



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:33 AM UTC

PDB ID : 6HLI / pdb_00006hli
Title : wild-type NuoEF from Aquifex aeolicus - reduced form bound to NAD+
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Deposited on : 2018-09-11
Resolution : 2.38 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	:	4-5-2 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

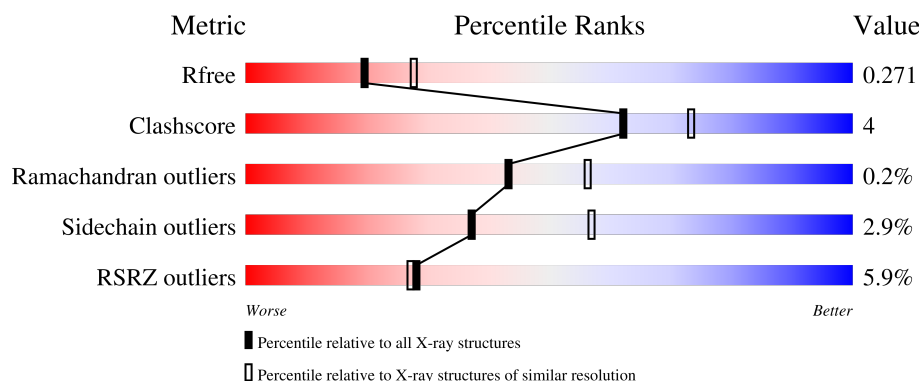
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	160	
1	C	160	
2	B	434	
2	D	434	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9517 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 NADH-quinone oxidoreductase subunit E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	155	Total	C	N	O	S	0	0	0
			1259	816	203	231	9			
1	C	155	Total	C	N	O	S	0	0	0
			1259	816	203	231	9			

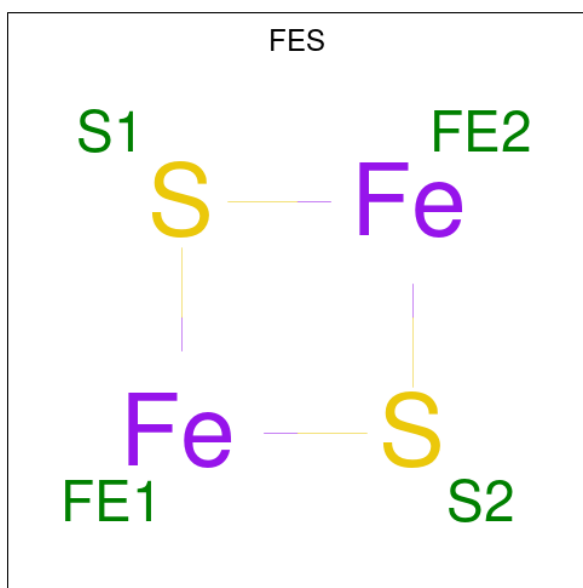
- Molecule 2 is a protein called NADH-quinone oxidoreductase subunit F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	417	Total	C	N	O	S	0	0	0
			3284	2111	547	613	13			
2	D	418	Total	C	N	O	S	0	0	0
			3293	2117	549	614	13			

There are 16 discrepancies between the modelled and reference sequences:

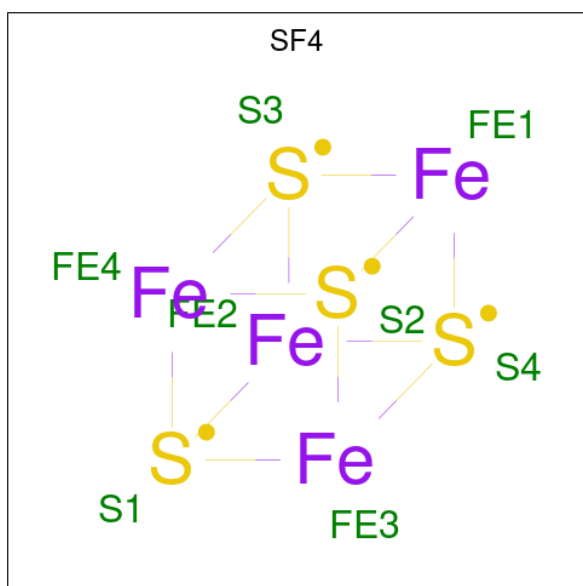
Chain	Residue	Modelled	Actual	Comment	Reference
B	427	ALA	-	expression tag	UNP O66841
B	428	GLY	-	expression tag	UNP O66841
B	429	HIS	-	expression tag	UNP O66841
B	430	HIS	-	expression tag	UNP O66841
B	431	HIS	-	expression tag	UNP O66841
B	432	HIS	-	expression tag	UNP O66841
B	433	HIS	-	expression tag	UNP O66841
B	434	HIS	-	expression tag	UNP O66841
D	427	ALA	-	expression tag	UNP O66841
D	428	GLY	-	expression tag	UNP O66841
D	429	HIS	-	expression tag	UNP O66841
D	430	HIS	-	expression tag	UNP O66841
D	431	HIS	-	expression tag	UNP O66841
D	432	HIS	-	expression tag	UNP O66841
D	433	HIS	-	expression tag	UNP O66841
D	434	HIS	-	expression tag	UNP O66841

- Molecule 3 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



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

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



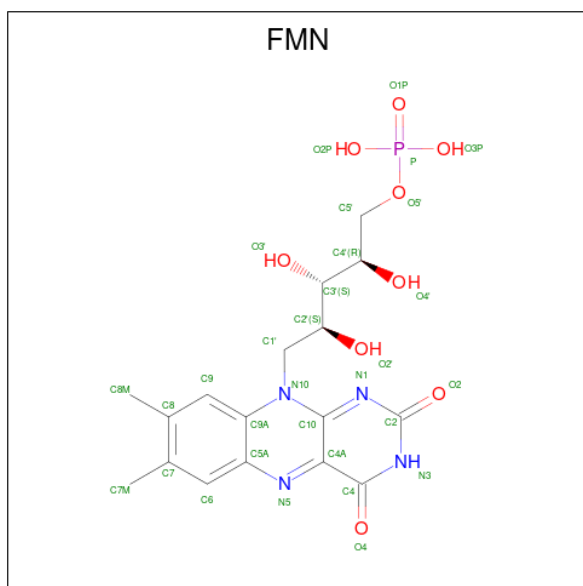
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	Fe	S	0	0
			8	4	4		

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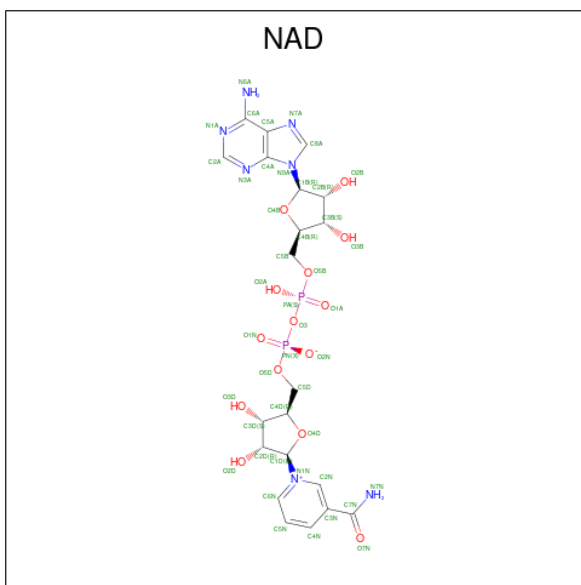
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 5 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
5	D	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 6 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

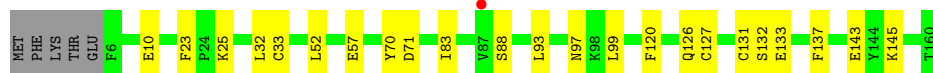
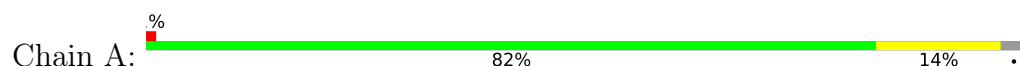
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	27	Total O 27 27	0	0
7	B	107	Total O 107 107	0	0
7	C	23	Total O 23 23	0	0
7	D	91	Total O 91 91	0	0

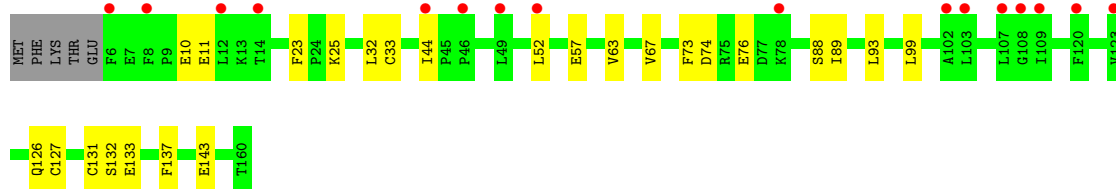
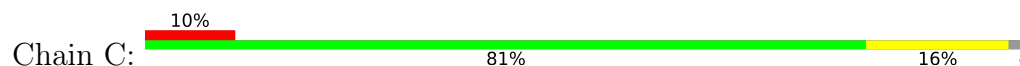
3 Residue-property plots [i](#)

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.

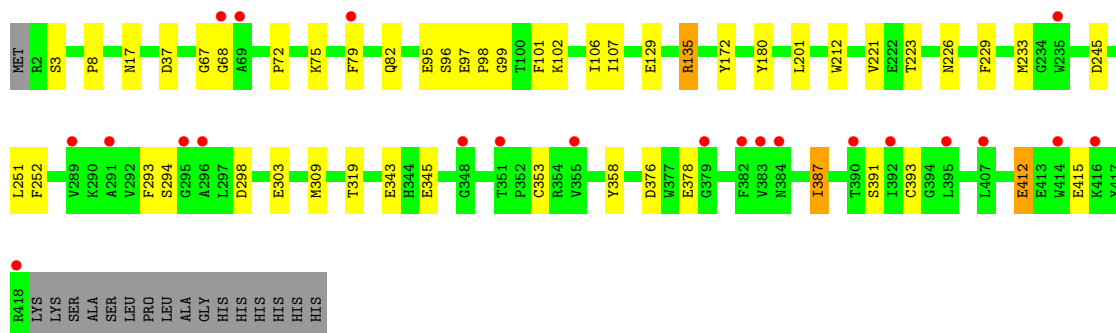
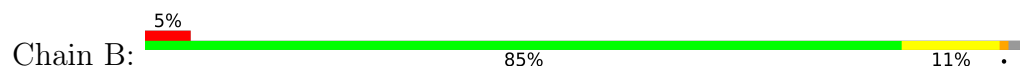
- Molecule 1: NADH-quinone oxidoreductase subunit E



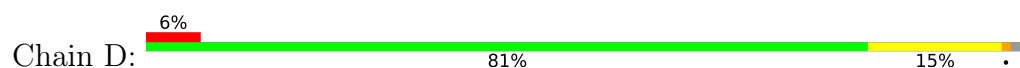
- Molecule 1: NADH-quinone oxidoreductase subunit E

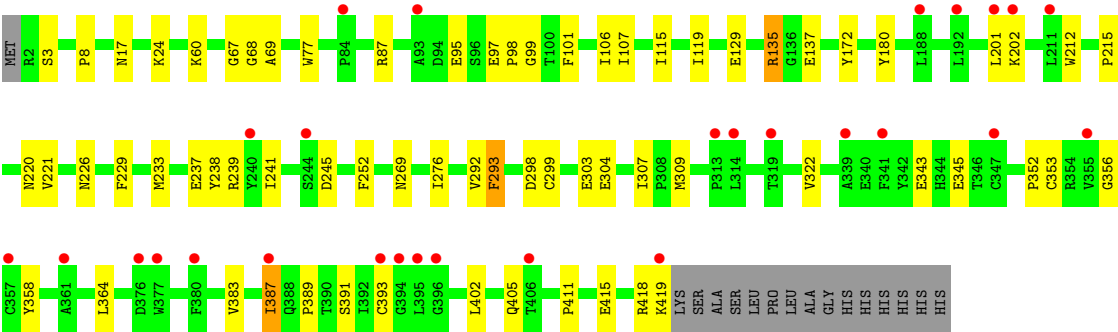


- Molecule 2: NADH-quinone oxidoreductase subunit F



- Molecule 2: NADH-quinone oxidoreductase subunit F





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.59Å 116.04Å 189.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.30 – 2.38 60.30 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (60.30-2.38) 99.8 (60.30-2.38)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.37Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.205 , 0.238 0.230 , 0.271	Depositor DCC
R_{free} test set	2841 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 76.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9517	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.03 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.3718e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: FMN, NAD, SF4, FES

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	0.79	0/1288	1.31	4/1740 (0.2%)
1	C	0.78	0/1288	1.31	4/1740 (0.2%)
2	B	0.88	0/3365	1.28	11/4553 (0.2%)
2	D	0.88	0/3374	1.31	14/4564 (0.3%)
All	All	0.85	0/9315	1.30	33/12597 (0.3%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	387	ILE	N-CA-C	-6.53	106.19	111.81
2	D	226	ASN	CA-C-N	6.52	125.55	120.33
2	D	226	ASN	C-N-CA	6.52	125.55	120.33
2	B	226	ASN	CA-C-N	6.52	125.55	120.33
2	B	226	ASN	C-N-CA	6.52	125.55	120.33
2	B	387	ILE	N-CA-C	-6.05	106.60	111.81
2	B	376	ASP	CA-CB-CG	6.04	118.64	112.60
2	D	293	PHE	CA-CB-CG	6.01	119.81	113.80
2	B	387	ILE	CA-C-N	5.92	127.50	120.09
2	B	387	ILE	C-N-CA	5.92	127.50	120.09
2	D	387	ILE	CA-C-N	5.90	127.46	120.09
2	D	387	ILE	C-N-CA	5.90	127.46	120.09
1	A	70	TYR	CA-C-N	5.60	129.56	120.60
1	A	70	TYR	C-N-CA	5.60	129.56	120.60
2	D	245	ASP	CA-CB-CG	5.50	118.11	112.60
2	B	293	PHE	CA-CB-CG	5.44	119.24	113.80
2	B	245	ASP	CA-CB-CG	5.37	117.97	112.60
1	C	89	ILE	N-CA-C	5.34	115.79	110.23
1	C	88	SER	CA-C-N	5.29	127.70	120.77
1	C	88	SER	C-N-CA	5.29	127.70	120.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	74	ASP	CA-CB-CG	5.25	117.85	112.60
1	A	88	SER	CA-C-N	5.22	127.61	120.77
1	A	88	SER	C-N-CA	5.22	127.61	120.77
2	B	37	ASP	CA-CB-CG	5.17	117.77	112.60
2	B	223	THR	CA-C-N	5.13	127.49	120.46
2	B	223	THR	C-N-CA	5.13	127.49	120.46
2	D	77	TRP	CA-C-N	5.13	127.15	120.28
2	D	77	TRP	C-N-CA	5.13	127.15	120.28
2	D	24	LYS	N-CA-C	-5.10	106.56	112.89
2	D	220	ASN	CA-C-N	5.09	127.16	120.60
2	D	220	ASN	C-N-CA	5.09	127.16	120.60
2	D	411	PRO	CA-C-N	5.01	127.69	120.38
2	D	411	PRO	C-N-CA	5.01	127.69	120.38

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	1259	0	1263	13	0
1	C	1259	0	1263	12	0
2	B	3284	0	3257	28	0
2	D	3293	0	3270	37	0
3	A	4	0	0	0	0
3	C	4	0	0	0	0
4	B	8	0	0	1	0
4	D	8	0	0	1	0
5	B	31	0	19	0	0
5	D	31	0	19	0	0
6	B	44	0	26	6	0
6	D	44	0	26	7	0
7	A	27	0	0	1	0
7	B	107	0	0	0	0
7	C	23	0	0	0	0
7	D	91	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9517	0	9143	76	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 (76) 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:97:GLU:HB2	6:B:502:NAD:C4N	2.22	0.70
2:D:202:LYS:HD2	7:D:625:HOH:O	1.94	0.67
2:B:68:GLY:HA3	6:B:502:NAD:H1D	1.78	0.66
2:B:294:SER:HB3	2:B:298:ASP:HB2	1.79	0.64
2:D:293:PHE:HE1	2:D:299:CYS:HG	1.44	0.64
2:D:135:ARG:HD3	2:D:137:GLU:HB2	1.84	0.60
2:B:229:PHE:O	2:B:233:MET:HG2	2.03	0.58
2:D:229:PHE:O	2:D:233:MET:HG2	2.04	0.58
2:D:68:GLY:HA3	6:D:502:NAD:H1D	1.86	0.58
2:D:180:TYR:CE1	6:D:502:NAD:H5N	2.38	0.58
2:D:97:GLU:HB2	6:D:502:NAD:C4N	2.34	0.57
2:D:298:ASP:HA	2:D:405:GLN:HE22	1.70	0.56
2:D:180:TYR:OH	6:D:502:NAD:H5N	2.05	0.56
2:B:79:PHE:O	2:B:82:GLN:HG2	2.06	0.55
2:D:67:GLY:O	6:D:502:NAD:H2N	2.07	0.54
2:D:391:SER:HB2	4:D:500:SF4:S1	2.48	0.53
2:B:412:GLU:H	2:B:412:GLU:CD	2.17	0.52
1:C:63:VAL:O	1:C:67:VAL:HG23	2.10	0.52
2:B:97:GLU:HB2	6:B:502:NAD:H4N	1.93	0.51
2:D:299:CYS:HB2	2:D:402:LEU:HG	1.93	0.51
2:D:180:TYR:CZ	6:D:502:NAD:H5N	2.46	0.51
2:D:418:ARG:HH21	2:D:419:LYS:HE2	1.76	0.51
2:B:391:SER:HB2	4:B:500:SF4:S1	2.51	0.50
2:D:238:TYR:O	2:D:241:ILE:HD12	2.11	0.50
2:B:180:TYR:CE1	6:B:502:NAD:H5N	2.47	0.49
2:B:107:ILE:HD11	2:B:221:VAL:HG21	1.92	0.49
1:A:33:CYS:HB3	1:A:52:LEU:HD11	1.95	0.49
2:D:87:ARG:HG2	2:D:215:PRO:HB2	1.95	0.49
1:A:143:GLU:HB3	2:B:8:PRO:HD3	1.94	0.49
1:C:126:GLN:HG3	2:D:345:GLU:OE2	2.12	0.49
1:C:33:CYS:HB3	1:C:52:LEU:HD11	1.96	0.48
2:D:276:ILE:HD11	2:D:292:VAL:HG21	1.97	0.47
2:B:96:SER:HA	2:B:135:ARG:NH1	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:107:ILE:HD11	2:D:221:VAL:HG21	1.96	0.47
2:D:115:ILE:O	2:D:119:ILE:HG13	2.15	0.47
2:B:96:SER:HB2	2:B:180:TYR:HD1	1.80	0.46
1:A:97:ASN:HB2	7:A:320:HOH:O	2.14	0.46
2:D:292:VAL:HG11	2:D:322:VAL:HG13	1.96	0.46
2:B:129:GLU:HG3	2:B:172:TYR:HE2	1.81	0.46
1:C:44:ILE:HG13	1:C:73:PHE:HB3	1.97	0.46
2:B:252:PHE:CE1	2:B:309:MET:HB3	2.51	0.46
1:C:132:SER:HB3	2:D:101:PHE:CZ	2.51	0.45
1:A:83:ILE:HD11	1:A:120:PHE:CD1	2.51	0.45
2:B:67:GLY:O	6:B:502:NAD:H1D	2.17	0.45
1:C:133:GLU:HG3	2:D:8:PRO:HB2	1.99	0.45
1:C:127:CYS:SG	2:D:98:PRO:HA	2.57	0.44
2:D:252:PHE:CE1	2:D:309:MET:HB3	2.52	0.44
1:C:131:CYS:SG	2:D:99:GLY:HA2	2.57	0.44
1:A:126:GLN:HG3	2:B:345:GLU:OE2	2.17	0.44
1:A:132:SER:HB3	2:B:101:PHE:CZ	2.53	0.43
1:A:127:CYS:SG	2:B:98:PRO:HA	2.58	0.43
1:C:99:LEU:HD21	1:C:137:PHE:HB3	2.00	0.43
1:C:23:PHE:HE2	1:C:32:LEU:HD12	1.84	0.43
2:B:353:CYS:HA	2:B:387:ILE:HG23	2.01	0.43
2:D:129:GLU:HG3	2:D:172:TYR:HE2	1.83	0.43
1:A:133:GLU:HG3	2:B:8:PRO:HB2	1.99	0.43
2:D:353:CYS:HA	2:D:387:ILE:HG23	2.01	0.43
2:D:415:GLU:HG2	2:D:418:ARG:NH2	2.34	0.43
2:B:106:ILE:HD11	2:B:251:LEU:HD11	2.01	0.43
2:B:180:TYR:OH	6:B:502:NAD:H5N	2.18	0.42
1:A:131:CYS:SG	2:B:99:GLY:HA2	2.60	0.42
1:C:25:LYS:HD2	2:D:212:TRP:CZ2	2.55	0.42
2:D:69:ALA:HB3	6:D:502:NAD:H51A	2.02	0.42
2:B:343:GLU:HG3	2:B:358:TYR:HA	2.01	0.42
1:A:23:PHE:HE2	1:A:32:LEU:HD12	1.85	0.41
1:C:143:GLU:HB3	2:D:8:PRO:HD3	2.02	0.41
2:D:356:GLY:HA2	2:D:383:VAL:HG13	2.02	0.41
2:B:102:LYS:HE2	2:B:319:THR:HB	2.02	0.41
2:B:72:PRO:HB2	2:B:75:LYS:HB2	2.02	0.41
1:A:99:LEU:HD21	1:A:137:PHE:HB3	2.02	0.41
2:D:60:LYS:O	2:D:239:ARG:HD2	2.20	0.41
2:D:343:GLU:HG3	2:D:358:TYR:HA	2.01	0.41
1:A:25:LYS:HD2	2:B:212:TRP:CZ2	2.57	0.41
1:A:133:GLU:HB3	1:A:145:LYS:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:304:GLU:O	2:D:307:ILE:HG12	2.21	0.40
2:D:352:PRO:HG2	2:D:389:PRO:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	153/160 (96%)	145 (95%)	8 (5%)	0	100	100
1	C	153/160 (96%)	146 (95%)	7 (5%)	0	100	100
2	B	415/434 (96%)	403 (97%)	11 (3%)	1 (0%)	43	56
2	D	416/434 (96%)	402 (97%)	13 (3%)	1 (0%)	43	56
All	All	1137/1188 (96%)	1096 (96%)	39 (3%)	2 (0%)	43	56

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	95	GLU
2	D	95	GLU

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	141/146 (97%)	137 (97%)	4 (3%)	38	58
1	C	141/146 (97%)	136 (96%)	5 (4%)	32	50
2	B	343/357 (96%)	334 (97%)	9 (3%)	40	60
2	D	344/357 (96%)	334 (97%)	10 (3%)	37	57
All	All	969/1006 (96%)	941 (97%)	28 (3%)	37	57

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU
1	A	57	GLU
1	A	71	ASP
1	A	93	LEU
2	B	3	SER
2	B	17	ASN
2	B	135	ARG
2	B	201	LEU
2	B	303	GLU
2	B	378	GLU
2	B	393	CYS
2	B	412	GLU
2	B	415	GLU
1	C	10	GLU
1	C	11	GLU
1	C	57	GLU
1	C	76	GLU
1	C	93	LEU
2	D	3	SER
2	D	17	ASN
2	D	106	ILE
2	D	135	ARG
2	D	201	LEU
2	D	237	GLU
2	D	269	ASN
2	D	303	GLU
2	D	364	LEU
2	D	393	CYS

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	39	ASN
2	B	41	GLN
2	B	83	ASN
2	B	220	ASN
2	B	269	ASN
1	C	17	GLN
1	C	39	ASN
2	D	269	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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	$\# Z > 2$
3	FES	A	200	1	0,4,4	-	-	-		
3	FES	C	200	1	0,4,4	-	-	-		
4	SF4	B	500	2	0,12,12	-	-	-		
4	SF4	D	500	2	0,12,12	-	-	-		
6	NAD	D	502	-	46,48,48	0.41	0	64,73,73	0.72	1 (1%)
6	NAD	B	502	-	46,48,48	0.42	0	64,73,73	0.76	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FMN	B	501	-	33,33,33	0.49	0	48,50,50	0.64	1 (2%)
5	FMN	D	501	-	33,33,33	0.37	0	48,50,50	0.51	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	FES	A	200	1	-	-	0/1/1/1
3	FES	C	200	1	-	-	0/1/1/1
4	SF4	B	500	2	-	-	0/6/5/5
4	SF4	D	500	2	-	-	0/6/5/5
6	NAD	D	502	-	-	6/30/62/62	0/5/5/5
6	NAD	B	502	-	-	3/30/62/62	0/5/5/5
5	FMN	B	501	-	-	4/18/18/18	0/3/3/3
5	FMN	D	501	-	-	5/18/18/18	0/3/3/3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	502	NAD	C4D-O4D-C1D	-3.64	106.59	109.92
6	B	502	NAD	C4D-O4D-C1D	-3.36	106.85	109.92
5	B	501	FMN	O5'-P-O1P	2.82	114.05	106.44

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	501	FMN	C1'-C2'-C3'-C4'
6	B	502	NAD	C5B-O5B-PA-O1A
6	B	502	NAD	O4D-C1D-N1N-C6N
6	D	502	NAD	O4D-C1D-N1N-C6N
5	D	501	FMN	O2'-C2'-C3'-O3'
5	D	501	FMN	O2'-C2'-C3'-C4'
6	D	502	NAD	PN-O3-PA-O1A
5	B	501	FMN	N10-C1'-C2'-C3'
6	D	502	NAD	C5B-O5B-PA-O1A
6	D	502	NAD	C5B-O5B-PA-O2A

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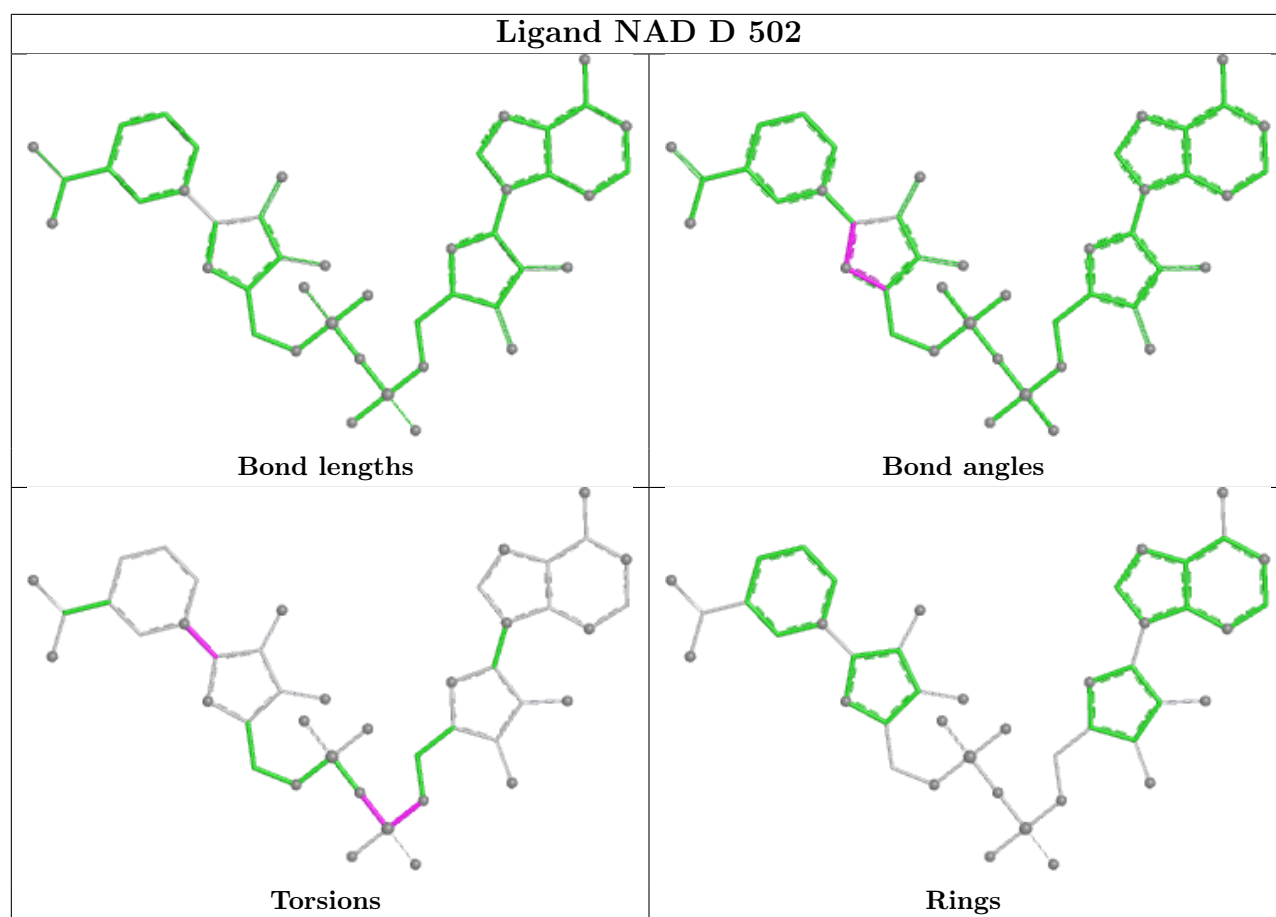
Mol	Chain	Res	Type	Atoms
6	D	502	NAD	C5B-O5B-PA-O3
5	B	501	FMN	C5'-O5'-P-O1P
6	B	502	NAD	O4D-C1D-N1N-C2N
6	D	502	NAD	O4D-C1D-N1N-C2N
5	B	501	FMN	C4'-C5'-O5'-P
5	D	501	FMN	C4'-C5'-O5'-P
5	D	501	FMN	C1'-C2'-C3'-O3'
5	B	501	FMN	N10-C1'-C2'-O2'

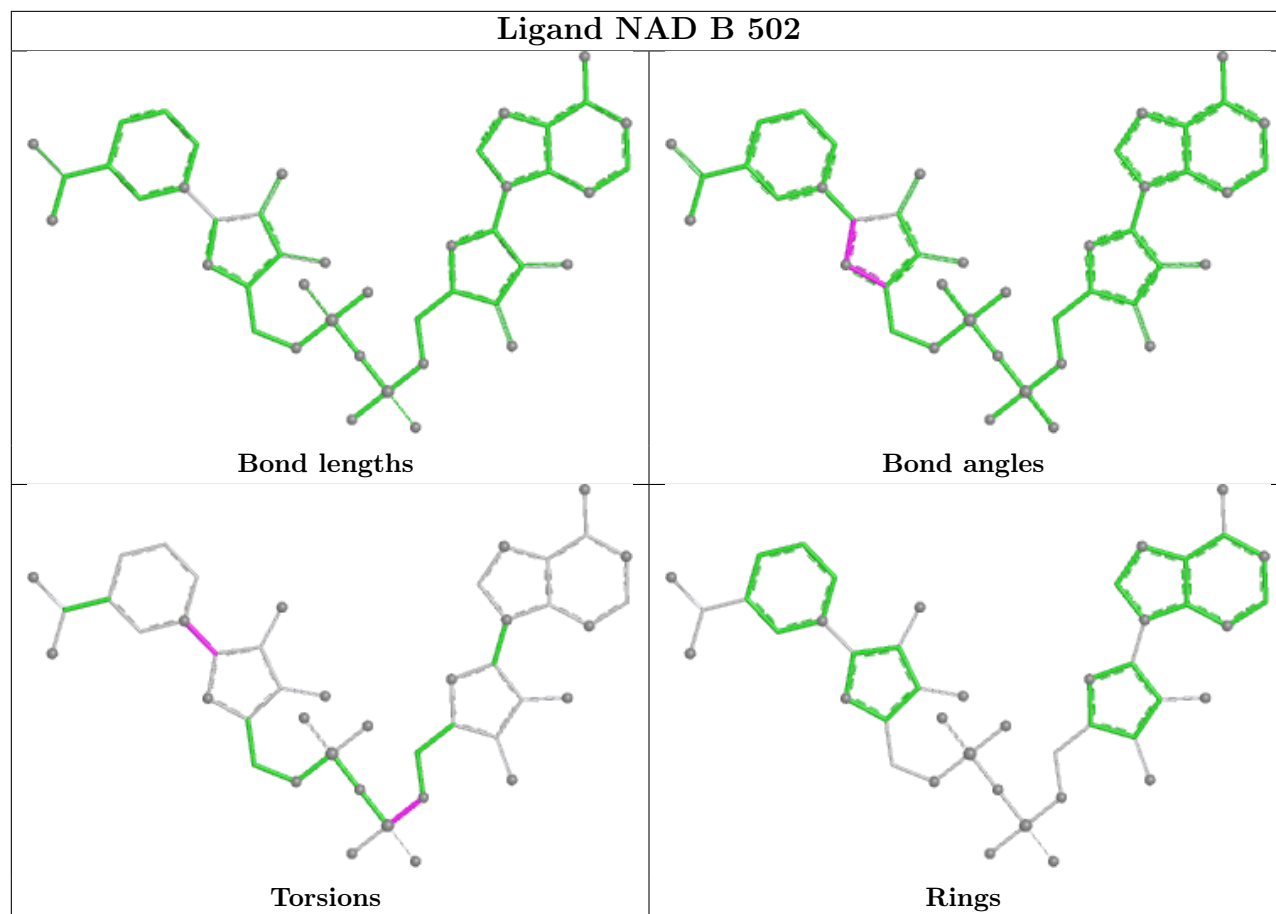
There are no ring outliers.

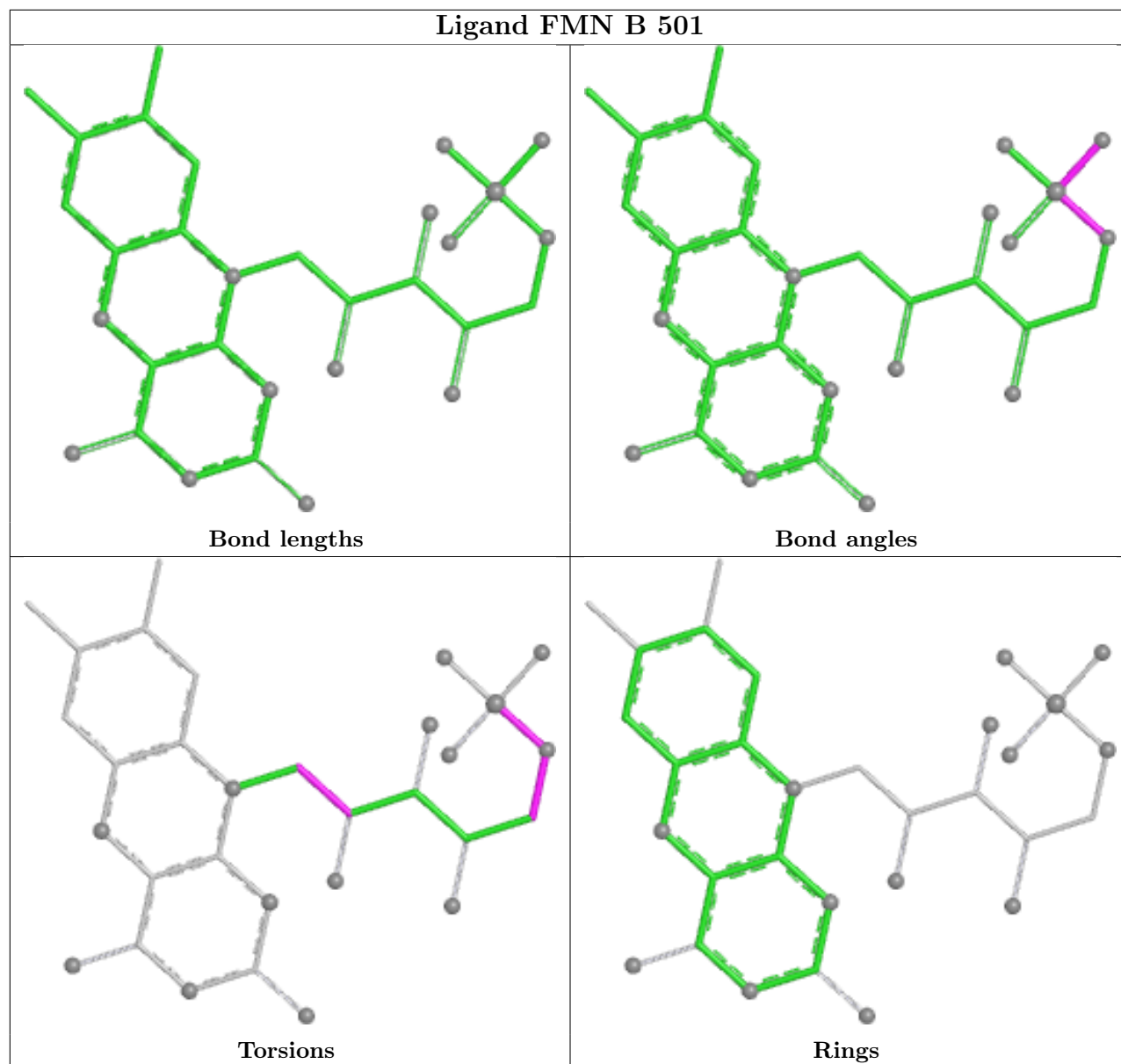
4 monomers are involved in 15 short contacts:

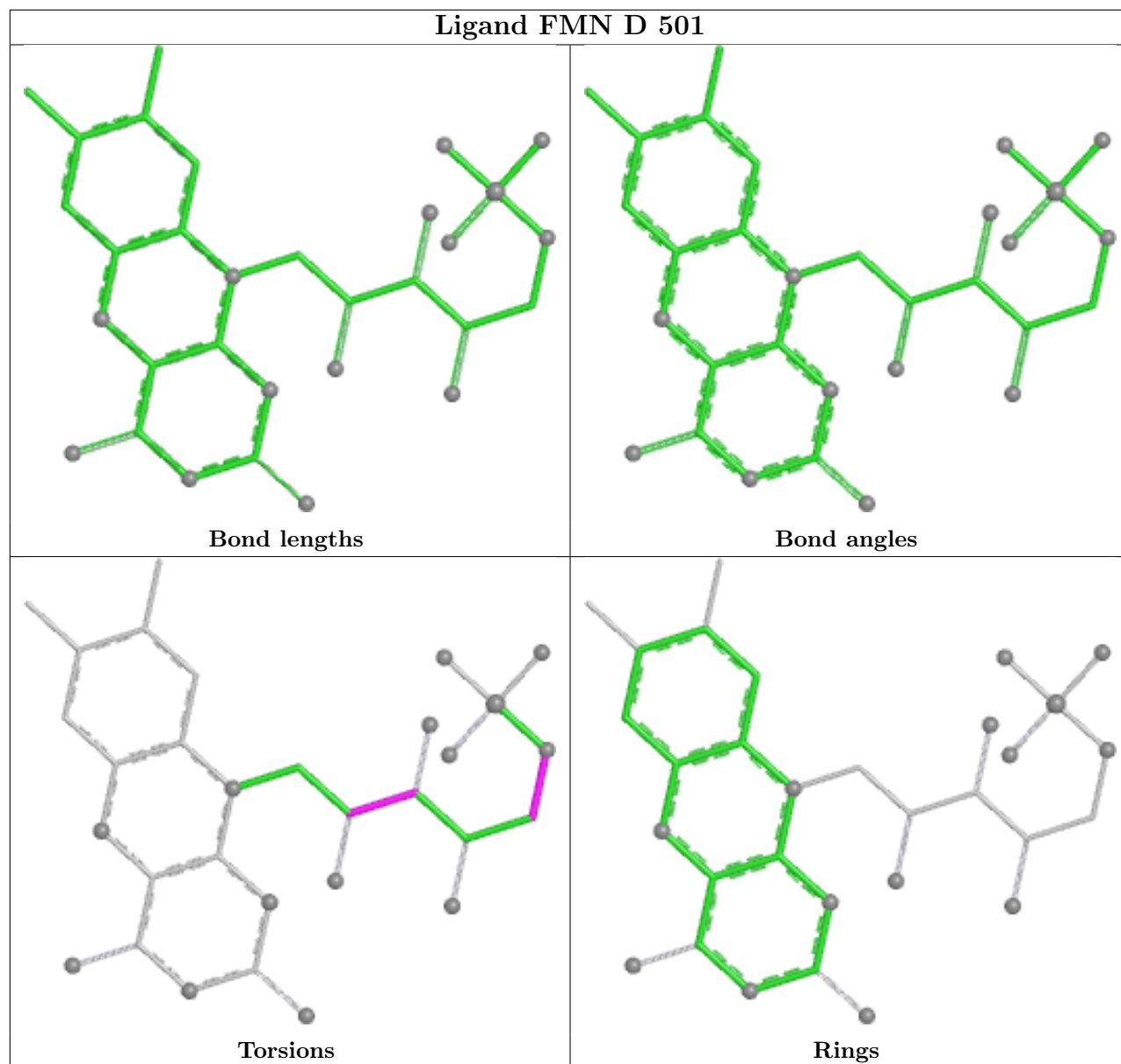
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	500	SF4	1	0
4	D	500	SF4	1	0
6	D	502	NAD	7	0
6	B	502	NAD	6	0

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	155/160 (96%)	0.46	1 (0%) 85 85	31, 49, 71, 84	0
1	C	155/160 (96%)	1.05	16 (10%) 12 10	37, 63, 90, 110	0
2	B	417/434 (96%)	0.40	22 (5%) 32 31	24, 44, 73, 117	0
2	D	418/434 (96%)	0.53	28 (6%) 24 22	23, 45, 74, 130	0
All	All	1145/1188 (96%)	0.54	67 (5%) 28 27	23, 48, 78, 130	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	395	LEU	5.3
2	D	393	CYS	4.7
2	D	240	TYR	4.4
2	D	357	CYS	4.1
2	B	379	GLY	3.9
2	D	394	GLY	3.6
2	D	380	PHE	3.6
2	B	407	LEU	3.5
2	B	418	ARG	3.4
1	A	87	VAL	3.3
2	D	377	TRP	3.3
2	D	244	SER	3.1
2	B	348	GLY	3.0
1	C	103	LEU	2.9
1	C	109	ILE	2.9
2	D	361	ALA	2.9
2	B	79	PHE	2.8
2	B	289	VAL	2.7
1	C	49	LEU	2.7
2	D	211	LEU	2.6
2	D	341	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
2	D	376	ASP	2.6
1	C	44	ILE	2.6
1	C	52	LEU	2.6
2	B	295	GLY	2.6
2	B	355	VAL	2.6
2	D	406	THR	2.6
1	C	107	LEU	2.6
2	D	314	LEU	2.6
2	B	68	GLY	2.6
2	D	347	CYS	2.5
2	B	416	LYS	2.5
2	D	419	LYS	2.5
1	C	120	PHE	2.5
2	B	235	TRP	2.5
2	D	355	VAL	2.5
1	C	46	PRO	2.5
1	C	12	LEU	2.4
1	C	108	GLY	2.4
2	D	313	PRO	2.3
2	B	351	THR	2.3
2	D	319	THR	2.3
2	D	201	LEU	2.3
2	D	93	ALA	2.3
2	B	414	TRP	2.3
2	B	296	ALA	2.3
2	B	390	THR	2.3
2	B	392	ILE	2.3
2	B	69	ALA	2.2
1	C	14	THR	2.2
2	B	383	VAL	2.2
1	C	8	PHE	2.2
2	B	395	LEU	2.2
2	B	382	PHE	2.2
2	D	192	LEU	2.1
2	D	387	ILE	2.1
2	D	339	ALA	2.1
1	C	78	LYS	2.1
2	D	202	LYS	2.1
1	C	102	ALA	2.1
1	C	123	VAL	2.1
2	B	291	ALA	2.1
2	B	384	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	6	PHE	2.1
2	D	188	LEU	2.0
2	D	84	PRO	2.0
2	D	396	GLY	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

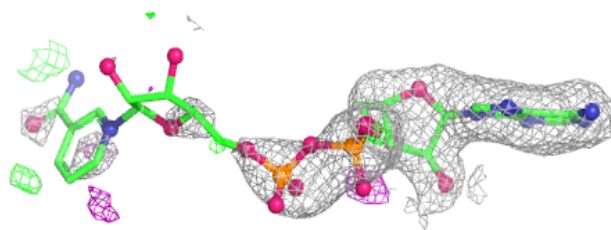
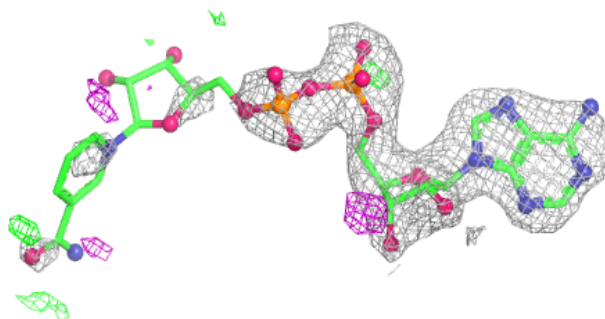
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
6	NAD	D	502	44/44	0.83	0.17	39,54,66,72	26
6	NAD	B	502	44/44	0.89	0.14	23,33,45,47	26
4	SF4	D	500	8/8	0.96	0.10	44,49,51,52	0
5	FMN	D	501	31/31	0.96	0.08	26,39,42,46	0
5	FMN	B	501	31/31	0.97	0.07	27,37,40,43	0
4	SF4	B	500	8/8	0.98	0.05	43,46,50,51	0
3	FES	C	200	4/4	0.98	0.04	28,37,38,40	0
3	FES	A	200	4/4	0.99	0.03	39,40,41,42	0

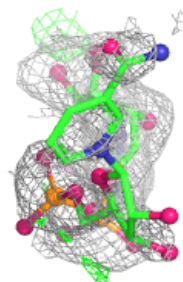
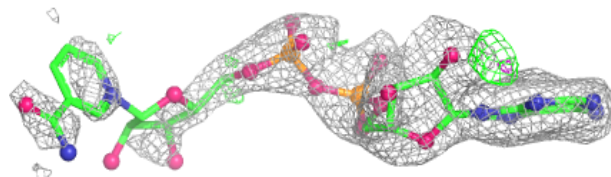
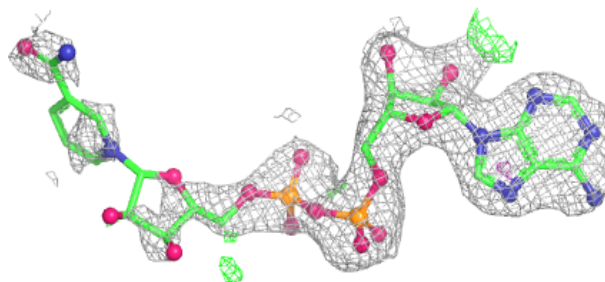
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD 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)

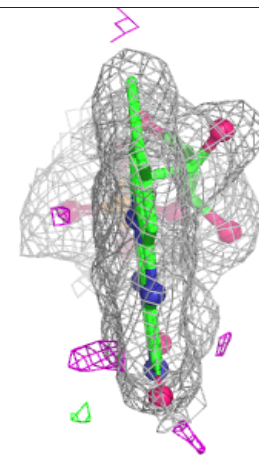
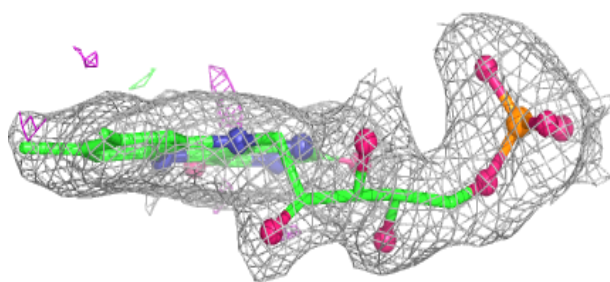
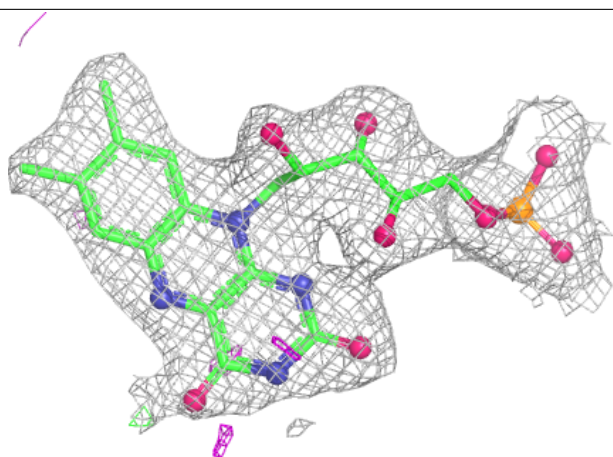
**Electron density around NAD B 502:**

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

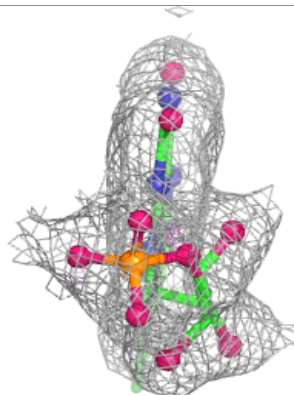
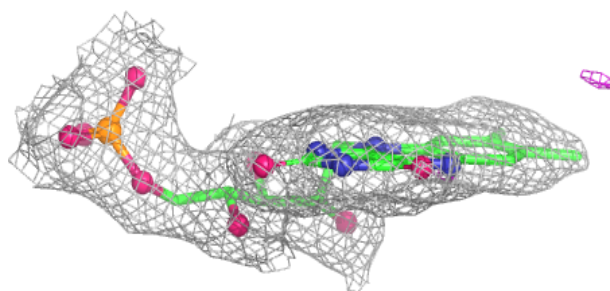
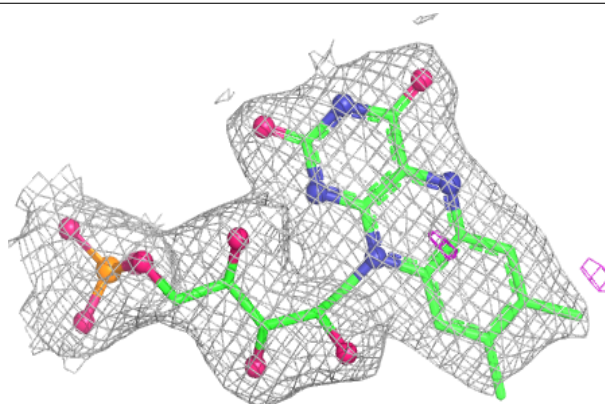


Electron density around FMN D 501:

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

**Electron density around FMN 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.