLU:H	1:A:68:GLU:CD	2.24	0.41
1:A:110:ASP:O	1:A:111:SER:HB2	2.21	0.41
1:A:132:THR:HA	2:A:5500:NAD:O3D	2.21	0.41
1:B:12:ALA:HA	1:B:13:PRO:HD3	1.86	0.41
1:B:147:ALA:HB2	1:B:209:LEU:HD11	2.01	0.41
1:B:200:ARG:HH22	1:B:219:LYS:HZ3	1.69	0.41
1:B:217:TRP:CD1	1:B:221:HIS:CD2	3.08	0.41
1:A:91:THR:OG1	1:A:93:ASP:OD1	2.39	0.40
1:A:151:GLN:HG2	1:A:152:ILE:N	2.36	0.40
1:A:285:THR:CG2	1:A:288:VAL:HG12	2.51	0.40
1:A:247:GLU:OE1	4:A:5903:HOH:O	2.22	0.40
1:A:277:TRP:HB3	1:A:282:VAL:CG2	2.50	0.40
1:A:443:LEU:O	1:A:447:ARG:HG3	2.21	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	490/493 (99%)	479 (98%)	9 (2%)	2 (0%)	30	19
1	B	490/493 (99%)	472 (96%)	16 (3%)	2 (0%)	30	19
All	All	980/986 (99%)	951 (97%)	25 (3%)	4 (0%)	30	19

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	111	SER
1	B	111	SER
1	B	148	HIS
1	A	93	ASP

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	410/400 (102%)	407 (99%)	3 (1%)	81	79
1	B	410/400 (102%)	407 (99%)	3 (1%)	81	79
All	All	820/800 (102%)	814 (99%)	6 (1%)	81	79

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	ARG
1	A	303	HIS

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Mol	Chain	Res	Type
1	A	425	THR
1	B	91	THR
1	B	272	ASN
1	B	303	HIS

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	191	ASN
1	A	446	GLN
1	B	148	HIS
1	B	191	ASN
1	B	211	ASN
1	B	221	HIS
1	B	243	GLN
1	B	272	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	B	6500	-	42,48,48	2.41	12 (28%)	50,73,73	1.76	8 (16%)
3	MTL	A	5600	-	11,11,11	0.99	1 (9%)	14,14,14	1.03	0
2	NAD	A	5500	-	42,48,48	2.59	12 (28%)	50,73,73	1.78	10 (20%)
3	MTL	B	6600	-	11,11,11	1.12	1 (9%)	14,14,14	1.03	1 (7%)

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	B	6500	-	-	1/26/62/62	0/5/5/5
3	MTL	A	5600	-	-	0/16/16/16	-
2	NAD	A	5500	-	-	1/26/62/62	0/5/5/5
3	MTL	B	6600	-	-	0/16/16/16	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	5500	NAD	C2N-N1N	8.76	1.44	1.35
2	B	6500	NAD	C2N-N1N	6.81	1.42	1.35
2	B	6500	NAD	O4D-C1D	6.52	1.49	1.40
2	A	5500	NAD	O4D-C1D	6.30	1.49	1.40
2	A	5500	NAD	C2A-N3A	5.21	1.40	1.32
2	A	5500	NAD	O7N-C7N	5.17	1.33	1.24
2	B	6500	NAD	C2A-N3A	5.12	1.40	1.32
2	A	5500	NAD	C4A-N3A	4.64	1.41	1.35
2	B	6500	NAD	C2A-N1A	4.57	1.42	1.33
2	A	5500	NAD	C2A-N1A	4.18	1.41	1.33
2	B	6500	NAD	C4N-C3N	4.12	1.45	1.39
2	B	6500	NAD	O7N-C7N	3.61	1.30	1.24
2	A	5500	NAD	C4N-C3N	3.55	1.44	1.39
2	B	6500	NAD	C6N-N1N	3.37	1.43	1.35
2	B	6500	NAD	O4B-C4B	3.27	1.52	1.45
2	B	6500	NAD	C4A-N3A	3.06	1.39	1.35
2	A	5500	NAD	C6N-N1N	2.96	1.42	1.35
2	B	6500	NAD	O4B-C1B	2.63	1.44	1.40
2	A	5500	NAD	O4B-C1B	2.57	1.44	1.40
2	B	6500	NAD	C2N-C3N	2.47	1.42	1.39
2	A	5500	NAD	O4B-C4B	2.41	1.50	1.45
3	A	5600	MTL	C6-C5	2.37	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	5500	NAD	C1B-N9A	-2.18	1.44	1.49
2	B	6500	NAD	C5B-C4B	2.07	1.57	1.51
3	B	6600	MTL	C6-C5	2.07	1.57	1.52
2	A	5500	NAD	C2N-C3N	2.06	1.42	1.39

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	5500	NAD	C6N-N1N-C1D	-5.93	108.10	119.73
2	B	6500	NAD	C6N-N1N-C1D	-5.89	108.16	119.73
2	B	6500	NAD	N3A-C2A-N1A	-4.55	122.50	128.67
2	A	5500	NAD	N3A-C2A-N1A	-4.21	122.95	128.67
2	A	5500	NAD	C2N-N1N-C1D	4.11	128.21	119.13
2	B	6500	NAD	C2N-N1N-C1D	4.06	128.10	119.13
2	B	6500	NAD	C5A-C6A-N6A	4.06	126.50	120.31
2	A	5500	NAD	C5A-C6A-N6A	3.64	125.86	120.31
2	A	5500	NAD	C1B-N9A-C4A	-3.26	120.91	126.64
2	B	6500	NAD	O2A-PA-O1A	2.65	124.77	112.44
2	A	5500	NAD	O2A-PA-O1A	2.63	124.69	112.44
2	A	5500	NAD	O3B-C3B-C4B	-2.59	103.64	111.08
2	A	5500	NAD	O5B-C5B-C4B	-2.43	100.71	108.99
2	B	6500	NAD	C2N-C3N-C4N	2.31	120.95	118.26
2	B	6500	NAD	C6N-N1N-C2N	2.16	123.72	121.88
2	B	6500	NAD	O5B-C5B-C4B	-2.11	101.80	108.99
2	A	5500	NAD	C6N-N1N-C2N	2.11	123.67	121.88
3	B	6600	MTL	O1-C1-C2	-2.07	106.81	111.16
2	A	5500	NAD	O4D-C4D-C5D	2.01	115.78	109.33

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	6500	NAD	O4B-C4B-C5B-O5B
2	A	5500	NAD	O4B-C4B-C5B-O5B

There are no ring outliers.

4 monomers are involved in 9 short contacts:

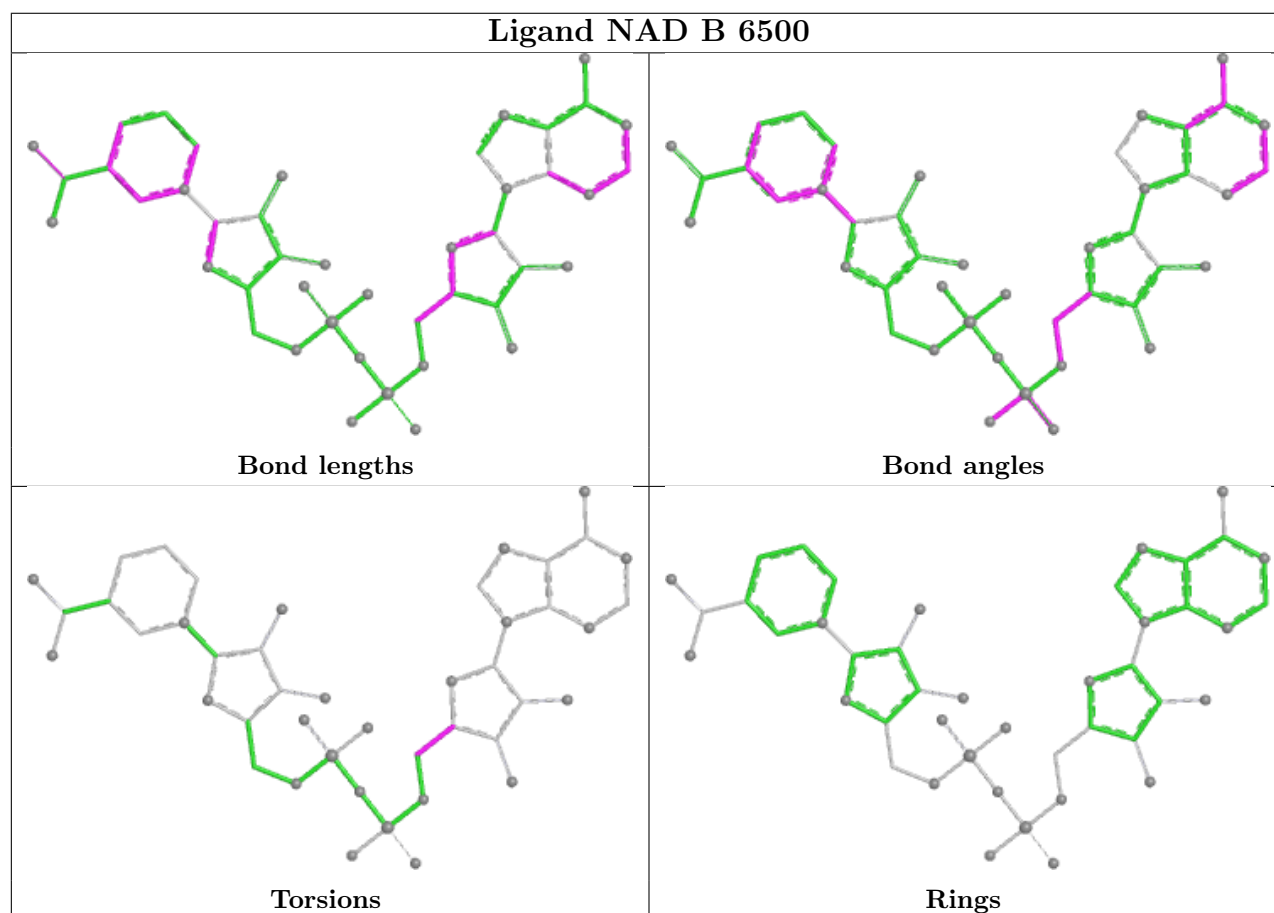
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	6500	NAD	5	0
3	A	5600	MTL	2	0

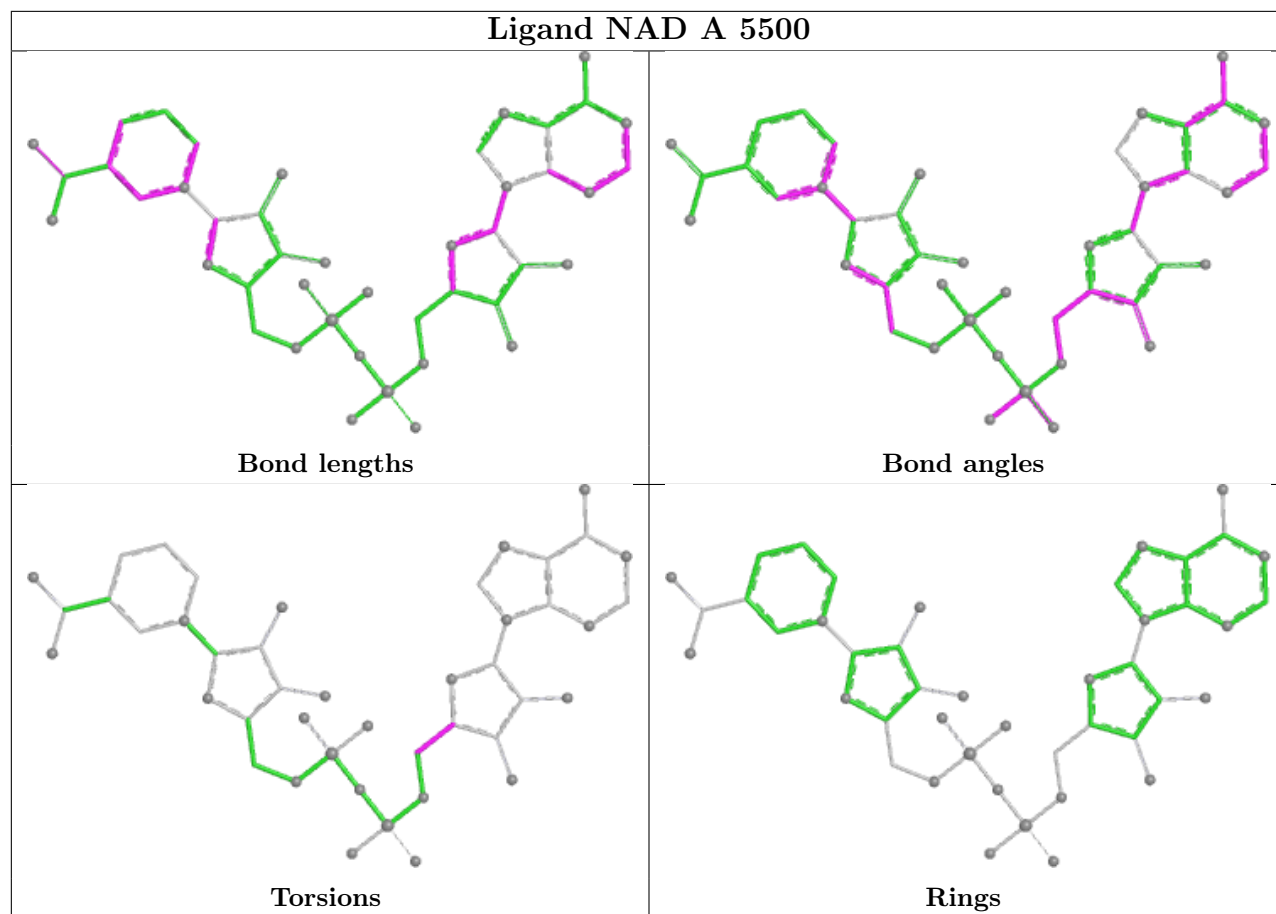
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	5500	NAD	4	0
3	B	6600	MTL	2	0

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 [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	481/493 (97%)	-0.47	5 (1%) 79 78	10, 18, 34, 56	0
1	B	481/493 (97%)	-0.25	11 (2%) 61 59	12, 21, 38, 61	0
All	All	962/986 (97%)	-0.36	16 (1%) 69 67	10, 19, 37, 61	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	492	PRO	4.9
1	B	93	ASP	3.9
1	B	148	HIS	3.5
1	A	492	PRO	3.4
1	B	147	ALA	3.2
1	A	93	ASP	2.9
1	A	92	ASP	2.9
1	A	490	LYS	2.8
1	B	211	ASN	2.6
1	B	143	GLY	2.6
1	B	91	THR	2.5
1	B	149	LEU	2.4
1	A	425	THR	2.4
1	B	92	ASP	2.1
1	B	209	LEU	2.0
1	B	490	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

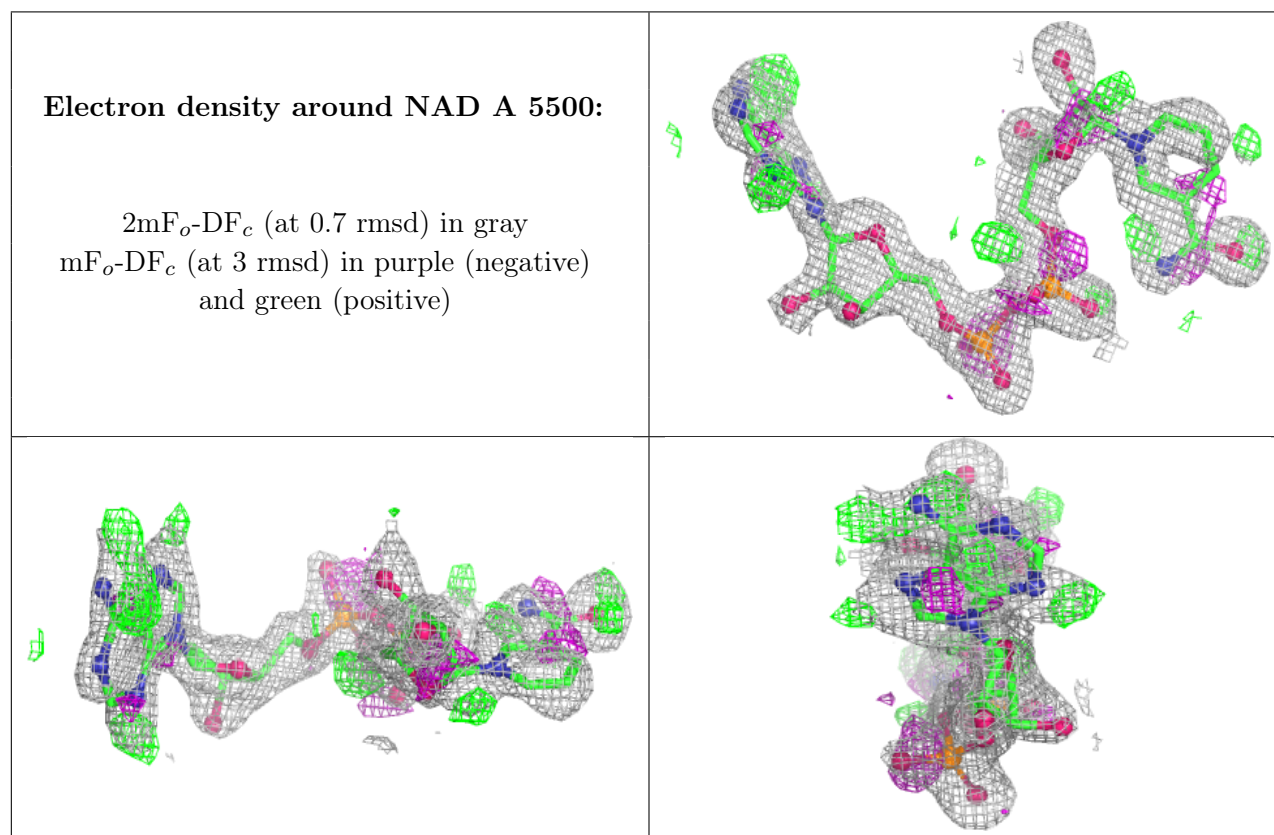
There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

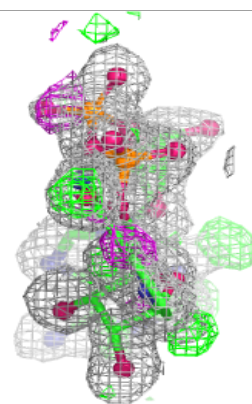
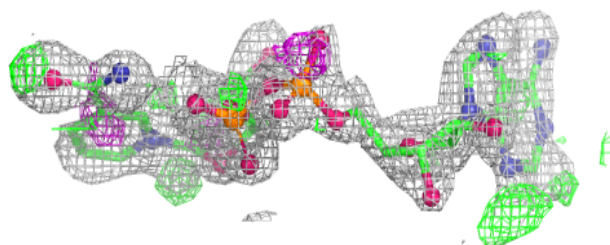
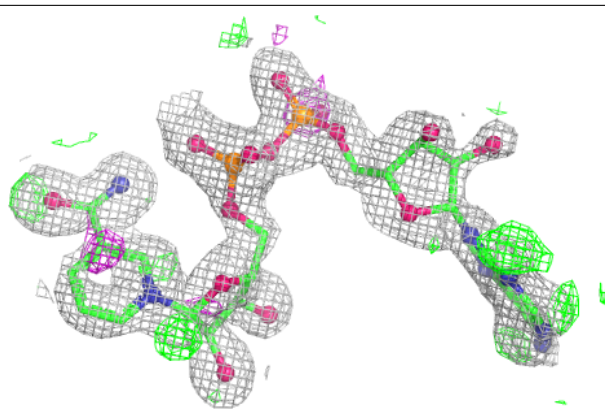
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAD	A	5500	44/44	0.84	0.14	7,30,35,36	44
2	NAD	B	6500	44/44	0.84	0.14	13,29,35,37	44
3	MTL	B	6600	12/12	0.97	0.04	12,14,16,16	0
3	MTL	A	5600	12/12	0.98	0.04	10,11,13,14	0

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 6500:

$2mF_o-DF_c$ (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.