



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:12 PM UTC

PDB ID : 5CPF / pdb_00005cpf
Title : Compensation of the effect of isoleucine to alanine mutation by designed inhibition in the InhA enzyme
Authors : Li, H.-J.; Lai, C.-T.; Pan, P.; Yu, W.; Shah, S.; Bommineni, G.R.; Perrone, V.; Garcia-Diaz, M.; Tonge, P.J.; Simmerling, C.
Deposited on : 2015-07-21
Resolution : 3.41 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

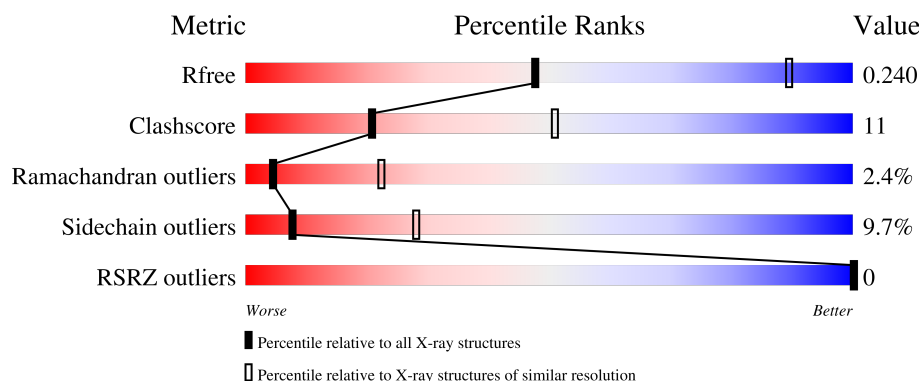
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.41 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	289	
1	B	289	
1	C	289	
1	D	289	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8002 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 Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			1916	1210	339	358	9			
1	B	267	Total	C	N	O	S	0	0	0
			1940	1232	337	361	10			
1	C	267	Total	C	N	O	S	0	0	0
			1919	1213	339	358	9			
1	D	267	Total	C	N	O	S	0	0	0
			1943	1233	338	362	10			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P9WGR0
A	-18	GLY	-	expression tag	UNP P9WGR0
A	-17	SER	-	expression tag	UNP P9WGR0
A	-16	SER	-	expression tag	UNP P9WGR0
A	-15	HIS	-	expression tag	UNP P9WGR0
A	-14	HIS	-	expression tag	UNP P9WGR0
A	-13	HIS	-	expression tag	UNP P9WGR0
A	-12	HIS	-	expression tag	UNP P9WGR0
A	-11	HIS	-	expression tag	UNP P9WGR0
A	-10	HIS	-	expression tag	UNP P9WGR0
A	-9	SER	-	expression tag	UNP P9WGR0
A	-8	SER	-	expression tag	UNP P9WGR0
A	-7	GLY	-	expression tag	UNP P9WGR0
A	-6	LEU	-	expression tag	UNP P9WGR0
A	-5	VAL	-	expression tag	UNP P9WGR0
A	-4	PRO	-	expression tag	UNP P9WGR0
A	-3	ARG	-	expression tag	UNP P9WGR0
A	-2	GLY	-	expression tag	UNP P9WGR0
A	-1	SER	-	expression tag	UNP P9WGR0
A	0	HIS	-	expression tag	UNP P9WGR0
A	215	ALA	ILE	engineered mutation	UNP P9WGR0

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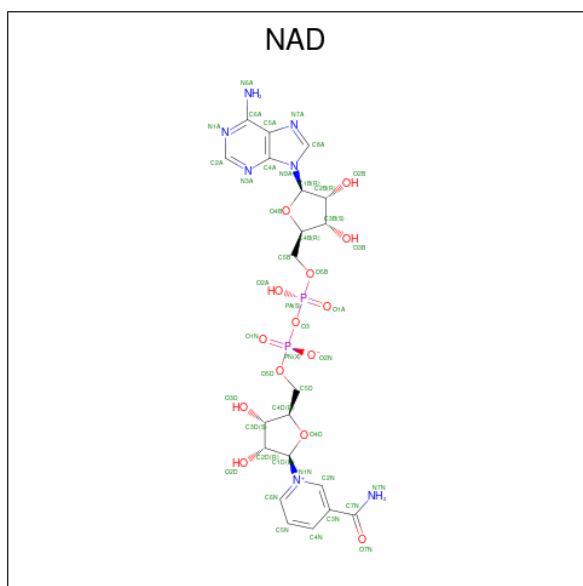
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP P9WGR0
B	-18	GLY	-	expression tag	UNP P9WGR0
B	-17	SER	-	expression tag	UNP P9WGR0
B	-16	SER	-	expression tag	UNP P9WGR0
B	-15	HIS	-	expression tag	UNP P9WGR0
B	-14	HIS	-	expression tag	UNP P9WGR0
B	-13	HIS	-	expression tag	UNP P9WGR0
B	-12	HIS	-	expression tag	UNP P9WGR0
B	-11	HIS	-	expression tag	UNP P9WGR0
B	-10	HIS	-	expression tag	UNP P9WGR0
B	-9	SER	-	expression tag	UNP P9WGR0
B	-8	SER	-	expression tag	UNP P9WGR0
B	-7	GLY	-	expression tag	UNP P9WGR0
B	-6	LEU	-	expression tag	UNP P9WGR0
B	-5	VAL	-	expression tag	UNP P9WGR0
B	-4	PRO	-	expression tag	UNP P9WGR0
B	-3	ARG	-	expression tag	UNP P9WGR0
B	-2	GLY	-	expression tag	UNP P9WGR0
B	-1	SER	-	expression tag	UNP P9WGR0
B	0	HIS	-	expression tag	UNP P9WGR0
B	215	ALA	ILE	engineered mutation	UNP P9WGR0
C	-19	MET	-	initiating methionine	UNP P9WGR0
C	-18	GLY	-	expression tag	UNP P9WGR0
C	-17	SER	-	expression tag	UNP P9WGR0
C	-16	SER	-	expression tag	UNP P9WGR0
C	-15	HIS	-	expression tag	UNP P9WGR0
C	-14	HIS	-	expression tag	UNP P9WGR0
C	-13	HIS	-	expression tag	UNP P9WGR0
C	-12	HIS	-	expression tag	UNP P9WGR0
C	-11	HIS	-	expression tag	UNP P9WGR0
C	-10	HIS	-	expression tag	UNP P9WGR0
C	-9	SER	-	expression tag	UNP P9WGR0
C	-8	SER	-	expression tag	UNP P9WGR0
C	-7	GLY	-	expression tag	UNP P9WGR0
C	-6	LEU	-	expression tag	UNP P9WGR0
C	-5	VAL	-	expression tag	UNP P9WGR0
C	-4	PRO	-	expression tag	UNP P9WGR0
C	-3	ARG	-	expression tag	UNP P9WGR0
C	-2	GLY	-	expression tag	UNP P9WGR0
C	-1	SER	-	expression tag	UNP P9WGR0
C	0	HIS	-	expression tag	UNP P9WGR0
C	215	ALA	ILE	engineered mutation	UNP P9WGR0

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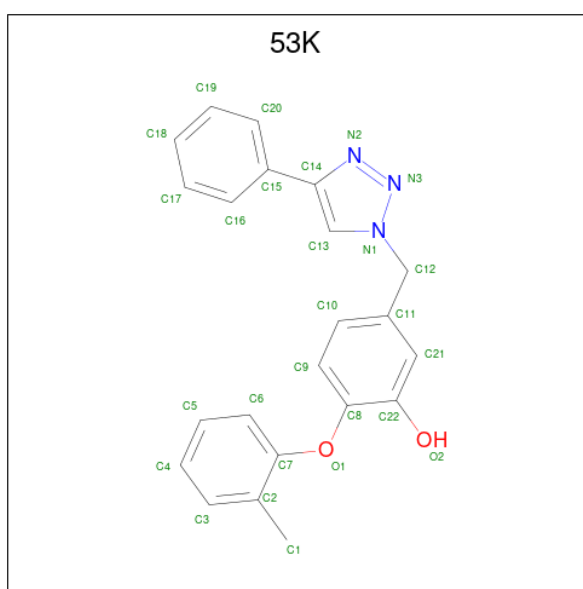
Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	initiating methionine	UNP P9WGR0
D	-18	GLY	-	expression tag	UNP P9WGR0
D	-17	SER	-	expression tag	UNP P9WGR0
D	-16	SER	-	expression tag	UNP P9WGR0
D	-15	HIS	-	expression tag	UNP P9WGR0
D	-14	HIS	-	expression tag	UNP P9WGR0
D	-13	HIS	-	expression tag	UNP P9WGR0
D	-12	HIS	-	expression tag	UNP P9WGR0
D	-11	HIS	-	expression tag	UNP P9WGR0
D	-10	HIS	-	expression tag	UNP P9WGR0
D	-9	SER	-	expression tag	UNP P9WGR0
D	-8	SER	-	expression tag	UNP P9WGR0
D	-7	GLY	-	expression tag	UNP P9WGR0
D	-6	LEU	-	expression tag	UNP P9WGR0
D	-5	VAL	-	expression tag	UNP P9WGR0
D	-4	PRO	-	expression tag	UNP P9WGR0
D	-3	ARG	-	expression tag	UNP P9WGR0
D	-2	GLY	-	expression tag	UNP P9WGR0
D	-1	SER	-	expression tag	UNP P9WGR0
D	0	HIS	-	expression tag	UNP P9WGR0
D	215	ALA	ILE	engineered mutation	UNP P9WGR0

- 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	
			44	21	7	14	2	
2	B	1	Total	C	N	O	P	
			44	21	7	14	2	
2	C	1	Total	C	N	O	P	
			44	21	7	14	2	
2	D	1	Total	C	N	O	P	
			44	21	7	14	2	

- Molecule 3 is 2-(2-methylphenoxy)-5-[(4-phenyl-1H-1,2,3-triazol-1-yl)methyl]phenol (CCD ID: 53K) (formula: C₂₂H₁₉N₃O₂).

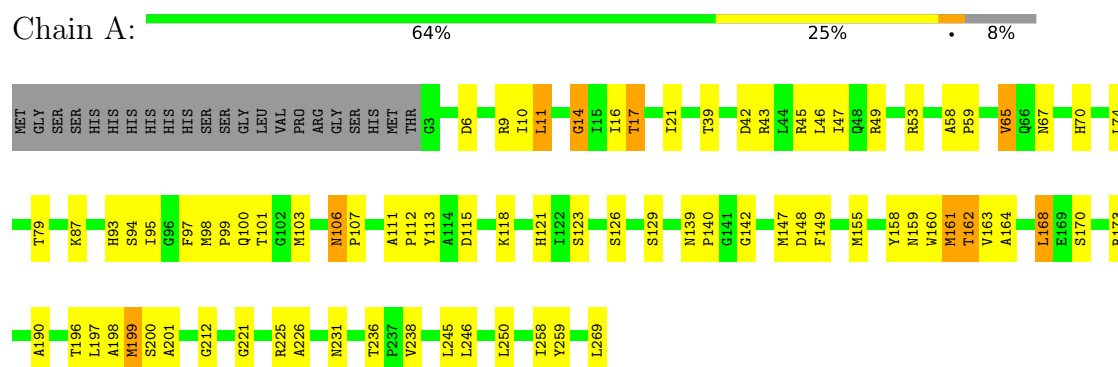


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O		
			27	22	3	2	0	0
3	B	1	Total	C	N	O		
			27	22	3	2	0	0
3	C	1	Total	C	N	O		
			27	22	3	2	0	0
3	D	1	Total	C	N	O		
			27	22	3	2	0	0

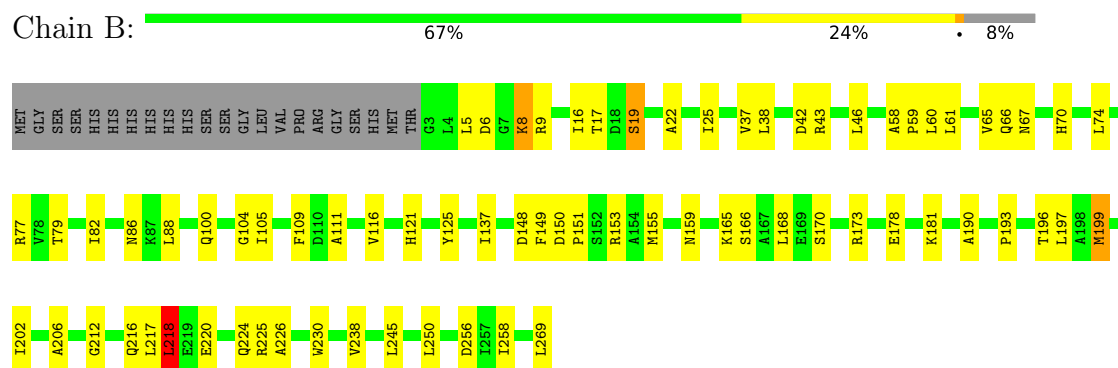
3 Residue-property plots [i](#)

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.

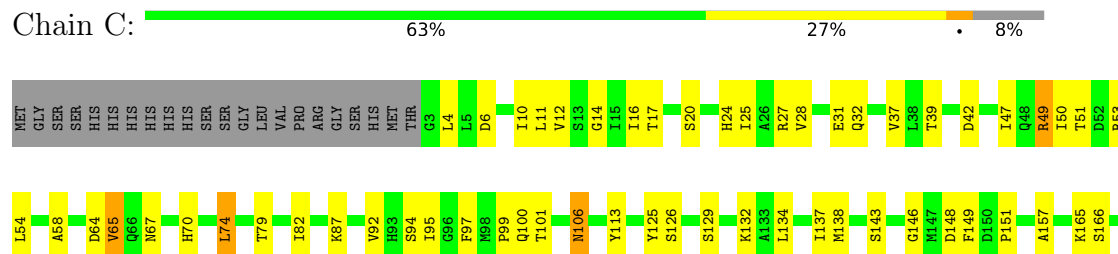
• Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



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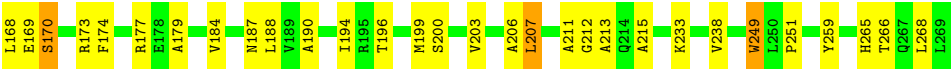
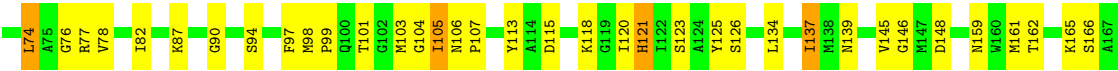
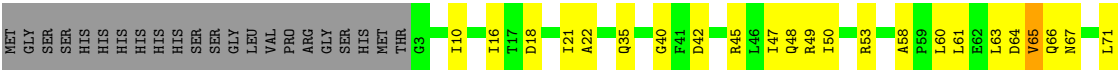


• Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]





● Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.01Å 97.56Å 184.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.78 – 3.41 48.78 – 3.41	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.78-3.41) 99.4 (48.78-3.41)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 3.40Å)	Xtriage
Refinement program	PHENIX 1.7.2 _869	Depositor
R, R_{free}	0.182 , 0.250 0.175 , 0.240	Depositor DCC
R_{free} test set	1196 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	71.1	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8002	wwPDB-VP
Average B, all atoms (Å ²)	59.0	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 39.19 % 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 3.2943e-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: 53K, 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	0.57	0/1954	1.00	4/2659 (0.2%)
1	B	0.56	0/1978	0.99	4/2689 (0.1%)
1	C	0.59	0/1957	1.02	7/2665 (0.3%)
1	D	0.58	0/1981	0.98	4/2694 (0.1%)
All	All	0.58	0/7870	1.00	19/10707 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	175	VAL	N-CA-C	-6.06	103.80	112.35
1	B	212	GLY	N-CA-C	-6.03	107.18	114.66
1	C	58	ALA	CA-C-N	5.97	125.93	119.78
1	C	58	ALA	C-N-CA	5.97	125.93	119.78
1	B	19	SER	N-CA-C	-5.86	106.75	114.31
1	A	14	GLY	N-CA-C	5.86	123.92	114.90
1	A	226	ALA	CA-C-N	5.53	125.36	119.28
1	A	226	ALA	C-N-CA	5.53	125.36	119.28
1	A	155	MET	N-CA-C	5.47	113.14	108.22
1	D	249	TRP	N-CA-C	5.43	119.77	113.20
1	C	106	ASN	CA-C-N	5.28	125.27	119.89
1	C	106	ASN	C-N-CA	5.28	125.27	119.89
1	D	58	ALA	CA-C-N	5.27	125.27	119.90
1	D	58	ALA	C-N-CA	5.27	125.27	119.90
1	B	226	ALA	CA-C-N	5.22	125.02	119.28
1	B	226	ALA	C-N-CA	5.22	125.02	119.28
1	D	18	ASP	N-CA-C	-5.21	106.88	113.18
1	C	236	THR	CA-C-N	-5.18	113.20	118.85
1	C	236	THR	C-N-CA	-5.18	113.20	118.85

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	1916	0	1855	48	1
1	B	1940	0	1924	33	0
1	C	1919	0	1866	51	1
1	D	1943	0	1927	57	0
2	A	44	0	26	3	0
2	B	44	0	26	0	0
2	C	44	0	26	2	0
2	D	44	0	26	2	0
3	A	27	0	18	4	0
3	B	27	0	18	2	0
3	C	27	0	18	4	0
3	D	27	0	18	2	0
All	All	8002	0	7748	175	1

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 11.

All (175) 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:C:99:PRO:O	1:C:101:THR:N	2.18	0.77
1:A:158:TYR:HD2	1:A:162:THR:HG1	1.34	0.76
1:D:207:LEU:HB3	1:D:211:ALA:HB2	1.70	0.74
1:C:16:ILE:HG23	1:C:17:THR:HG23	1.72	0.72
1:B:155:MET:HE2	3:B:301:53K:H6	1.74	0.70
1:A:98:MET:HE2	1:A:161:MET:HE3	1.75	0.68
1:D:65:VAL:HG21	1:D:126:SER:HB2	1.78	0.66
1:C:99:PRO:C	1:C:101:THR:H	2.04	0.64
2:D:302:NAD:H52N	2:D:302:NAD:H52A	1.79	0.63
1:D:48:GLN:HG2	1:D:60:LEU:HD23	1.78	0.63
1:C:143:SER:HB2	1:C:185:ARG:NH2	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ILE:HD13	1:A:246:LEU:HD13	1.80	0.62
1:B:67:ASN:CG	1:B:70:HIS:HD2	2.08	0.62
1:B:217:LEU:O	1:B:218:LEU:HB2	1.99	0.62
1:D:49:ARG:HH21	1:D:53:ARG:HH12	1.48	0.62
1:D:148:ASP:HA	1:D:165:LYS:HD2	1.82	0.61
1:B:61:LEU:HD22	1:B:77:ARG:HB3	1.81	0.61
1:C:227:PRO:HB3	1:D:177:ARG:HA	1.84	0.60
1:A:196:THR:O	1:A:198:ALA:N	2.34	0.60
1:C:126:SER:HA	1:C:129:SER:HB2	1.82	0.60
1:A:126:SER:HA	1:A:129:SER:HB2	1.84	0.59
1:A:245:LEU:HD11	1:A:258:ILE:HD12	1.84	0.59
1:C:157:ALA:HB3	3:C:301:53K:H18	1.84	0.58
1:A:97:PHE:HA	3:A:301:53K:H10	1.84	0.58
1:D:148:ASP:O	1:D:190:ALA:HA	2.04	0.58
1:B:16:ILE:HG23	1:B:17:THR:HG23	1.86	0.58
1:B:196:THR:H	1:B:199:MET:HB3	1.69	0.57
2:C:300:NAD:O2D	3:C:301:53K:O2	2.17	0.57
1:D:97:PHE:HE2	1:D:118:LYS:HD3	1.70	0.56
1:A:159:ASN:HA	1:D:174:PHE:CE1	2.39	0.56
1:C:193:PRO:HD2	1:C:230:TRP:NE1	2.20	0.56
1:A:221:GLY:O	1:A:225:ARG:HG3	2.06	0.55
1:D:179:ALA:HB1	1:D:184:VAL:HB	1.89	0.55
2:A:300:NAD:O2D	3:A:301:53K:O2	2.12	0.55
1:D:97:PHE:H	1:D:123:SER:HB3	1.71	0.55
1:B:148:ASP:OD1	1:B:149:PHE:N	2.40	0.55
1:D:66:GLN:HG2	1:D:121:HIS:CE1	2.41	0.55
1:B:109:PHE:HB3	1:C:132:LYS:HD2	1.88	0.55
1:B:148:ASP:HA	1:B:165:LYS:HD2	1.89	0.55
1:A:16:ILE:HG23	1:A:17:THR:HG23	1.89	0.54
1:A:199:MET:HB2	3:A:301:53K:H4	1.89	0.54
1:B:178:GLU:HA	1:B:181:LYS:HE2	1.88	0.54
1:C:148:ASP:HA	1:C:165:LYS:HE3	1.89	0.54
1:C:148:ASP:OD1	1:C:149:PHE:N	2.41	0.54
1:C:178:GLU:OE1	1:C:181:LYS:NZ	2.34	0.54
1:B:22:ALA:HA	1:B:25:ILE:HD12	1.90	0.53
1:B:100:GLN:HG2	1:B:206:ALA:HB2	1.90	0.53
1:D:99:PRO:HG2	1:D:115:ASP:HB3	1.89	0.53
1:D:194:ILE:N	2:D:302:NAD:O7N	2.33	0.53
1:C:47:ILE:O	1:C:51:THR:HG23	2.09	0.53
1:A:103:MET:SD	3:A:301:53K:H2	2.48	0.53
1:B:199:MET:HE1	3:B:301:53K:N2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:194:ILE:O	1:D:196:THR:HG23	2.10	0.52
1:C:148:ASP:O	1:C:190:ALA:HA	2.09	0.52
1:D:10:ILE:HG12	1:D:90:GLY:HA3	1.91	0.52
1:D:50:ILE:O	1:D:53:ARG:HB2	2.09	0.52
1:C:240:LYS:HD3	1:D:251:PRO:HG3	1.92	0.52
1:A:21:ILE:HG21	1:A:147:MET:HE2	1.92	0.51
1:A:11:LEU:HD11	1:A:39:THR:HG23	1.91	0.51
1:A:163:VAL:HG22	1:D:170:SER:HB3	1.92	0.51
1:D:104:GLY:HA3	1:D:206:ALA:O	2.10	0.51
1:A:99:PRO:C	1:A:101:THR:H	2.19	0.51
1:C:143:SER:HB2	1:C:185:ARG:HH21	1.75	0.51
1:A:162:THR:HG21	1:D:174:PHE:HE2	1.76	0.51
1:D:65:VAL:CG2	1:D:126:SER:HB2	2.41	0.50
1:B:220:GLU:O	1:B:224:GLN:HB2	2.12	0.50
1:C:11:LEU:HD11	1:C:39:THR:HG23	1.93	0.50
1:A:67:ASN:CG	1:A:70:HIS:HD2	2.20	0.50
1:D:64:ASP:HB3	1:D:67:ASN:HB2	1.93	0.50
1:A:14:GLY:HA3	1:A:94:SER:O	2.13	0.49
1:A:199:MET:O	1:A:201:ALA:N	2.45	0.49
1:A:161:MET:HE2	1:A:164:ALA:HB3	1.94	0.49
1:C:27:ARG:O	1:C:31:GLU:HG3	2.12	0.49
1:A:123:SER:OG	1:A:161:MET:HE1	2.12	0.49
1:A:158:TYR:HD2	1:A:162:THR:OG1	1.95	0.49
1:B:151:PRO:O	1:C:173:ARG:NH1	2.46	0.49
1:C:64:ASP:HB3	1:C:67:ASN:HB2	1.95	0.49
1:C:240:LYS:HG2	1:D:249:TRP:CZ3	2.48	0.49
1:C:194:ILE:HG22	1:C:196:THR:HG23	1.95	0.48
1:A:113:TYR:CZ	1:D:121:HIS:HB2	2.48	0.48
1:C:221:GLY:O	1:C:225:ARG:HG3	2.13	0.48
1:D:203:VAL:O	1:D:207:LEU:HB2	2.13	0.48
1:D:87:LYS:HD2	1:D:137:ILE:C	2.39	0.48
1:D:165:LYS:O	1:D:169:GLU:HG3	2.14	0.48
1:B:148:ASP:O	1:B:190:ALA:HA	2.14	0.48
1:C:193:PRO:HD2	1:C:230:TRP:HE1	1.78	0.48
1:A:112:PRO:HD2	1:A:115:ASP:HB2	1.96	0.47
1:D:97:PHE:HA	3:D:301:53K:H10	1.96	0.47
1:D:40:GLY:HA3	1:D:47:ILE:HD13	1.96	0.47
1:D:66:GLN:HE22	1:D:118:LYS:HG3	1.79	0.47
1:D:121:HIS:HA	1:D:125:TYR:HB3	1.97	0.47
1:D:145:VAL:HA	1:D:187:ASN:O	2.15	0.47
1:A:113:TYR:CE2	1:D:121:HIS:HB2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:VAL:O	1:D:82:ILE:HG12	2.14	0.47
1:C:4:LEU:HB3	1:C:32:GLN:HB3	1.96	0.46
1:D:63:LEU:HD11	1:D:71:LEU:HD23	1.98	0.46
1:B:70:HIS:O	1:B:77:ARG:NH2	2.48	0.46
1:C:205:GLY:HA2	1:C:208:GLY:O	2.16	0.46
1:A:65:VAL:HB	2:A:300:NAD:N1A	2.30	0.46
1:B:9:ARG:NH2	1:B:86:ASN:HB3	2.31	0.46
1:A:106:ASN:HA	1:A:107:PRO:HD3	1.89	0.45
1:C:256:ASP:OD2	1:D:259:TYR:HB2	2.16	0.45
1:D:211:ALA:C	1:D:213:ALA:H	2.24	0.45
1:A:170:SER:HA	1:A:173:ARG:NH1	2.31	0.45
1:C:70:HIS:O	1:C:74:LEU:HB2	2.16	0.45
1:B:245:LEU:HD11	1:B:258:ILE:HG13	1.99	0.45
1:C:14:GLY:HA3	1:C:94:SER:O	2.16	0.45
1:A:245:LEU:HD21	1:A:258:ILE:HD11	1.98	0.45
1:D:162:THR:O	1:D:166:SER:HB2	2.17	0.45
1:A:99:PRO:O	1:A:101:THR:N	2.50	0.44
1:D:22:ALA:HB2	1:D:94:SER:HB3	1.99	0.44
1:D:74:LEU:HD13	1:D:134:LEU:HD21	1.99	0.44
1:D:212:GLY:HA2	1:D:215:ALA:HB3	1.98	0.44
1:C:173:ARG:HD2	1:D:266:THR:HB	1.99	0.44
1:C:266:THR:HB	1:D:173:ARG:HD2	1.98	0.44
1:C:4:LEU:HD11	1:C:247:SER:HB2	1.99	0.44
1:C:65:VAL:HB	2:C:300:NAD:N1A	2.32	0.44
1:C:97:PHE:HA	3:C:301:53K:H10	1.98	0.44
1:A:93:HIS:O	1:A:93:HIS:ND1	2.50	0.44
1:A:148:ASP:OD1	1:A:149:PHE:N	2.51	0.44
1:A:190:ALA:HB3	1:A:259:TYR:CD2	2.53	0.44
1:B:88:LEU:HD23	1:B:88:LEU:HA	1.68	0.44
1:D:106:ASN:HA	1:D:107:PRO:HD3	1.91	0.44
1:A:159:ASN:O	1:A:162:THR:HB	2.18	0.44
1:D:87:LYS:HB2	1:D:139:ASN:OD1	2.18	0.43
1:A:98:MET:HE3	1:A:160:TRP:C	2.42	0.43
2:A:300:NAD:H52A	2:A:300:NAD:H52N	2.00	0.43
1:D:146:GLY:HA3	1:D:188:LEU:HD23	2.00	0.43
1:A:168:LEU:HD22	1:A:168:LEU:O	2.18	0.43
1:B:5:LEU:O	1:B:8:LYS:HB2	2.19	0.43
1:C:134:LEU:HD23	1:C:137:ILE:HD11	2.00	0.43
1:D:199:MET:HE2	3:D:301:53K:N3	2.34	0.43
1:B:104:GLY:HA3	1:B:206:ALA:O	2.19	0.43
1:B:153:ARG:NE	1:D:265:HIS:O	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:GLY:O	1:C:188:LEU:HA	2.19	0.43
1:D:98:MET:HE3	1:D:161:MET:HA	2.00	0.43
1:A:121:HIS:HB2	1:D:113:TYR:CE2	2.54	0.42
1:B:38:LEU:O	1:B:60:LEU:HA	2.18	0.42
1:A:43:ARG:O	1:A:47:ILE:HG13	2.19	0.42
1:C:24:HIS:O	1:C:28:VAL:HG23	2.19	0.42
1:A:67:ASN:HB3	1:A:70:HIS:HB2	2.00	0.42
1:C:138:MET:HE2	1:C:138:MET:HB3	1.91	0.42
1:B:66:GLN:HG2	1:B:121:HIS:CE1	2.53	0.42
1:B:67:ASN:HB3	1:B:70:HIS:CD2	2.54	0.42
1:C:25:ILE:HG21	1:C:92:VAL:HG11	2.02	0.42
1:B:121:HIS:HB2	1:C:113:TYR:CZ	2.55	0.42
1:D:22:ALA:HB2	1:D:94:SER:CB	2.50	0.41
1:A:111:ALA:HA	1:A:112:PRO:HD3	1.78	0.41
1:C:245:LEU:HD21	1:C:258:ILE:HG13	2.02	0.41
1:D:103:MET:HE2	1:D:203:VAL:HG22	2.01	0.41
1:D:105:ILE:HG23	1:D:207:LEU:HD23	2.03	0.41
1:A:250:LEU:HD23	1:A:250:LEU:HA	1.77	0.41
1:B:111:ALA:O	1:C:125:TYR:OH	2.37	0.41
1:C:10:ILE:HG22	1:C:12:VAL:HG23	2.03	0.41
1:D:199:MET:HE3	1:D:199:MET:HB2	1.92	0.41
1:A:140:PRO:C	1:A:142:GLY:H	2.29	0.41
1:B:193:PRO:HD2	1:B:230:TRP:CD1	2.55	0.41
1:C:196:THR:H	1:C:199:MET:HB3	1.86	0.41
1:C:247:SER:OG	1:C:248:ASP:N	2.54	0.41
1:D:268:LEU:HA	1:D:268:LEU:HD12	1.67	0.41
1:B:58:ALA:HA	1:B:59:PRO:HD3	1.94	0.41
1:A:97:PHE:HE2	1:A:118:LYS:HD3	1.86	0.41
1:B:125:TYR:CE1	1:C:113:TYR:HB2	2.56	0.41
1:B:218:LEU:HD12	1:B:218:LEU:HA	1.90	0.41
1:C:51:THR:HA	1:C:54:LEU:HD12	2.02	0.41
1:C:87:LYS:HG2	1:C:137:ILE:C	2.46	0.41
1:A:87:LYS:HB2	1:A:139:ASN:OD1	2.21	0.40
1:C:157:ALA:HB3	3:C:301:53K:C18	2.50	0.40
1:A:161:MET:O	1:A:161:MET:HG3	2.21	0.40
1:C:250:LEU:HD23	1:C:250:LEU:HA	1.93	0.40
1:B:170:SER:HA	1:B:173:ARG:NH1	2.36	0.40
1:A:93:HIS:CE1	1:A:95:ILE:HB	2.57	0.40
1:C:11:LEU:HD12	1:C:37:VAL:O	2.22	0.40
1:A:58:ALA:HA	1:A:59:PRO:HD3	1.84	0.40
1:D:49:ARG:HH21	1:D:53:ARG:NH1	2.16	0.40

All (1) 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)
1:A:212:GLY:O	1:C:49:ARG:NH2[3_555]	2.16	0.04

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	265/289 (92%)	236 (89%)	23 (9%)	6 (2%)	5	23
1	B	265/289 (92%)	234 (88%)	23 (9%)	8 (3%)	3	18
1	C	265/289 (92%)	244 (92%)	17 (6%)	4 (2%)	8	30
1	D	265/289 (92%)	244 (92%)	14 (5%)	7 (3%)	4	21
All	All	1060/1156 (92%)	958 (90%)	77 (7%)	25 (2%)	4	22

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	8	LYS
1	B	218	LEU
1	C	100	GLN
1	A	100	GLN
1	A	197	LEU
1	A	200	SER
1	B	159	ASN
1	B	216	GLN
1	C	199	MET
1	D	120	ILE
1	D	121	HIS
1	A	6	ASP
1	A	42	ASP
1	D	16	ILE
1	D	159	ASN

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Mol	Chain	Res	Type
1	A	199	MET
1	B	42	ASP
1	D	42	ASP
1	B	150	ASP
1	C	42	ASP
1	C	151	PRO
1	D	76	GLY
1	B	137	ILE
1	D	137	ILE
1	B	202	ILE

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	180/221 (81%)	162 (90%)	18 (10%)	7	26
1	B	190/221 (86%)	169 (89%)	21 (11%)	6	21
1	C	181/221 (82%)	163 (90%)	18 (10%)	7	26
1	D	191/221 (86%)	176 (92%)	15 (8%)	11	36
All	All	742/884 (84%)	670 (90%)	72 (10%)	8	28

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	11	LEU
1	A	17	THR
1	A	45	ARG
1	A	46	LEU
1	A	49	ARG
1	A	53	ARG
1	A	65	VAL
1	A	74	LEU
1	A	79	THR
1	A	106	ASN

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Mol	Chain	Res	Type
1	A	161	MET
1	A	162	THR
1	A	168	LEU
1	A	231	ASN
1	A	236	THR
1	A	238	VAL
1	A	269	LEU
1	B	6	ASP
1	B	19	SER
1	B	37	VAL
1	B	43	ARG
1	B	46	LEU
1	B	65	VAL
1	B	74	LEU
1	B	79	THR
1	B	82	ILE
1	B	105	ILE
1	B	116	VAL
1	B	166	SER
1	B	168	LEU
1	B	197	LEU
1	B	199	MET
1	B	218	LEU
1	B	225	ARG
1	B	238	VAL
1	B	250	LEU
1	B	256	ASP
1	B	269	LEU
1	C	6	ASP
1	C	20	SER
1	C	49	ARG
1	C	50	ILE
1	C	53	ARG
1	C	65	VAL
1	C	74	LEU
1	C	79	THR
1	C	82	ILE
1	C	95	ILE
1	C	106	ASN
1	C	166	SER
1	C	194	ILE
1	C	218	LEU

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Mol	Chain	Res	Type
1	C	236	THR
1	C	238	VAL
1	C	241	THR
1	C	247	SER
1	D	21	ILE
1	D	35	GLN
1	D	45	ARG
1	D	61	LEU
1	D	65	VAL
1	D	74	LEU
1	D	77	ARG
1	D	101	THR
1	D	105	ILE
1	D	168	LEU
1	D	170	SER
1	D	200	SER
1	D	207	LEU
1	D	233	LYS
1	D	238	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	48	GLN
1	A	70	HIS
1	A	121	HIS
1	A	172	ASN
1	B	35	GLN
1	B	70	HIS
1	C	48	GLN
1	C	187	ASN
1	D	30	GLN
1	D	35	GLN
1	D	172	ASN
1	D	231	ASN

5.3.3 RNA ⓘ

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](#)

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$
2	NAD	B	300	-	46,48,48	3.69	22 (47%)	64,73,73	1.82	17 (26%)
2	NAD	C	300	-	46,48,48	3.73	21 (45%)	64,73,73	1.95	17 (26%)
3	53K	B	301	-	30,30,30	1.61	3 (10%)	41,41,41	1.94	8 (19%)
2	NAD	A	300	-	46,48,48	3.80	21 (45%)	64,73,73	2.00	18 (28%)
3	53K	A	301	-	30,30,30	1.66	3 (10%)	41,41,41	1.40	7 (17%)
3	53K	D	301	-	30,30,30	1.53	3 (10%)	41,41,41	1.47	5 (12%)
2	NAD	D	302	-	46,48,48	3.59	21 (45%)	64,73,73	1.85	14 (21%)
3	53K	C	301	-	30,30,30	1.62	3 (10%)	41,41,41	1.24	4 (9%)

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	300	-	-	14/30/62/62	0/5/5/5
2	NAD	C	300	-	-	7/30/62/62	0/5/5/5
3	53K	B	301	-	-	4/12/12/12	0/4/4/4
2	NAD	A	300	-	-	11/30/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	53K	A	301	-	-	4/12/12/12	0/4/4/4
3	53K	D	301	-	-	2/12/12/12	0/4/4/4
2	NAD	D	302	-	-	8/30/62/62	0/5/5/5
3	53K	C	301	-	-	4/12/12/12	0/4/4/4

All (97) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	300	NAD	C2N-N1N	11.07	1.47	1.35
2	C	300	NAD	C2N-N1N	10.76	1.46	1.35
2	A	300	NAD	C2N-N1N	10.57	1.46	1.35
2	D	302	NAD	C2N-N1N	9.54	1.45	1.35
2	A	300	NAD	PA-O3	8.06	1.68	1.59
2	A	300	NAD	C7N-N7N	7.89	1.47	1.33
2	C	300	NAD	C7N-N7N	7.73	1.47	1.33
2	A	300	NAD	C4N-C3N	7.48	1.50	1.39
2	D	302	NAD	PA-O3	7.46	1.67	1.59
2	D	302	NAD	C7N-N7N	7.36	1.46	1.33
2	B	300	NAD	C7N-N7N	7.29	1.46	1.33
2	A	300	NAD	PN-O3	7.27	1.67	1.59
2	D	302	NAD	PN-O3	7.11	1.67	1.59
2	B	300	NAD	C4N-C3N	7.04	1.50	1.39
2	B	300	NAD	C8A-N7A	7.03	1.45	1.31
2	C	300	NAD	C4N-C3N	7.01	1.50	1.39
2	C	300	NAD	C8A-N7A	6.94	1.45	1.31
2	B	300	NAD	PA-O3	6.77	1.66	1.59
2	C	300	NAD	PA-O3	6.72	1.66	1.59
2	D	302	NAD	C8A-N7A	6.66	1.44	1.31
2	D	302	NAD	C4N-C3N	6.64	1.49	1.39
2	A	300	NAD	C8A-N7A	6.64	1.44	1.31
2	B	300	NAD	C6N-N1N	6.23	1.49	1.35
2	B	300	NAD	PN-O3	6.20	1.66	1.59
2	D	302	NAD	C6N-N1N	6.12	1.49	1.35
2	C	300	NAD	PN-O3	6.08	1.66	1.59
2	A	300	NAD	C6N-N1N	6.05	1.49	1.35
2	C	300	NAD	C6N-N1N	5.92	1.48	1.35
3	B	301	53K	C15-C14	-5.60	1.39	1.47
3	A	301	53K	N2-N3	5.54	1.41	1.32
3	C	301	53K	N2-N3	5.47	1.40	1.32
3	C	301	53K	C15-C14	-5.27	1.39	1.47
2	A	300	NAD	C2N-C3N	5.27	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	53K	N2-N3	5.27	1.40	1.32
3	A	301	53K	C15-C14	-5.27	1.39	1.47
3	D	301	53K	C15-C14	-5.22	1.39	1.47
2	B	300	NAD	C6A-N6A	5.18	1.47	1.34
2	C	300	NAD	C6A-N6A	5.16	1.47	1.34
2	A	300	NAD	C6A-N6A	5.11	1.47	1.34
2	D	302	NAD	C6A-N6A	5.09	1.47	1.34
2	C	300	NAD	C2A-N1A	5.03	1.42	1.33
3	D	301	53K	N2-N3	5.02	1.40	1.32
2	C	300	NAD	C2N-C3N	5.01	1.46	1.39
2	B	300	NAD	C2A-N1A	4.95	1.42	1.33
2	D	302	NAD	C2A-N1A	4.93	1.42	1.33
2	B	300	NAD	PA-O1A	4.83	1.67	1.50
2	D	302	NAD	PA-O1A	4.81	1.67	1.50
2	B	300	NAD	C2N-C3N	4.64	1.46	1.39
2	C	300	NAD	PA-O1A	4.61	1.66	1.50
2	A	300	NAD	PA-O1A	4.48	1.66	1.50
2	B	300	NAD	C2A-N3A	4.47	1.41	1.33
2	C	300	NAD	C2A-N3A	4.45	1.41	1.33
2	A	300	NAD	C2A-N1A	4.44	1.41	1.33
2	D	302	NAD	C2A-N3A	4.42	1.41	1.33
2	C	300	NAD	PN-O1N	4.06	1.64	1.50
2	D	302	NAD	C2N-C3N	4.05	1.45	1.39
2	A	300	NAD	C2A-N3A	4.04	1.41	1.33
2	D	302	NAD	PN-O1N	4.00	1.64	1.50
2	A	300	NAD	PN-O1N	3.98	1.64	1.50
2	A	300	NAD	C5N-C4N	3.89	1.45	1.38
2	C	300	NAD	C3N-C7N	3.89	1.56	1.50
2	B	300	NAD	PN-O1N	3.84	1.64	1.50
3	A	301	53K	N1-N3	3.70	1.40	1.34
2	A	300	NAD	C6N-C5N	3.63	1.46	1.38
2	A	300	NAD	C3N-C7N	3.61	1.56	1.50
2	C	300	NAD	C5N-C4N	3.54	1.45	1.38
3	C	301	53K	N1-N3	3.53	1.40	1.34
2	C	300	NAD	C4A-N9A	3.50	1.45	1.37
2	B	300	NAD	C5N-C4N	3.48	1.44	1.38
2	C	300	NAD	C6N-C5N	3.45	1.45	1.38
2	A	300	NAD	C4A-N9A	3.33	1.44	1.37
2	D	302	NAD	C6N-C5N	3.32	1.45	1.38
2	B	300	NAD	C3B-C4B	3.30	1.61	1.53
2	A	300	NAD	C3B-C4B	3.28	1.61	1.53
2	D	302	NAD	C3B-C4B	3.27	1.61	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	300	NAD	C3B-C4B	3.17	1.61	1.53
2	B	300	NAD	C6N-C5N	3.14	1.45	1.38
2	B	300	NAD	C4A-N9A	3.13	1.44	1.37
2	D	302	NAD	C5N-C4N	3.12	1.44	1.38
2	B	300	NAD	C3N-C7N	2.99	1.55	1.50
2	C	300	NAD	O7N-C7N	-2.99	1.18	1.24
2	D	302	NAD	C4A-N9A	2.89	1.43	1.37
2	B	300	NAD	O7N-C7N	-2.87	1.18	1.24
2	A	300	NAD	O7N-C7N	-2.77	1.19	1.24
2	D	302	NAD	O7N-C7N	-2.60	1.19	1.24
3	D	301	53K	N1-N3	2.55	1.38	1.34
2	D	302	NAD	C3N-C7N	2.53	1.54	1.50
3	B	301	53K	N1-N3	2.53	1.38	1.34
2	D	302	NAD	C3D-C4D	2.34	1.58	1.53
2	C	300	NAD	PA-O5B	2.21	1.68	1.59
2	A	300	NAD	C3D-C4D	2.21	1.58	1.53
2	C	300	NAD	C3D-C4D	2.18	1.58	1.53
2	D	302	NAD	C5A-N7A	2.14	1.43	1.39
2	B	300	NAD	C5A-N7A	2.07	1.42	1.39
2	B	300	NAD	C3D-C4D	2.07	1.58	1.53
2	B	300	NAD	C8A-N9A	2.03	1.41	1.37
2	A	300	NAD	PN-O5D	2.00	1.67	1.59

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	NAD	N3A-C2A-N1A	-7.11	117.83	128.58
3	B	301	53K	C15-C14-C13	-6.82	119.23	129.65
2	C	300	NAD	N3A-C2A-N1A	-6.03	119.45	128.58
2	B	300	NAD	N3A-C2A-N1A	-5.98	119.54	128.58
2	D	302	NAD	N3A-C2A-N1A	-5.79	119.81	128.58
2	C	300	NAD	C4B-O4B-C1B	-5.43	97.47	109.47
2	A	300	NAD	C4B-O4B-C1B	-5.22	97.95	109.47
2	D	302	NAD	N9A-C8A-N7A	-4.93	106.94	113.94
3	D	301	53K	C12-N1-N3	-4.87	113.18	120.03
3	D	301	53K	C12-N1-C13	4.78	135.52	128.65
3	B	301	53K	N1-N3-N2	-4.60	102.89	107.02
2	A	300	NAD	N9A-C8A-N7A	-4.31	107.81	113.94
3	A	301	53K	C15-C14-C13	-4.28	123.11	129.65
2	C	300	NAD	C5A-C4A-N3A	-4.24	120.88	126.72
3	B	301	53K	C12-N1-N3	-4.13	114.22	120.03
2	C	300	NAD	N9A-C8A-N7A	-4.08	108.14	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	NAD	C5A-C4A-N3A	-3.98	121.24	126.72
2	B	300	NAD	N9A-C8A-N7A	-3.95	108.33	113.94
2	D	302	NAD	C4B-O4B-C1B	-3.91	100.84	109.47
3	C	301	53K	C12-N1-C13	3.84	134.16	128.65
2	D	302	NAD	C4A-N9A-C8A	3.78	109.70	105.74
2	B	300	NAD	C5A-C4A-N3A	-3.71	121.61	126.72
2	A	300	NAD	C2A-N3A-C4A	3.53	120.44	111.83
2	D	302	NAD	C5A-C4A-N3A	-3.51	121.89	126.72
2	C	300	NAD	O4B-C1B-N9A	3.44	114.69	108.09
3	C	301	53K	C12-N1-N3	-3.39	115.27	120.03
3	B	301	53K	C12-N1-C13	3.38	133.51	128.65
2	C	300	NAD	C3N-C7N-N7N	3.19	121.67	117.74
2	B	300	NAD	C4B-O4B-C1B	-3.16	102.49	109.47
3	B	301	53K	C15-C14-N2	-3.15	116.98	122.05
2	D	302	NAD	O3D-C3D-C4D	3.15	120.13	111.08
2	C	300	NAD	C2A-N3A-C4A	3.12	119.45	111.83
2	B	300	NAD	C6N-N1N-C1D	-3.04	113.75	119.73
2	B	300	NAD	C2A-N3A-C4A	2.96	119.07	111.83
2	C	300	NAD	O7N-C7N-N7N	-2.93	118.39	122.62
3	A	301	53K	N1-N3-N2	-2.85	104.46	107.02
2	A	300	NAD	C3N-C7N-N7N	2.83	121.23	117.74
3	A	301	53K	C12-N1-N3	-2.83	116.05	120.03
2	D	302	NAD	C2A-N3A-C4A	2.81	118.69	111.83
2	B	300	NAD	O4D-C4D-C3D	-2.81	99.58	105.15
3	A	301	53K	C12-N1-C13	2.76	132.61	128.65
2	B	300	NAD	O7N-C7N-N7N	-2.74	118.66	122.62
2	C	300	NAD	C5A-N7A-C8A	2.74	107.75	103.45
2	B	300	NAD	O2B-C2B-C1B	2.69	119.38	110.10
3	A	301	53K	C8-O1-C7	2.69	124.51	118.09
2	D	302	NAD	C5A-N7A-C8A	2.68	107.67	103.45
2	A	300	NAD	O3D-C3D-C4D	2.68	118.78	111.08
3	C	301	53K	C15-C14-C13	-2.65	125.60	129.65
3	D	301	53K	N1-N3-N2	-2.65	104.64	107.02
3	D	301	53K	C15-C14-C13	-2.64	125.61	129.65
2	C	300	NAD	O4B-C4B-C3B	-2.60	99.99	105.15
2	D	302	NAD	O2B-C2B-C1B	2.59	119.04	110.10
3	B	301	53K	C14-C13-N1	2.59	107.37	105.28
2	A	300	NAD	C5A-N7A-C8A	2.59	107.52	103.45
2	C	300	NAD	O3D-C3D-C4D	2.59	118.51	111.08
2	C	300	NAD	O4B-C4B-C5B	2.58	117.60	109.33
3	B	301	53K	C10-C11-C21	2.53	122.03	118.55
2	D	302	NAD	C2N-C3N-C4N	2.46	121.12	118.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300	NAD	O5B-C5B-C4B	2.44	117.32	108.99
2	A	300	NAD	C4A-N9A-C8A	2.44	108.30	105.74
2	B	300	NAD	C3N-C7N-N7N	2.41	120.71	117.74
2	D	302	NAD	C3B-C2B-C1B	-2.40	96.92	101.46
2	D	302	NAD	O4B-C1B-C2B	-2.39	101.50	106.62
3	B	301	53K	C22-C21-C11	-2.37	118.36	120.81
2	A	300	NAD	O5D-C5D-C4D	2.36	117.03	108.99
2	A	300	NAD	O2B-C2B-C1B	2.36	118.22	110.10
2	C	300	NAD	C4A-C5A-N7A	-2.35	107.89	110.58
2	A	300	NAD	O2N-PN-O3	2.32	113.56	107.27
2	C	300	NAD	C4A-N9A-C8A	2.31	108.17	105.74
2	C	300	NAD	N3A-C4A-N9A	2.30	131.08	127.17
2	D	302	NAD	N3A-C4A-N9A	2.28	131.04	127.17
2	A	300	NAD	N3A-C4A-N9A	2.27	131.03	127.17
2	B	300	NAD	C5N-C4N-C3N	2.25	122.58	120.36
2	B	300	NAD	O2N-PN-O3	2.24	113.34	107.27
2	B	300	NAD	C2N-N1N-C1D	2.23	124.05	119.13
2	A	300	NAD	O4B-C1B-C2B	-2.22	101.87	106.62
2	B	300	NAD	C4A-N9A-C8A	2.21	108.06	105.74
3	C	301	53K	C8-O1-C7	2.15	123.22	118.09
3	A	301	53K	C1-C2-C7	-2.12	117.94	120.94
2	A	300	NAD	O7N-C7N-N7N	-2.12	119.55	122.62
2	A	300	NAD	O4B-C4B-C3B	-2.12	100.95	105.15
2	A	300	NAD	O3D-C3D-C2D	2.09	118.52	111.82
2	A	300	NAD	C2A-N1A-C6A	2.07	122.13	118.73
2	C	300	NAD	C5A-C4A-N9A	2.05	108.05	105.81
2	D	302	NAD	O2N-PN-O3	2.05	112.81	107.27
3	A	301	53K	C14-C13-N1	2.05	106.93	105.28
2	B	300	NAD	C5A-N7A-C8A	2.05	106.67	103.45
3	D	301	53K	C14-C13-N1	2.02	106.91	105.28
2	B	300	NAD	O4B-C4B-C3B	-2.02	101.15	105.15
2	B	300	NAD	N3A-C4A-N9A	2.01	130.58	127.17

There are no chirality outliers.

All (54) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	300	NAD	O4D-C1D-N1N-C2N
2	A	300	NAD	O4D-C1D-N1N-C6N
2	B	300	NAD	C5D-O5D-PN-O1N
2	B	300	NAD	O4D-C1D-N1N-C2N
2	B	300	NAD	O4D-C1D-N1N-C6N

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Mol	Chain	Res	Type	Atoms
2	B	300	NAD	C2D-C1D-N1N-C2N
2	B	300	NAD	C2D-C1D-N1N-C6N
2	C	300	NAD	O4B-C4B-C5B-O5B
2	C	300	NAD	O4D-C4D-C5D-O5D
2	C	300	NAD	O4D-C1D-N1N-C2N
2	C	300	NAD	O4D-C1D-N1N-C6N
2	D	302	NAD	C5D-O5D-PN-O3
2	D	302	NAD	C5D-O5D-PN-O1N
2	D	302	NAD	C5D-O5D-PN-O2N
2	D	302	NAD	O4D-C1D-N1N-C2N
3	A	301	53K	N2-C14-C15-C16
3	B	301	53K	N2-C14-C15-C20
3	C	301	53K	C13-C14-C15-C20
3	C	301	53K	C13-C14-C15-C16
3	C	301	53K	N2-C14-C15-C16
3	A	301	53K	N2-C14-C15-C20
3	B	301	53K	N2-C14-C15-C16
3	C	301	53K	N2-C14-C15-C20
3	A	301	53K	C13-C14-C15-C20
3	A	301	53K	C13-C14-C15-C16
2	A	300	NAD	O4D-C4D-C5D-O5D
2	C	300	NAD	C3B-C4B-C5B-O5B
2	B	300	NAD	C2N-C3N-C7N-N7N
2	B	300	NAD	C4N-C3N-C7N-N7N
2	B	300	NAD	O4D-C4D-C5D-O5D
2	B	300	NAD	C3D-C4D-C5D-O5D
3	B	301	53K	C13-C14-C15-C16
2	C	300	NAD	C3D-C4D-C5D-O5D
3	B	301	53K	C13-C14-C15-C20
2	A	300	NAD	C3D-C4D-C5D-O5D
3	D	301	53K	N2-C14-C15-C20
2	A	300	NAD	PN-O3-PA-O5B
2	B	300	NAD	PN-O3-PA-O5B
2	D	302	NAD	PN-O3-PA-O5B
3	D	301	53K	N2-C14-C15-C16
2	A	300	NAD	O4B-C4B-C5B-O5B
2	A	300	NAD	PA-O3-PN-O1N
2	D	302	NAD	PA-O3-PN-O1N
2	A	300	NAD	C5B-O5B-PA-O1A
2	A	300	NAD	C5B-O5B-PA-O3
2	B	300	NAD	C5D-O5D-PN-O3
2	C	300	NAD	C5B-O5B-PA-O3

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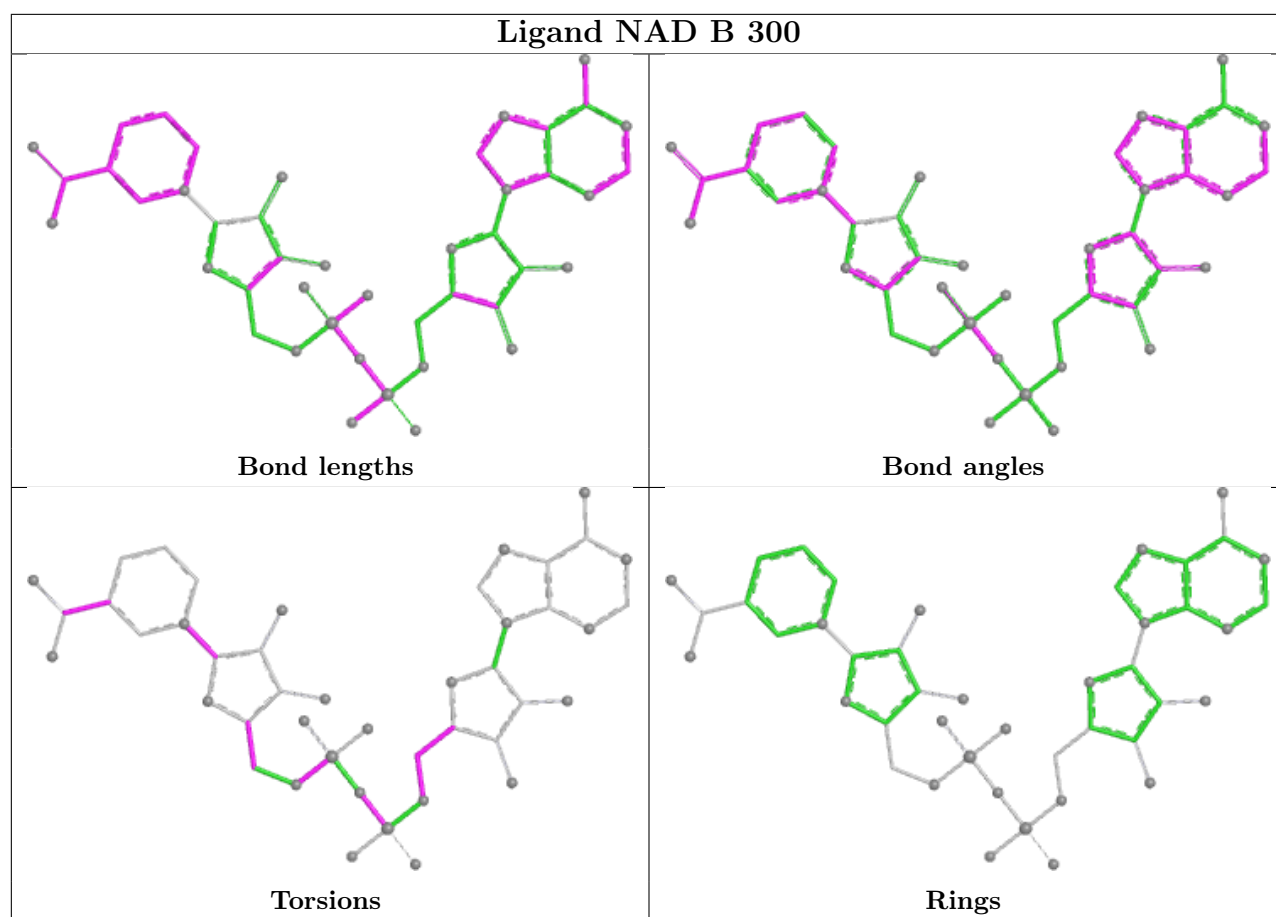
Mol	Chain	Res	Type	Atoms
2	A	300	NAD	C2D-C1D-N1N-C6N
2	B	300	NAD	O4B-C4B-C5B-O5B
2	D	302	NAD	O4D-C1D-N1N-C6N
2	B	300	NAD	C4B-C5B-O5B-PA
2	B	300	NAD	C4N-C3N-C7N-O7N
2	A	300	NAD	PA-O3-PN-O2N
2	D	302	NAD	PA-O3-PN-O2N

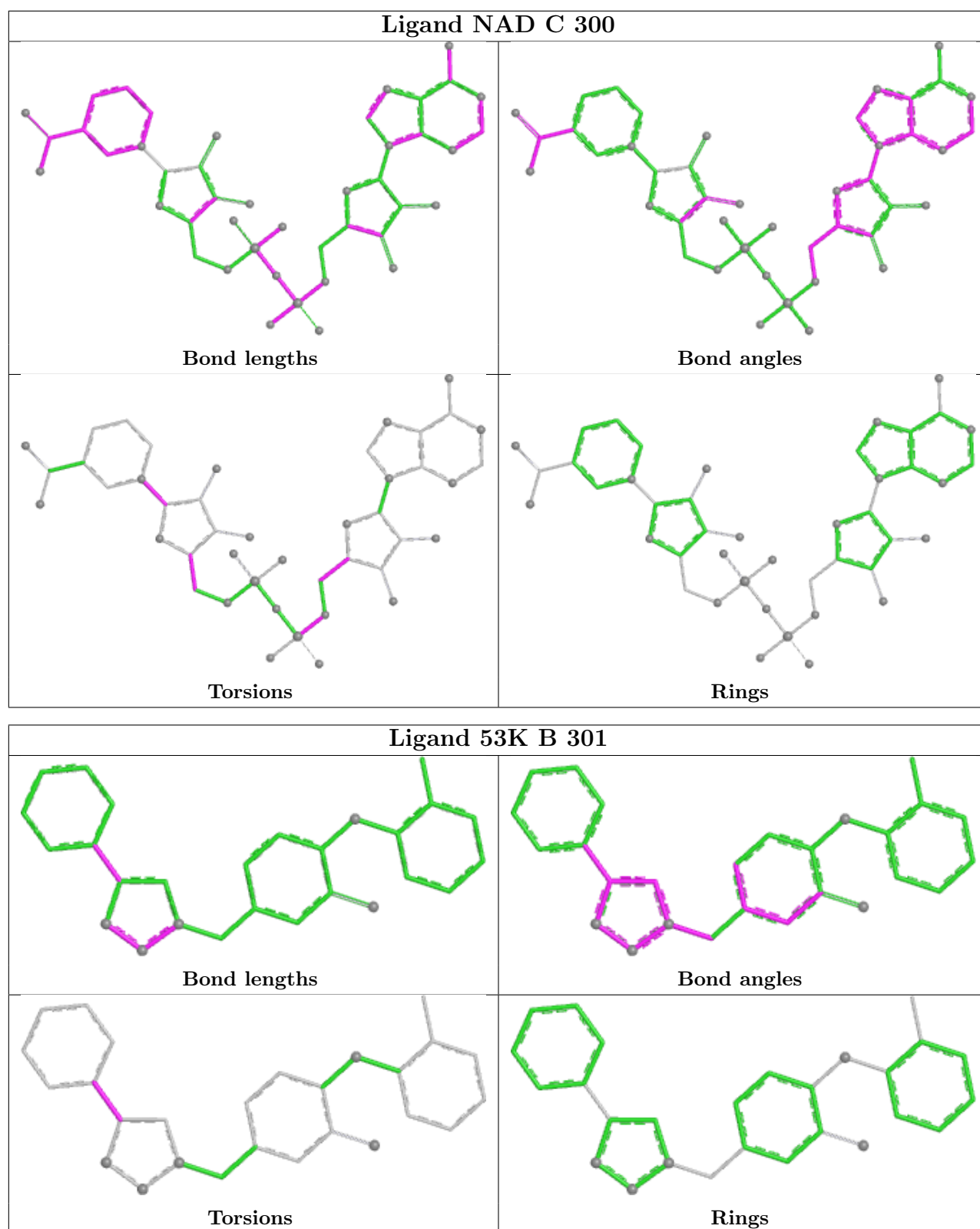
There are no ring outliers.

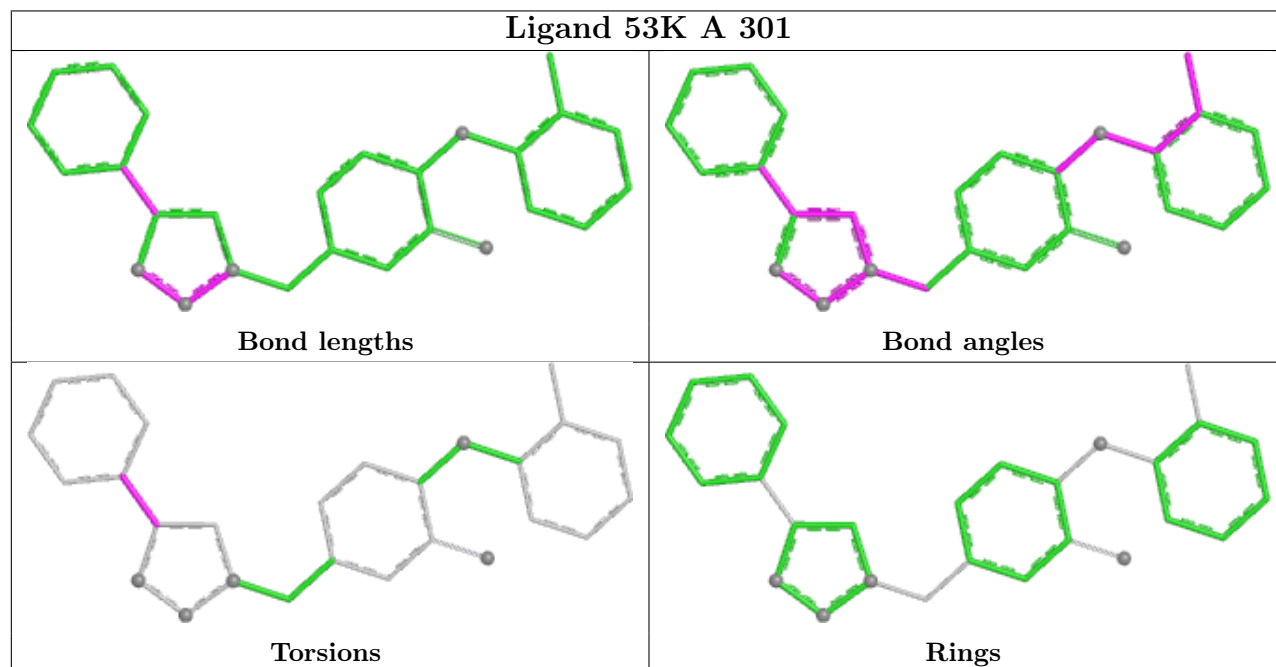
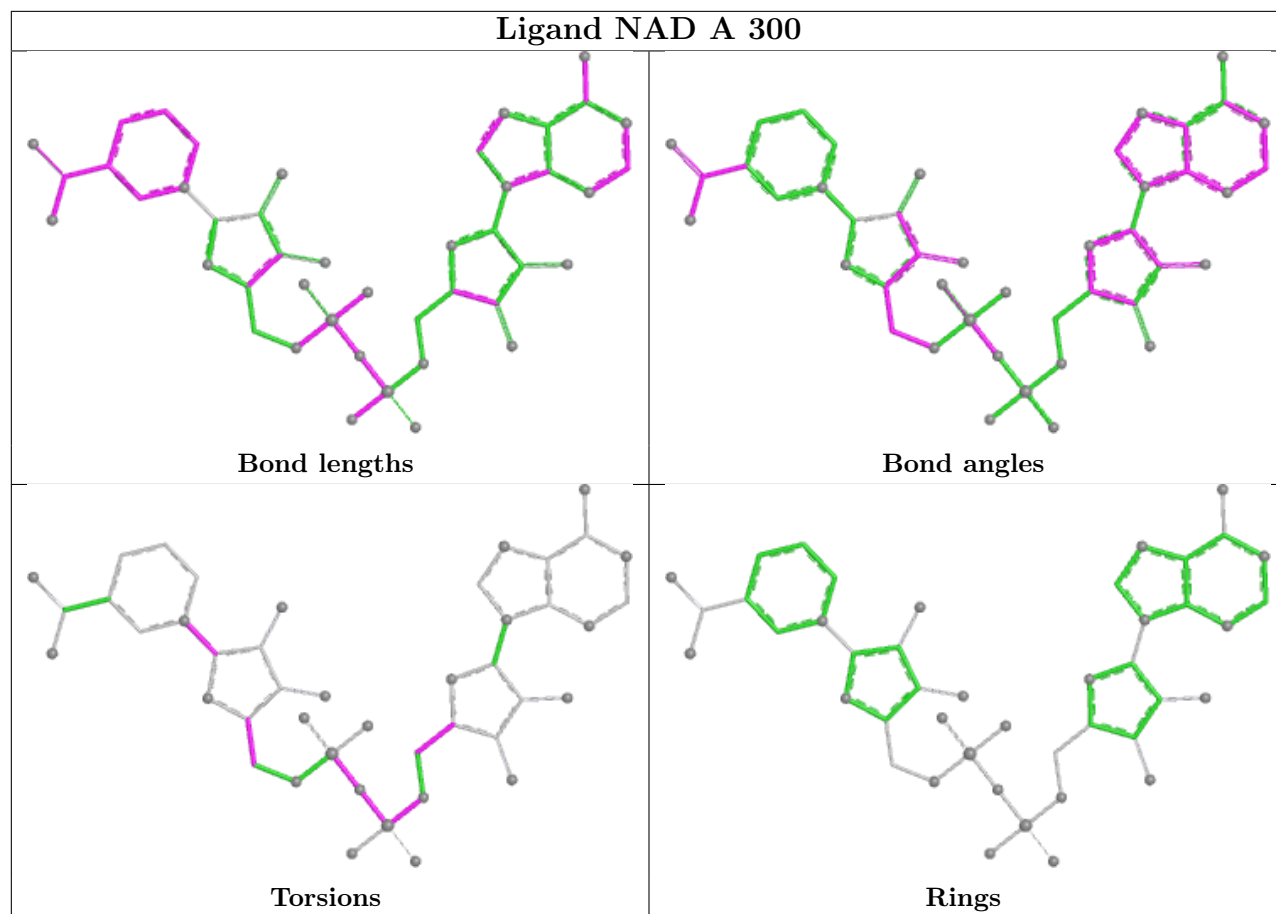
7 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	300	NAD	2	0
3	B	301	53K	2	0
2	A	300	NAD	3	0
3	A	301	53K	4	0
3	D	301	53K	2	0
2	D	302	NAD	2	0
3	C	301	53K	4	0

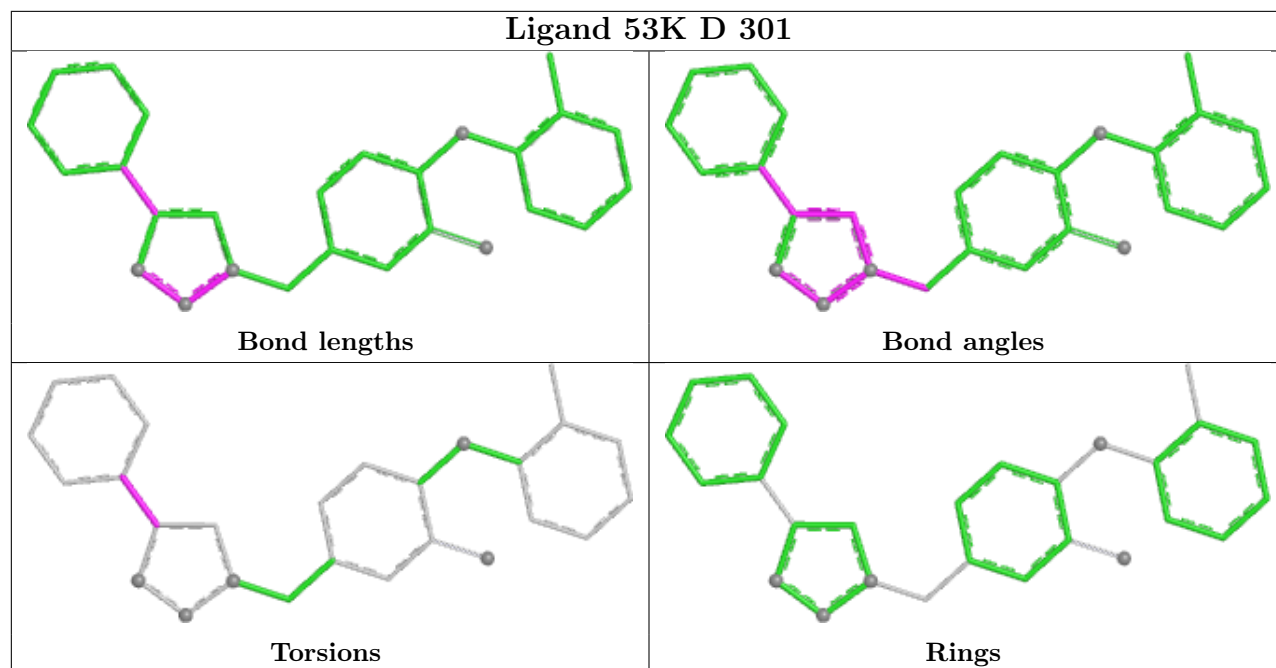
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



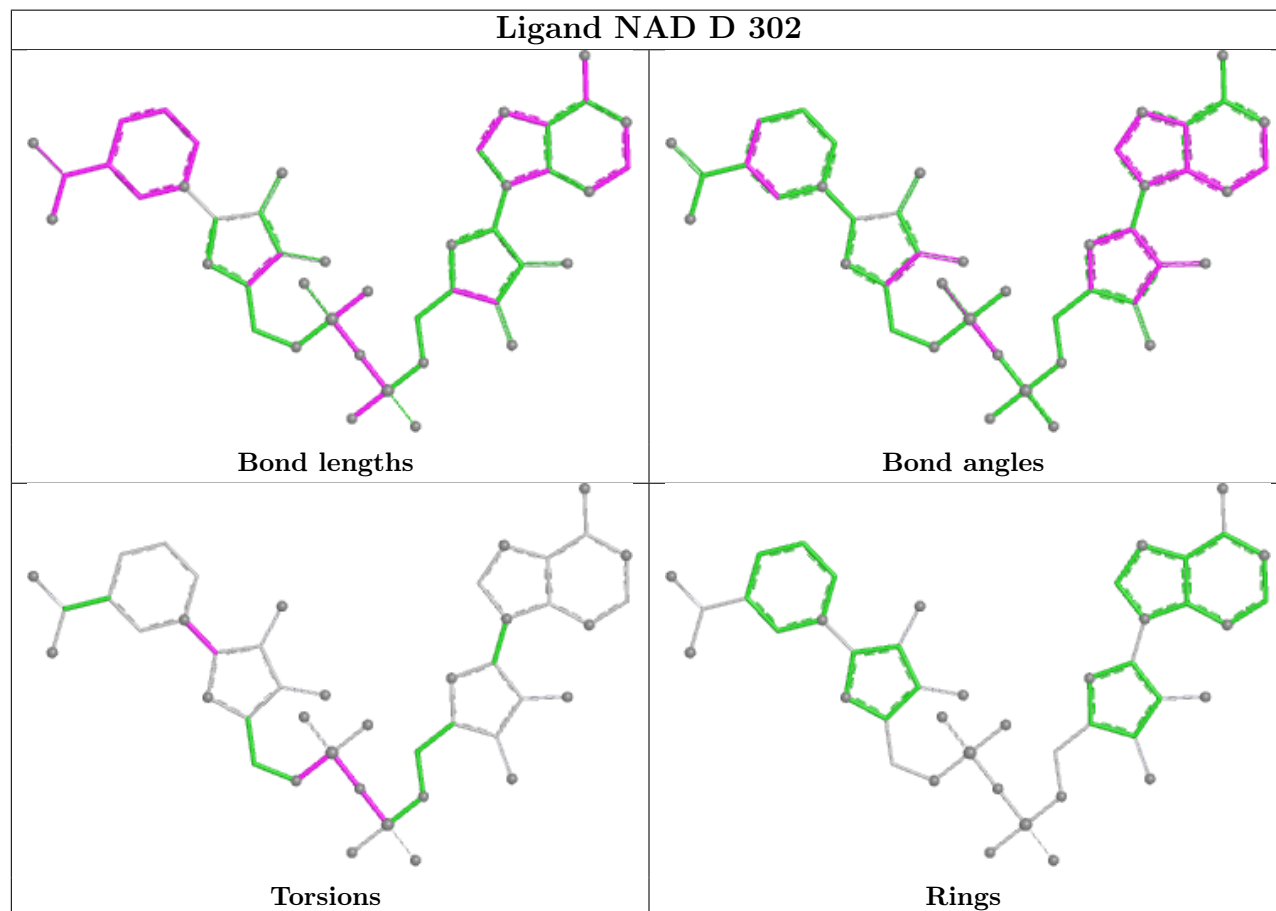


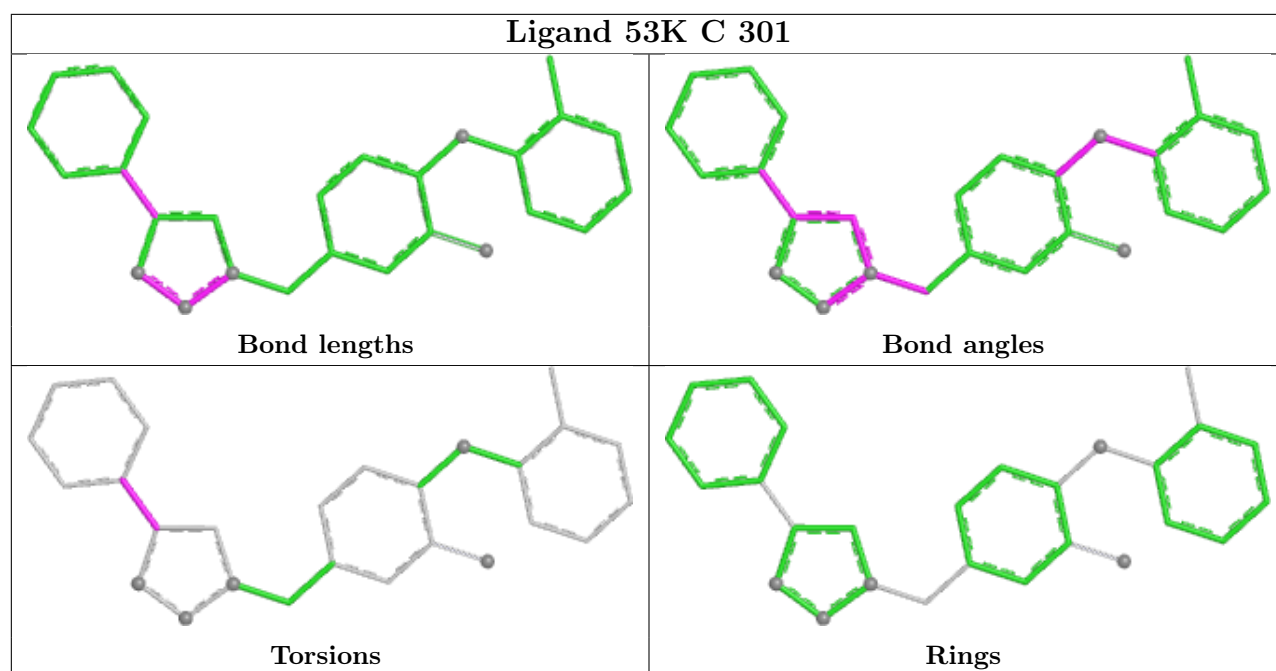


Ligand 53K D 301



Ligand NAD D 302





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	267/289 (92%)	-0.36	0 100 100	31, 59, 84, 98	1 (0%)
1	B	267/289 (92%)	-0.44	0 100 100	37, 55, 75, 91	1 (0%)
1	C	267/289 (92%)	-0.39	0 100 100	30, 58, 87, 98	1 (0%)
1	D	267/289 (92%)	-0.40	0 100 100	37, 58, 82, 96	0
All	All	1068/1156 (92%)	-0.39	0 100 100	30, 57, 83, 98	3 (0%)

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

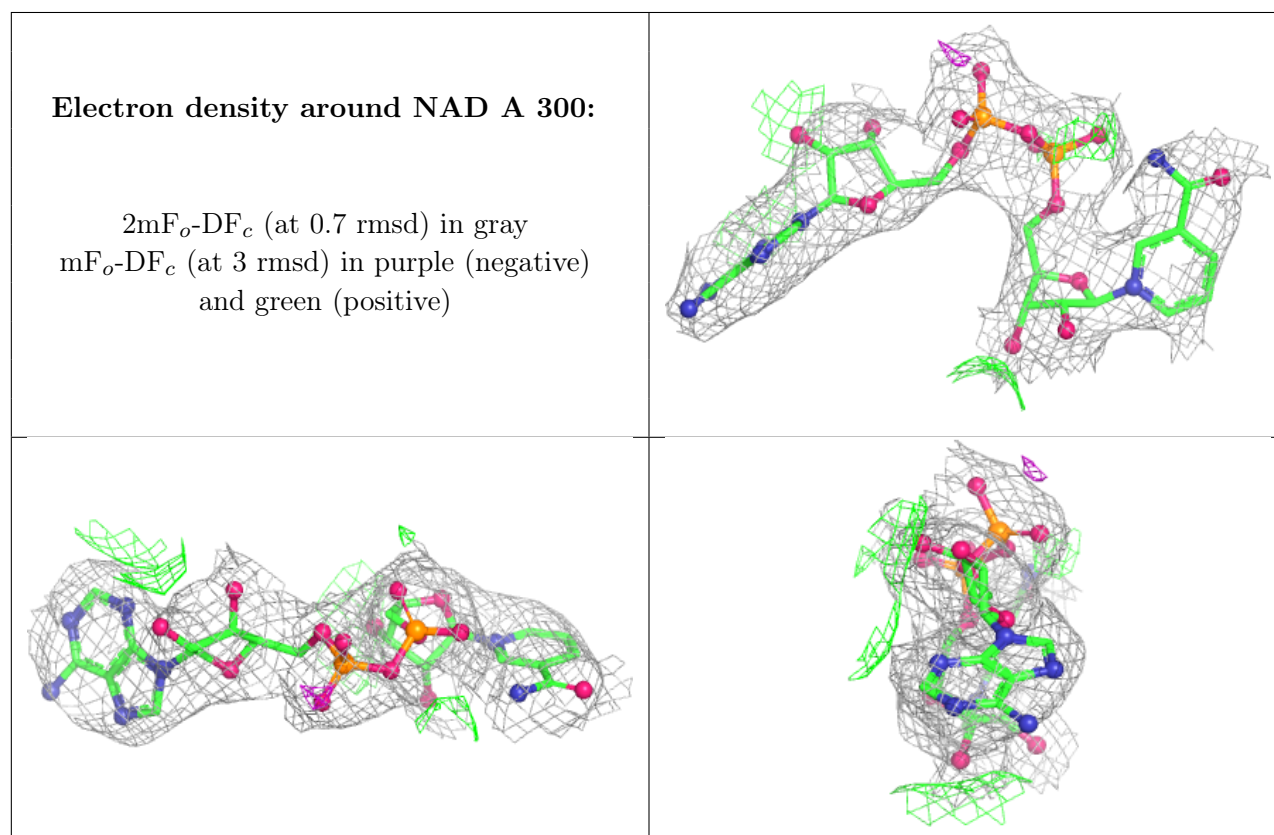
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAD	A	300	44/44	0.94	0.08	52,61,69,75	0
2	NAD	B	300	44/44	0.95	0.08	45,51,57,61	0
2	NAD	C	300	44/44	0.96	0.07	51,59,65,70	0
2	NAD	D	302	44/44	0.96	0.07	46,55,65,72	0

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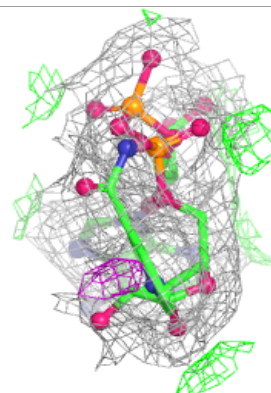
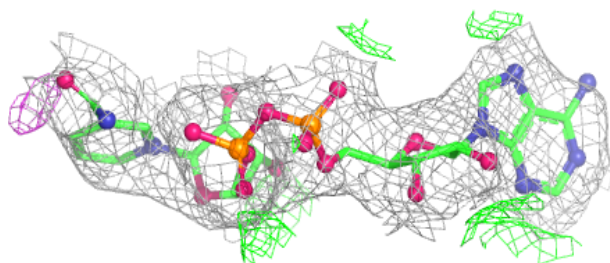
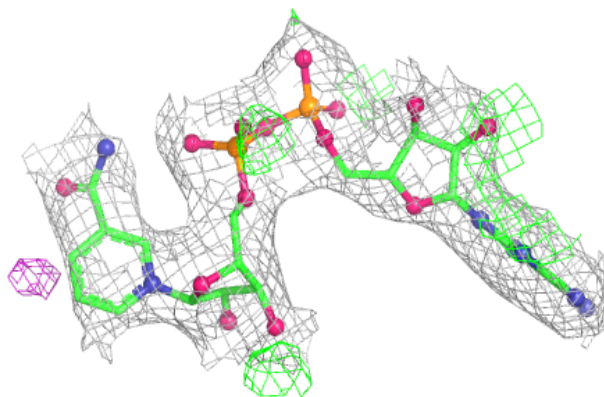
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	53K	A	301	27/27	0.96	0.10	64,74,82,84	0
3	53K	C	301	27/27	0.96	0.10	63,70,80,84	0
3	53K	B	301	27/27	0.97	0.09	45,54,59,62	0
3	53K	D	301	27/27	0.97	0.08	51,58,63,67	0

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.

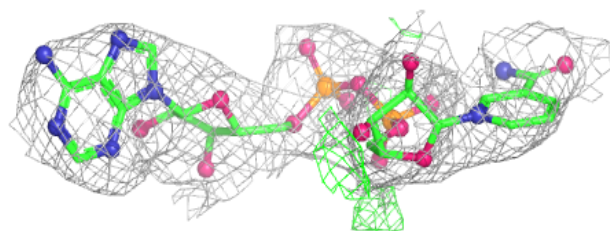
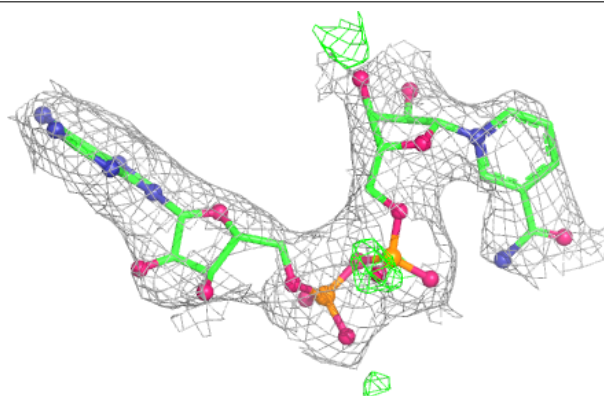


Electron density around NAD B 300:

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

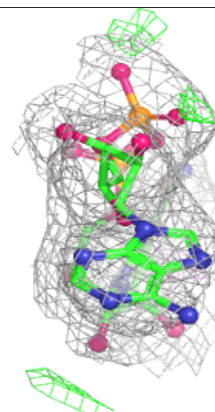
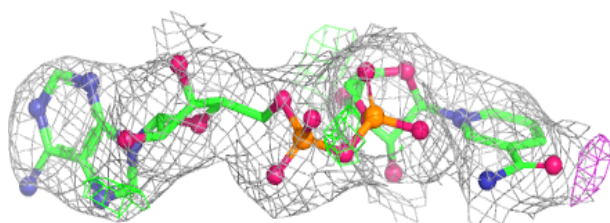
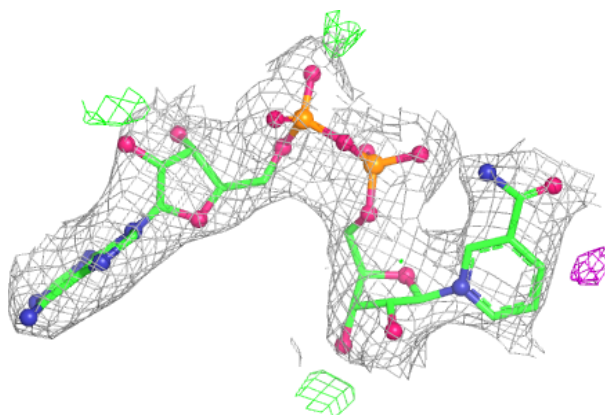
**Electron density around NAD C 300:**

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

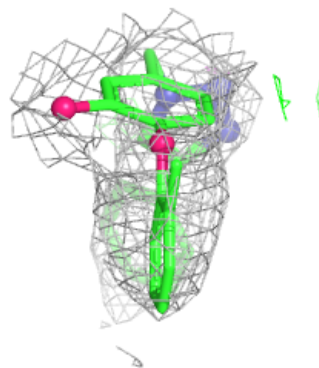
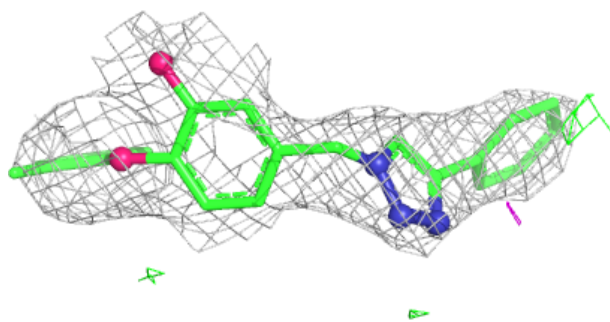
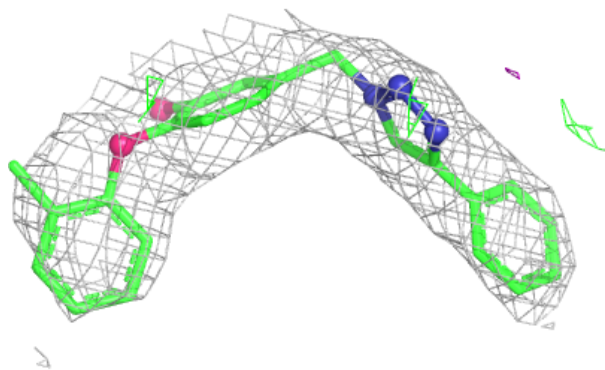


Electron density around NAD D 302:

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

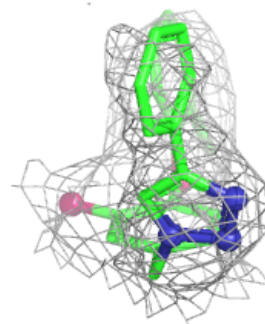
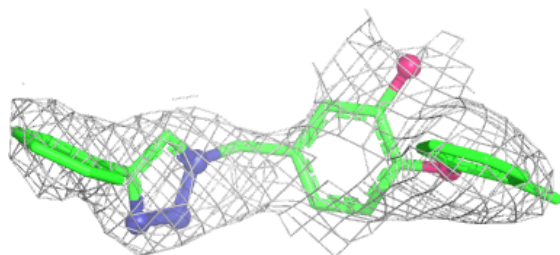
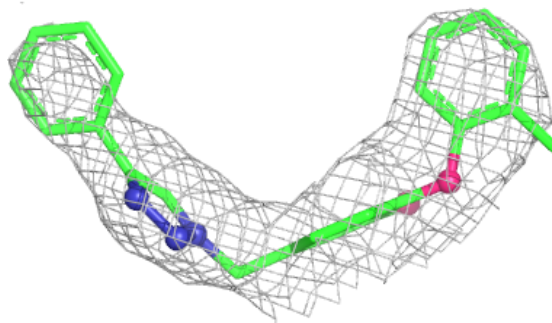
**Electron density around 53K A 301:**

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

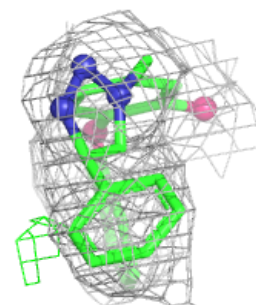
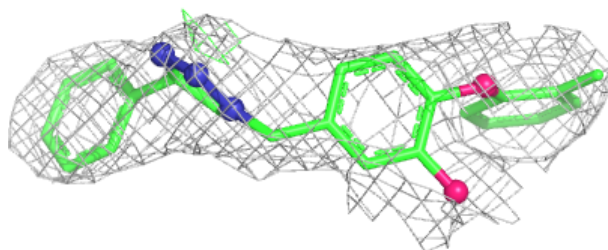
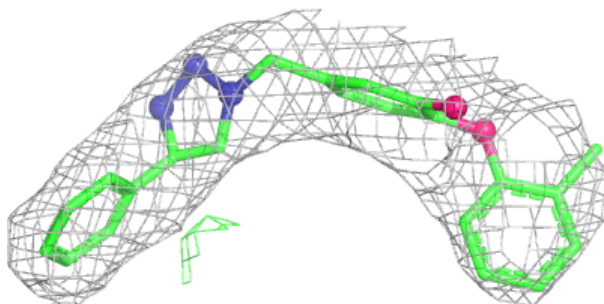


Electron density around 53K C 301:

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

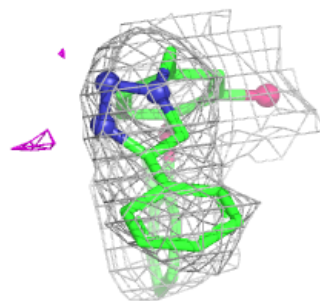
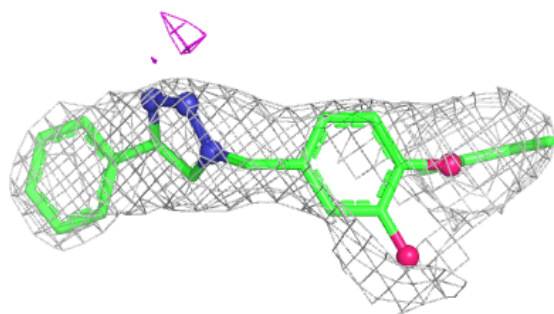
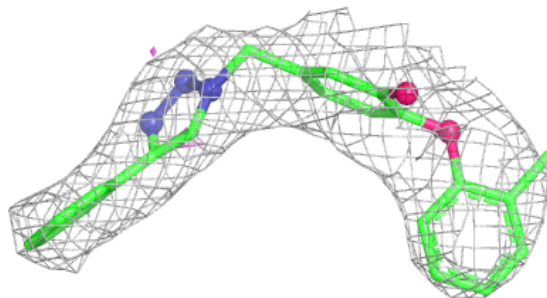
**Electron density around 53K B 301:**

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



Electron density around 53K D 301:

$2mF_o - DF_c$ (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.