



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:43 PM UTC

PDB ID : 1ZBQ / pdb_00001zbq
Title : Crystal Structure Of Human 17-Beta-Hydroxysteroid Dehydrogenase Type 4
In Complex With NAD
Authors : Lukacik, P.; Shafqat, N.; Kavanagh, K.; Bray, J.; von Delft, F.; Edwards,
A.; Arrowsmith, C.; Sundstrom, M.; Oppermann, U.; Structural Genomics
Consortium (SGC)
Deposited on : 2005-04-08
Resolution : 2.19 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

The following versions of software and data (see [references](#) ①) were used in the production of this report:

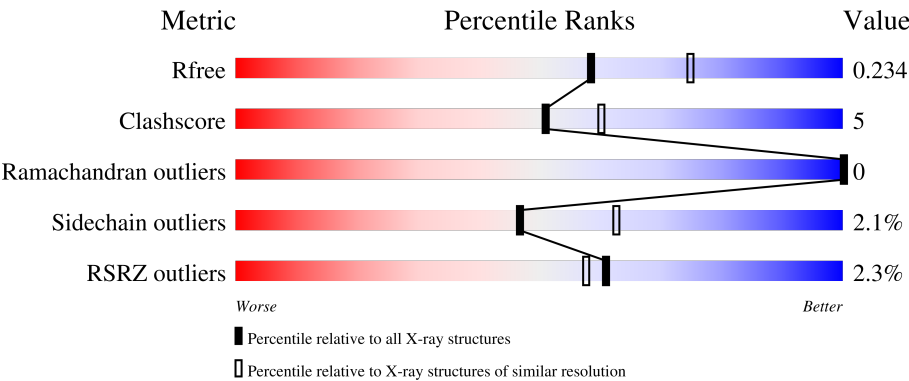
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	327	82%10%8%
1	B	327	85%8%8%
1	C	327	79%12%8%
1	D	327	2%81%11%8%
1	E	327	8%85%7%8%

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Mol	Chain	Length	Quality of chain
1	F	327	2% 84% 8% 8%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 14189 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 17-beta-hydroxysteroid dehydrogenase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	2	0
			2267	1430	409	419	9			
1	B	302	Total	C	N	O	S	0	1	0
			2264	1427	410	418	9			
1	C	302	Total	C	N	O	S	0	1	0
			2268	1428	409	422	9			
1	D	302	Total	C	N	O	S	0	3	0
			2289	1442	419	419	9			
1	E	302	Total	C	N	O	S	0	1	0
			2235	1403	408	415	9			
1	F	302	Total	C	N	O	S	0	2	0
			2259	1418	407	425	9			

There are 138 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	GLY	-	cloning artifact	UNP P51659
A	-19	SER	-	cloning artifact	UNP P51659
A	-18	SER	-	cloning artifact	UNP P51659
A	-17	HIS	-	expression tag	UNP P51659
A	-16	HIS	-	expression tag	UNP P51659
A	-15	HIS	-	expression tag	UNP P51659
A	-14	HIS	-	expression tag	UNP P51659
A	-13	HIS	-	expression tag	UNP P51659
A	-12	HIS	-	expression tag	UNP P51659
A	-11	SER	-	cloning artifact	UNP P51659
A	-10	SER	-	cloning artifact	UNP P51659
A	-9	GLY	-	cloning artifact	UNP P51659
A	-8	ARG	-	cloning artifact	UNP P51659
A	-7	GLU	-	cloning artifact	UNP P51659
A	-6	ASN	-	cloning artifact	UNP P51659
A	-5	LEU	-	cloning artifact	UNP P51659
A	-4	TYR	-	cloning artifact	UNP P51659

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	PHE	-	cloning artifact	UNP P51659
A	-2	GLN	-	cloning artifact	UNP P51659
A	-1	GLY	-	cloning artifact	UNP P51659
A	0	HIS	-	cloning artifact	UNP P51659
A	305	GLY	-	cloning artifact	UNP P51659
A	306	SER	-	cloning artifact	UNP P51659
B	-20	GLY	-	cloning artifact	UNP P51659
B	-19	SER	-	cloning artifact	UNP P51659
B	-18	SER	-	cloning artifact	UNP P51659
B	-17	HIS	-	expression tag	UNP P51659
B	-16	HIS	-	expression tag	UNP P51659
B	-15	HIS	-	expression tag	UNP P51659
B	-14	HIS	-	expression tag	UNP P51659
B	-13	HIS	-	expression tag	UNP P51659
B	-12	HIS	-	expression tag	UNP P51659
B	-11	SER	-	cloning artifact	UNP P51659
B	-10	SER	-	cloning artifact	UNP P51659
B	-9	GLY	-	cloning artifact	UNP P51659
B	-8	ARG	-	cloning artifact	UNP P51659
B	-7	GLU	-	cloning artifact	UNP P51659
B	-6	ASN	-	cloning artifact	UNP P51659
B	-5	LEU	-	cloning artifact	UNP P51659
B	-4	TYR	-	cloning artifact	UNP P51659
B	-3	PHE	-	cloning artifact	UNP P51659
B	-2	GLN	-	cloning artifact	UNP P51659
B	-1	GLY	-	cloning artifact	UNP P51659
B	0	HIS	-	cloning artifact	UNP P51659
B	305	GLY	-	cloning artifact	UNP P51659
B	306	SER	-	cloning artifact	UNP P51659
C	-20	GLY	-	cloning artifact	UNP P51659
C	-19	SER	-	cloning artifact	UNP P51659
C	-18	SER	-	cloning artifact	UNP P51659
C	-17	HIS	-	expression tag	UNP P51659
C	-16	HIS	-	expression tag	UNP P51659
C	-15	HIS	-	expression tag	UNP P51659
C	-14	HIS	-	expression tag	UNP P51659
C	-13	HIS	-	expression tag	UNP P51659
C	-12	HIS	-	expression tag	UNP P51659
C	-11	SER	-	cloning artifact	UNP P51659
C	-10	SER	-	cloning artifact	UNP P51659
C	-9	GLY	-	cloning artifact	UNP P51659
C	-8	ARG	-	cloning artifact	UNP P51659

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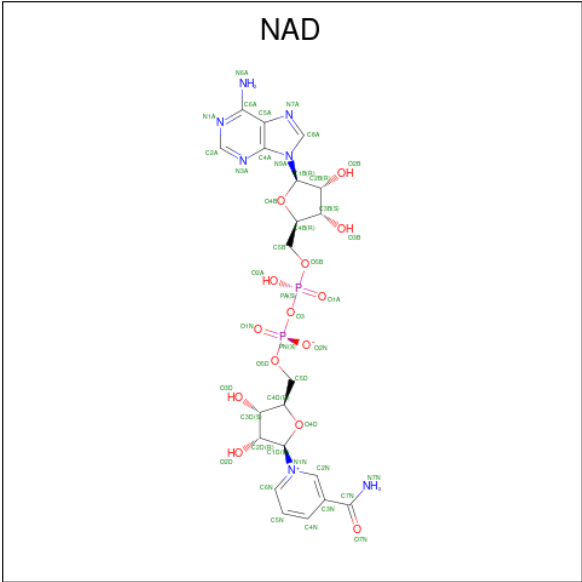
Chain	Residue	Modelled	Actual	Comment	Reference
C	-7	GLU	-	cloning artifact	UNP P51659
C	-6	ASN	-	cloning artifact	UNP P51659
C	-5	LEU	-	cloning artifact	UNP P51659
C	-4	TYR	-	cloning artifact	UNP P51659
C	-3	PHE	-	cloning artifact	UNP P51659
C	-2	GLN	-	cloning artifact	UNP P51659
C	-1	GLY	-	cloning artifact	UNP P51659
C	0	HIS	-	cloning artifact	UNP P51659
C	305	GLY	-	cloning artifact	UNP P51659
C	306	SER	-	cloning artifact	UNP P51659
D	-20	GLY	-	cloning artifact	UNP P51659
D	-19	SER	-	cloning artifact	UNP P51659
D	-18	SER	-	cloning artifact	UNP P51659
D	-17	HIS	-	expression tag	UNP P51659
D	-16	HIS	-	expression tag	UNP P51659
D	-15	HIS	-	expression tag	UNP P51659
D	-14	HIS	-	expression tag	UNP P51659
D	-13	HIS	-	expression tag	UNP P51659
D	-12	HIS	-	expression tag	UNP P51659
D	-11	SER	-	cloning artifact	UNP P51659
D	-10	SER	-	cloning artifact	UNP P51659
D	-9	GLY	-	cloning artifact	UNP P51659
D	-8	ARG	-	cloning artifact	UNP P51659
D	-7	GLU	-	cloning artifact	UNP P51659
D	-6	ASN	-	cloning artifact	UNP P51659
D	-5	LEU	-	cloning artifact	UNP P51659
D	-4	TYR	-	cloning artifact	UNP P51659
D	-3	PHE	-	cloning artifact	UNP P51659
D	-2	GLN	-	cloning artifact	UNP P51659
D	-1	GLY	-	cloning artifact	UNP P51659
D	0	HIS	-	cloning artifact	UNP P51659
D	305	GLY	-	cloning artifact	UNP P51659
D	306	SER	-	cloning artifact	UNP P51659
E	-20	GLY	-	cloning artifact	UNP P51659
E	-19	SER	-	cloning artifact	UNP P51659
E	-18	SER	-	cloning artifact	UNP P51659
E	-17	HIS	-	expression tag	UNP P51659
E	-16	HIS	-	expression tag	UNP P51659
E	-15	HIS	-	expression tag	UNP P51659
E	-14	HIS	-	expression tag	UNP P51659
E	-13	HIS	-	expression tag	UNP P51659
E	-12	HIS	-	expression tag	UNP P51659

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-11	SER	-	cloning artifact	UNP P51659
E	-10	SER	-	cloning artifact	UNP P51659
E	-9	GLY	-	cloning artifact	UNP P51659
E	-8	ARG	-	cloning artifact	UNP P51659
E	-7	GLU	-	cloning artifact	UNP P51659
E	-6	ASN	-	cloning artifact	UNP P51659
E	-5	LEU	-	cloning artifact	UNP P51659
E	-4	TYR	-	cloning artifact	UNP P51659
E	-3	PHE	-	cloning artifact	UNP P51659
E	-2	GLN	-	cloning artifact	UNP P51659
E	-1	GLY	-	cloning artifact	UNP P51659
E	0	HIS	-	cloning artifact	UNP P51659
E	305	GLY	-	cloning artifact	UNP P51659
E	306	SER	-	cloning artifact	UNP P51659
F	-20	GLY	-	cloning artifact	UNP P51659
F	-19	SER	-	cloning artifact	UNP P51659
F	-18	SER	-	cloning artifact	UNP P51659
F	-17	HIS	-	expression tag	UNP P51659
F	-16	HIS	-	expression tag	UNP P51659
F	-15	HIS	-	expression tag	UNP P51659
F	-14	HIS	-	expression tag	UNP P51659
F	-13	HIS	-	expression tag	UNP P51659
F	-12	HIS	-	expression tag	UNP P51659
F	-11	SER	-	cloning artifact	UNP P51659
F	-10	SER	-	cloning artifact	UNP P51659
F	-9	GLY	-	cloning artifact	UNP P51659
F	-8	ARG	-	cloning artifact	UNP P51659
F	-7	GLU	-	cloning artifact	UNP P51659
F	-6	ASN	-	cloning artifact	UNP P51659
F	-5	LEU	-	cloning artifact	UNP P51659
F	-4	TYR	-	cloning artifact	UNP P51659
F	-3	PHE	-	cloning artifact	UNP P51659
F	-2	GLN	-	cloning artifact	UNP P51659
F	-1	GLY	-	cloning artifact	UNP P51659
F	0	HIS	-	cloning artifact	UNP P51659
F	305	GLY	-	cloning artifact	UNP P51659
F	306	SER	-	cloning artifact	UNP P51659

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



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

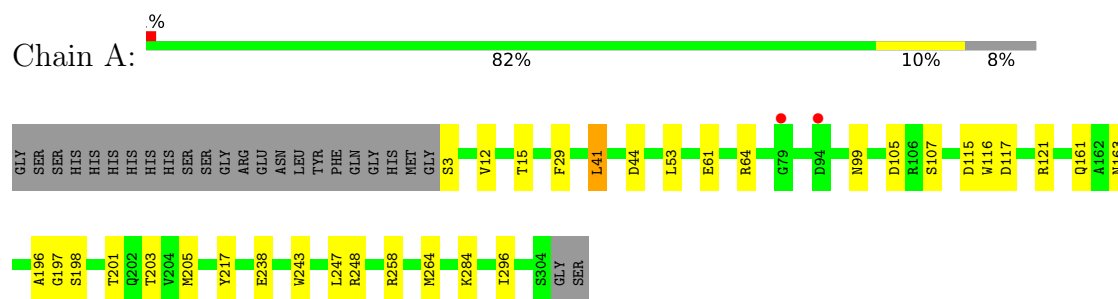
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	61	Total	O	0	0
			61	61		
3	B	67	Total	O	0	0
			67	67		
3	C	66	Total	O	0	0
			66	66		
3	D	64	Total	O	0	0
			64	64		
3	E	42	Total	O	0	0
			42	42		
3	F	43	Total	O	0	0
			43	43		

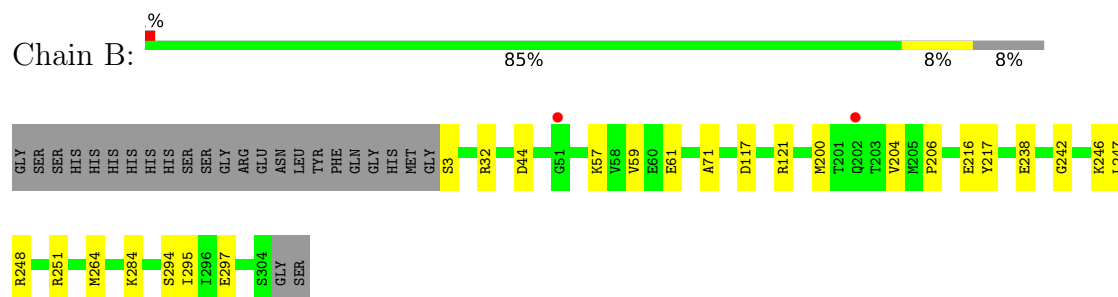
3 Residue-property plots [i](#)

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.

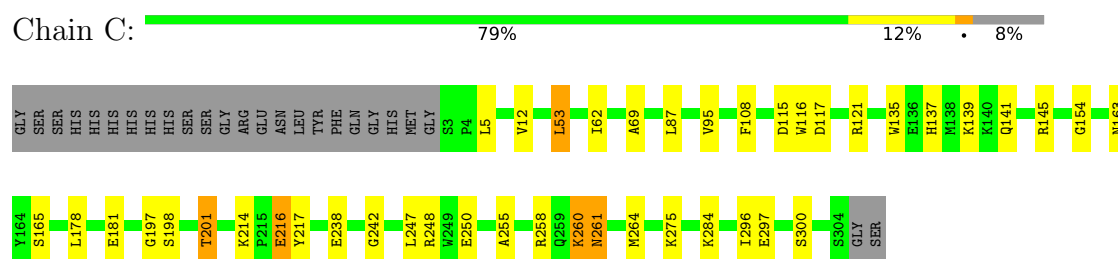
• Molecule 1: 17-beta-hydroxysteroid dehydrogenase 4



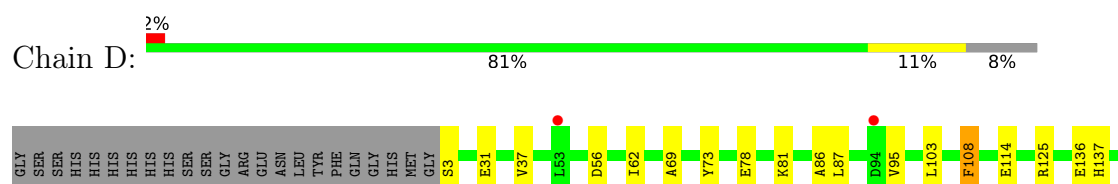
• Molecule 1: 17-beta-hydroxysteroid dehydrogenase 4



• Molecule 1: 17-beta-hydroxysteroid dehydrogenase 4

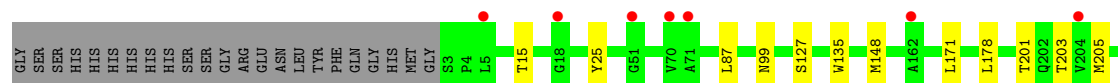
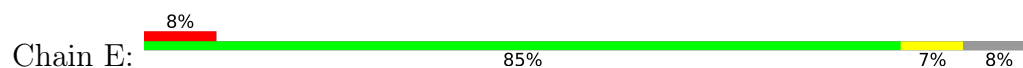


• Molecule 1: 17-beta-hydroxysteroid dehydrogenase 4

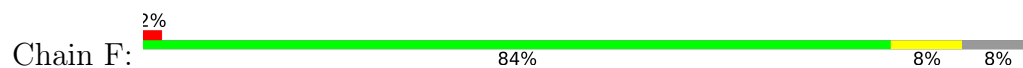




- Molecule 1: 17-beta-hydroxysteroid dehydrogenase 4



- Molecule 1: 17-beta-hydroxysteroid dehydrogenase 4



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.83Å 50.77Å 183.78Å 87.52° 87.07° 70.75°	Depositor
Resolution (Å)	50.57 – 2.19 50.57 – 2.19	Depositor EDS
% Data completeness (in resolution range)	76.2 (50.57-2.19) 76.2 (50.57-2.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.188 , 0.236 0.186 , 0.234	Depositor DCC
R_{free} test set	3383 reflections (3.80%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.123	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.019 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14189	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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: 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	0.72	0/2310	0.88	0/3126
1	B	0.75	0/2307	0.90	0/3120
1	C	0.75	0/2311	0.92	1/3128 (0.0%)
1	D	0.73	0/2339	0.88	1/3159 (0.0%)
1	E	0.65	0/2278	0.86	0/3089
1	F	0.74	2/2305 (0.1%)	0.87	0/3124
All	All	0.73	2/13850 (0.0%)	0.89	2/18746 (0.0%)

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

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	63	ARG	CZ-NH1	8.86	1.45	1.32
1	F	10	ARG	NE-CZ	6.95	1.40	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	108	PHE	N-CA-C	5.42	117.61	111.11
1	D	108	PHE	N-CA-C	5.12	117.25	111.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	260	LYS	Peptide

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	2267	0	2227	32	0
1	B	2264	0	2226	26	0
1	C	2268	0	2229	40	0
1	D	2289	0	2282	30	0
1	E	2235	0	2158	27	0
1	F	2259	0	2184	23	0
2	A	44	0	26	0	0
2	B	44	0	26	1	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
2	E	44	0	26	1	0
2	F	44	0	26	0	0
3	A	61	0	0	1	0
3	B	67	0	0	2	0
3	C	66	0	0	0	0
3	D	64	0	0	2	0
3	E	42	0	0	1	0
3	F	43	0	0	2	0
All	All	14189	0	13462	143	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 5.

All (143) 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:E:148:MET:HE2	1:E:171:LEU:HD22	1.40	1.03
1:E:258:ARG:HB2	1:E:264:MET:HE2	1.50	0.92
1:E:264:MET:CE	1:F:242:GLY:HA2	2.04	0.86
1:E:127:SER:HB3	1:E:148:MET:HE1	1.57	0.86
1:F:258:ARG:HB2	1:F:264:MET:HE2	1.60	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:ASP:O	1:A:121:ARG:HG3	1.79	0.83
1:E:148:MET:CE	1:E:171:LEU:HD22	2.08	0.83
1:A:217:TYR:HA	1:B:264:MET:HE1	1.62	0.82
1:A:217:TYR:HA	1:B:264:MET:CE	2.10	0.81
1:C:258:ARG:HB2	1:C:264:MET:HE2	1.63	0.81
1:C:242:GLY:HA2	1:D:264:MET:CE	2.14	0.77
1:A:217:TYR:CA	1:B:264:MET:HE1	2.17	0.75
1:B:206:PRO:HA	3:B:1063:HOH:O	1.86	0.74
1:A:217:TYR:CB	1:B:264:MET:HE1	2.18	0.74
1:A:198:SER:H	1:A:201:THR:HG22	1.51	0.74
1:E:264:MET:HE1	1:F:242:GLY:HA2	1.69	0.72
1:E:242:GLY:HA2	1:F:264:MET:CE	2.22	0.69
1:E:127:SER:CB	1:E:148:MET:HE1	2.23	0.69
1:E:264:MET:HE2	1:F:242:GLY:HA2	1.74	0.68
1:C:264:MET:CE	1:D:242:GLY:HA2	2.24	0.67
1:D:258:ARG:HB2	1:D:264:MET:HE2	1.78	0.66
1:C:115:ASP:O	1:C:163[B]:ASN:ND2	2.27	0.66
1:C:242:GLY:HA2	1:D:264:MET:HE2	1.76	0.66
1:A:217:TYR:HB3	1:B:264:MET:HE1	1.78	0.65
1:C:198:SER:H	1:C:201:THR:HG22	1.62	0.63
1:C:242:GLY:HA2	1:D:264:MET:HE1	1.81	0.62
1:E:250:GLU:HG3	1:E:284:LYS:HZ2	1.66	0.61
1:A:201:THR:HG23	1:A:205:MET:HE2	1.83	0.61
1:E:264:MET:HE1	1:F:217:TYR:HB3	1.84	0.59
1:A:196:ALA:O	1:A:201:THR:HG21	2.02	0.59
1:A:258:ARG:HD3	1:A:264:MET:HE2	1.86	0.58
1:E:258:ARG:HD3	1:E:264:MET:CE	2.33	0.58
1:C:117:ASP:O	1:C:121:ARG:HG3	2.04	0.58
1:C:264:MET:HE1	1:D:242:GLY:HA2	1.85	0.58
1:C:248:ARG:HG2	1:D:279:PHE:CE2	2.39	0.57
1:A:198:SER:H	1:A:201:THR:CG2	2.17	0.57
1:A:217:TYR:HA	1:B:264:MET:HE3	1.85	0.56
1:E:250:GLU:HG3	1:E:284:LYS:NZ	2.20	0.56
1:C:250:GLU:HG3	1:C:284:LYS:HZ3	1.71	0.56
1:E:217:TYR:HB3	1:F:264:MET:HE1	1.86	0.56
1:C:250:GLU:HG3	1:C:284:LYS:NZ	2.21	0.56
1:E:242:GLY:HA2	1:F:264:MET:HE1	1.87	0.56
1:C:264:MET:HE1	1:D:217:TYR:HB3	1.89	0.54
1:F:57:LYS:O	1:F:61:GLU:HG3	2.07	0.54
1:C:135:TRP:NE1	1:C:139:LYS:HD2	2.23	0.54
1:B:57:LYS:O	1:B:61:GLU:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:ILE:HD13	1:D:209:LEU:HD21	1.90	0.53
1:E:148:MET:HE3	3:E:1029:HOH:O	2.09	0.53
1:D:95:VAL:HG22	1:D:145:ARG:HB2	1.89	0.53
1:A:198:SER:N	1:A:201:THR:HG22	2.22	0.53
1:A:238:GLU:HG3	1:A:247:LEU:HD11	1.90	0.53
1:A:61:GLU:OE1	1:A:64:ARG:NH1	2.42	0.52
1:C:5:LEU:HD12	1:D:266:PRO:HB2	1.90	0.52
1:C:238:GLU:OE2	1:D:251:ARG:HD2	2.09	0.52
1:A:115:ASP:O	1:A:163[B]:ASN:ND2	2.40	0.52
1:B:117:ASP:O	1:B:121:ARG:HG3	2.10	0.52
1:E:248:ARG:CZ	1:E:284:LYS:HG2	2.40	0.52
1:E:258:ARG:HD3	1:E:264:MET:HE3	1.91	0.52
1:F:258:ARG:HD3	1:F:264:MET:CE	2.40	0.51
1:F:248:ARG:CZ	1:F:284:LYS:HG2	2.41	0.51
1:D:56:ASP:OD1	1:D:73:TYR:OH	2.29	0.51
1:B:248:ARG:CZ	1:B:284:LYS:HG2	2.41	0.51
1:A:15:THR:HG1	1:A:99:ASN:H	1.59	0.51
1:B:294:SER:HA	1:B:297:GLU:HG2	1.93	0.51
1:A:116:TRP:HA	1:A:163[B]:ASN:ND2	2.26	0.50
1:B:200:MET:HE1	2:B:1002:NAD:H2D	1.94	0.50
1:A:41:LEU:HB3	3:A:1011:HOH:O	2.12	0.49
1:F:117:ASP:O	1:F:121:ARG:HG2	2.12	0.49
1:D:37:VAL:HG21	1:D:86:ALA:HB2	1.94	0.49
1:B:247:LEU:HD12	1:B:247:LEU:N	2.26	0.49
1:D:103:LEU:HD11	1:D:200:MET:HG3	1.93	0.49
1:F:23[A]:ARG:HD3	3:F:1024:HOH:O	2.11	0.49
1:E:242:GLY:HA2	1:F:264:MET:HE2	1.93	0.49
1:F:250:GLU:HG3	1:F:284:LYS:HZ2	1.78	0.49
1:A:264:MET:CE	1:B:242:GLY:HA2	2.42	0.49
1:F:258:ARG:HD3	1:F:264:MET:HE2	1.94	0.49
1:A:264:MET:CE	1:B:217:TYR:HA	2.43	0.49
1:C:12:VAL:HG22	1:C:95:VAL:HB	1.96	0.48
1:E:270:LYS:HG3	1:F:224:TRP:CH2	2.49	0.48
1:D:87:LEU:HD11	1:D:137:HIS:CE1	2.49	0.47
1:C:247:LEU:CD1	1:D:251:ARG:HG3	2.45	0.47
1:A:264:MET:HE1	1:B:217:TYR:HB3	1.97	0.46
1:C:260:LYS:HA	1:C:261:ASN:HD22	1.80	0.46
1:A:197:GLY:HA2	1:A:201:THR:CG2	2.45	0.46
1:F:12:VAL:HG22	1:F:95:VAL:HB	1.96	0.46
1:E:25:TYR:CZ	1:E:222:VAL:HG21	2.51	0.46
1:C:217:TYR:HB3	1:D:264:MET:HE1	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:GLU:HG3	1:B:247:LEU:HD11	1.97	0.46
1:D:142:LYS:HE3	3:D:1046:HOH:O	2.15	0.46
1:A:197:GLY:HA2	1:A:201:THR:HG21	1.97	0.46
1:C:116:TRP:HA	1:C:163[B]:ASN:ND2	2.31	0.46
1:C:154:GLY:HA2	1:C:165:SER:HB2	1.98	0.45
1:B:59:VAL:HG21	1:B:71:ALA:HB2	1.98	0.45
1:B:44:ASP:OD1	1:B:44:ASP:C	2.60	0.45
1:C:178:LEU:HD21	1:D:108:PHE:CD1	2.52	0.45
1:E:255:ALA:HB2	1:E:275:LYS:HD3	1.99	0.45
1:D:125:ARG:HD3	3:D:1013:HOH:O	2.17	0.45
1:E:208:ASP:OD1	1:E:208:ASP:N	2.50	0.44
1:C:197:GLY:HA2	1:C:201:THR:HG21	2.00	0.44
1:F:41:LEU:HD13	1:F:75:SER:HB2	1.99	0.44
1:A:258:ARG:HD3	1:A:264:MET:CE	2.48	0.44
1:C:95:VAL:HG22	1:C:145:ARG:HB2	2.00	0.44
1:D:246:LYS:C	1:D:247:LEU:HD12	2.42	0.44
2:E:1005:NAD:O1N	2:E:1005:NAD:N7N	2.49	0.44
1:C:238:GLU:HG3	1:C:247:LEU:HD11	2.00	0.44
1:C:135:TRP:HE1	1:C:139:LYS:HD2	1.81	0.43
1:F:39:ASN:ND2	1:F:72:ASN:OD1	2.48	0.43
1:D:78:GLU:HB3	1:D:81:LYS:HG3	2.00	0.43
1:D:62:ILE:HG21	1:D:69:ALA:HB2	2.01	0.43
1:A:12:VAL:HG11	1:A:29:PHE:CD1	2.54	0.43
1:A:258:ARG:HB2	1:A:264:MET:HE2	2.00	0.43
1:C:264:MET:HE2	1:D:242:GLY:HA2	1.99	0.43
1:C:197:GLY:HA2	1:C:201:THR:CG2	2.49	0.43
1:C:115:ASP:C	1:C:163[B]:ASN:HD21	2.21	0.42
1:A:248:ARG:CZ	1:A:284:LYS:HG2	2.50	0.42
1:B:32[A]:ARG:HG2	1:B:32[A]:ARG:HH11	1.85	0.42
1:B:251:ARG:NH1	3:B:1026:HOH:O	2.52	0.42
1:E:15:THR:O	1:E:99:ASN:HB3	2.20	0.42
1:A:105:ASP:OD1	1:A:161:GLN:HG2	2.19	0.42
1:A:264:MET:HE1	1:B:217:TYR:CB	2.49	0.42
1:D:198:SER:O	1:D:202:GLN:HG3	2.19	0.42
1:D:103:LEU:CD1	1:D:200:MET:HG3	2.49	0.42
1:C:87:LEU:HD11	1:C:137:HIS:CE1	2.54	0.42
1:B:264:MET:HE3	1:B:264:MET:HB3	1.83	0.41
1:C:62:ILE:HG21	1:C:69:ALA:HB2	2.02	0.41
1:C:248:ARG:CZ	1:C:284:LYS:HG2	2.50	0.41
1:F:151:SER:HB2	3:F:1035:HOH:O	2.20	0.41
1:A:243:TRP:CE3	1:B:295:ILE:HD12	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:25:TYR:CE1	1:E:222:VAL:HG21	2.55	0.41
1:A:44:ASP:OD1	1:A:44:ASP:C	2.64	0.41
1:B:246:LYS:C	1:B:247:LEU:HD12	2.46	0.41
1:C:53:LEU:HD23	1:C:53:LEU:H	1.86	0.41
1:C:248:ARG:HG2	1:D:279:PHE:CZ	2.55	0.41
1:D:247:LEU:HD12	1:D:247:LEU:N	2.36	0.41
1:E:135:TRP:CZ3	1:E:178:LEU:HD22	2.56	0.41
1:E:201:THR:HB	1:E:205:MET:HE2	2.02	0.41
1:C:141:GLN:HE21	1:C:141:GLN:HB2	1.67	0.40
1:C:181:GLU:HG2	1:D:159:PHE:CD1	2.56	0.40
1:B:294:SER:O	1:B:297:GLU:HG2	2.21	0.40
1:F:15:THR:HA	1:F:39:ASN:HB3	2.02	0.40
1:C:255:ALA:HB2	1:C:275:LYS:HD3	2.04	0.40
1:F:214:LYS:HA	1:F:215:PRO:HD2	1.88	0.40
1:C:214:LYS:HB3	1:C:216:GLU:HG2	2.04	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	302/327 (92%)	295 (98%)	7 (2%)	0	100	100
1	B	301/327 (92%)	297 (99%)	4 (1%)	0	100	100
1	C	301/327 (92%)	296 (98%)	5 (2%)	0	100	100
1	D	303/327 (93%)	296 (98%)	7 (2%)	0	100	100
1	E	301/327 (92%)	294 (98%)	7 (2%)	0	100	100
1	F	302/327 (92%)	296 (98%)	6 (2%)	0	100	100
All	All	1810/1962 (92%)	1774 (98%)	36 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	222/258 (86%)	216 (97%)	6 (3%)	39	53
1	B	222/258 (86%)	219 (99%)	3 (1%)	59	75
1	C	225/258 (87%)	219 (97%)	6 (3%)	39	53
1	D	229/258 (89%)	222 (97%)	7 (3%)	35	48
1	E	216/258 (84%)	213 (99%)	3 (1%)	59	75
1	F	221/258 (86%)	218 (99%)	3 (1%)	59	75
All	All	1335/1548 (86%)	1307 (98%)	28 (2%)	47	63

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

Mol	Chain	Res	Type
1	A	3	SER
1	A	41	LEU
1	A	53	LEU
1	A	107	SER
1	A	203	THR
1	A	296	ILE
1	B	3	SER
1	B	204	VAL
1	B	216	GLU
1	C	53	LEU
1	C	201	THR
1	C	216	GLU
1	C	261	ASN
1	C	297	GLU
1	C	300	SER
1	D	3	SER
1	D	31	GLU
1	D	114	GLU
1	D	136	GLU
1	D	203	THR
1	D	204	VAL
1	D	286	GLN

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Mol	Chain	Res	Type
1	E	87	LEU
1	E	203	THR
1	E	270	LYS
1	F	284	LYS
1	F	286	GLN
1	F	301	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	120	HIS
1	A	158	ASN
1	A	161	GLN
1	A	272	ASN
1	B	120	HIS
1	B	158	ASN
1	B	286	GLN
1	C	120	HIS
1	C	137	HIS
1	C	141	GLN
1	C	158	ASN
1	C	161	GLN
1	C	261	ASN
1	D	98	ASN
1	D	120	HIS
1	D	137	HIS
1	D	262	HIS
1	D	286	GLN
1	E	137	HIS
1	E	262	HIS
1	F	98	ASN
1	F	120	HIS
1	F	141	GLN
1	F	262	HIS
1	F	272	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](#)

6 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	1002	-	46,48,48	1.71	7 (15%)	64,73,73	1.68	11 (17%)
2	NAD	F	1006	-	46,48,48	1.78	6 (13%)	64,73,73	1.73	13 (20%)
2	NAD	D	1004	-	46,48,48	1.82	7 (15%)	64,73,73	1.74	13 (20%)
2	NAD	A	1001	-	46,48,48	1.83	7 (15%)	64,73,73	1.73	10 (15%)
2	NAD	C	1003	-	46,48,48	1.68	5 (10%)	64,73,73	1.73	11 (17%)
2	NAD	E	1005	-	46,48,48	1.72	4 (8%)	64,73,73	1.70	9 (14%)

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	1002	-	-	5/30/62/62	0/5/5/5
2	NAD	F	1006	-	-	2/30/62/62	0/5/5/5
2	NAD	D	1004	-	-	4/30/62/62	0/5/5/5
2	NAD	A	1001	-	-	1/30/62/62	0/5/5/5
2	NAD	C	1003	-	-	5/30/62/62	0/5/5/5
2	NAD	E	1005	-	-	5/30/62/62	0/5/5/5

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	NAD	O7N-C7N	9.17	1.41	1.24
2	C	1003	NAD	O7N-C7N	8.92	1.40	1.24
2	F	1006	NAD	O7N-C7N	8.90	1.40	1.24
2	B	1002	NAD	O7N-C7N	8.90	1.40	1.24
2	E	1005	NAD	O7N-C7N	8.83	1.40	1.24
2	D	1004	NAD	O7N-C7N	8.44	1.39	1.24
2	A	1001	NAD	C2A-N3A	3.42	1.40	1.33
2	D	1004	NAD	PA-O3	3.41	1.63	1.59
2	A	1001	NAD	C2A-N1A	3.17	1.39	1.33
2	F	1006	NAD	C2A-N3A	3.06	1.39	1.33
2	F	1006	NAD	C2A-N1A	3.03	1.39	1.33
2	D	1004	NAD	C2N-N1N	2.91	1.38	1.35
2	F	1006	NAD	C2N-N1N	2.89	1.38	1.35
2	B	1002	NAD	PN-O3	2.83	1.62	1.59
2	E	1005	NAD	C2A-N3A	2.82	1.38	1.33
2	D	1004	NAD	C2A-N1A	2.78	1.38	1.33
2	A	1001	NAD	PA-O3	2.65	1.62	1.59
2	D	1004	NAD	C8A-N7A	2.57	1.36	1.31
2	A	1001	NAD	C2N-N1N	2.54	1.37	1.35
2	C	1003	NAD	C2A-N1A	2.52	1.38	1.33
2	D	1004	NAD	C2A-N3A	2.49	1.38	1.33
2	E	1005	NAD	C2A-N1A	2.47	1.38	1.33
2	A	1001	NAD	C8A-N7A	2.46	1.36	1.31
2	F	1006	NAD	PA-O3	2.45	1.62	1.59
2	B	1002	NAD	C5A-N7A	-2.43	1.34	1.39
2	E	1005	NAD	PA-O3	2.37	1.62	1.59
2	D	1004	NAD	C5N-C4N	2.31	1.42	1.38
2	C	1003	NAD	C2A-N3A	2.30	1.38	1.33
2	A	1001	NAD	PN-O3	2.29	1.62	1.59
2	F	1006	NAD	C8A-N7A	2.29	1.36	1.31
2	C	1003	NAD	C5A-N7A	-2.21	1.35	1.39
2	B	1002	NAD	C8A-N7A	2.20	1.35	1.31
2	C	1003	NAD	PA-O3	-2.17	1.57	1.59
2	B	1002	NAD	C2A-N3A	2.14	1.37	1.33
2	B	1002	NAD	C2A-N1A	2.11	1.37	1.33
2	B	1002	NAD	PA-O2A	-2.01	1.46	1.55

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1005	NAD	N3A-C2A-N1A	-6.83	118.25	128.58
2	B	1002	NAD	N3A-C2A-N1A	-6.70	118.44	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1003	NAD	N3A-C2A-N1A	-6.29	119.06	128.58
2	F	1006	NAD	N3A-C2A-N1A	-6.29	119.06	128.58
2	A	1001	NAD	N3A-C2A-N1A	-5.84	119.74	128.58
2	D	1004	NAD	N3A-C2A-N1A	-5.81	119.79	128.58
2	A	1001	NAD	N9A-C8A-N7A	-4.79	107.14	113.94
2	F	1006	NAD	C5A-C4A-N3A	-4.61	120.37	126.72
2	D	1004	NAD	N9A-C8A-N7A	-4.56	107.47	113.94
2	C	1003	NAD	N9A-C8A-N7A	-4.37	107.74	113.94
2	E	1005	NAD	C5A-C4A-N3A	-4.25	120.86	126.72
2	F	1006	NAD	O4B-C1B-N9A	4.22	116.20	108.09
2	B	1002	NAD	N9A-C8A-N7A	-4.20	107.98	113.94
2	A	1001	NAD	C5A-C4A-N3A	-4.04	121.16	126.72
2	C	1003	NAD	C5A-C4A-N3A	-3.91	121.34	126.72
2	A	1001	NAD	O4B-C1B-N9A	3.88	115.54	108.09
2	A	1001	NAD	C5A-N7A-C8A	3.87	109.54	103.45
2	E	1005	NAD	C2A-N3A-C4A	3.85	121.23	111.83
2	E	1005	NAD	O4B-C1B-N9A	3.84	115.46	108.09
2	E	1005	NAD	N9A-C8A-N7A	-3.81	108.53	113.94
2	F	1006	NAD	C2A-N3A-C4A	3.72	120.92	111.83
2	D	1004	NAD	C3N-C7N-N7N	3.66	122.25	117.74
2	B	1002	NAD	O4B-C1B-N9A	3.65	115.10	108.09
2	F	1006	NAD	N9A-C8A-N7A	-3.63	108.78	113.94
2	C	1003	NAD	C3N-C7N-N7N	3.57	122.14	117.74
2	B	1002	NAD	C5A-C4A-N3A	-3.57	121.80	126.72
2	B	1002	NAD	C5A-N7A-C8A	3.51	108.97	103.45
2	D	1004	NAD	C5A-N7A-C8A	3.47	108.90	103.45
2	C	1003	NAD	C5A-N7A-C8A	3.46	108.89	103.45
2	D	1004	NAD	C5A-C4A-N3A	-3.43	122.00	126.72
2	C	1003	NAD	C2A-N3A-C4A	3.37	120.05	111.83
2	B	1002	NAD	C2A-N3A-C4A	3.36	120.04	111.83
2	A	1001	NAD	C2A-N3A-C4A	3.33	119.97	111.83
2	F	1006	NAD	C5A-N7A-C8A	3.18	108.44	103.45
2	D	1004	NAD	C2A-N3A-C4A	3.13	119.48	111.83
2	F	1006	NAD	C3N-C7N-N7N	3.08	121.53	117.74
2	A	1001	NAD	C4A-C5A-N7A	-3.04	107.10	110.58
2	E	1005	NAD	C5A-N7A-C8A	3.00	108.16	103.45
2	F	1006	NAD	N3A-C4A-N9A	2.99	132.25	127.17
2	C	1003	NAD	O2N-PN-O1N	2.94	126.12	112.44
2	E	1005	NAD	N3A-C4A-N9A	2.75	131.84	127.17
2	D	1004	NAD	O7N-C7N-C3N	-2.73	116.26	119.60
2	D	1004	NAD	C4A-N9A-C8A	2.66	108.53	105.74
2	B	1002	NAD	O2N-PN-O3	2.64	114.42	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1003	NAD	N3A-C4A-N9A	2.61	131.61	127.17
2	D	1004	NAD	C5N-C4N-C3N	-2.59	117.82	120.36
2	C	1003	NAD	O4B-C1B-N9A	2.58	113.04	108.09
2	F	1006	NAD	O2A-PA-O3	2.57	114.22	107.27
2	A	1001	NAD	C4A-N9A-C8A	2.56	108.43	105.74
2	A	1001	NAD	C4A-N9A-C1B	-2.45	120.90	126.63
2	C	1003	NAD	O7N-C7N-C3N	-2.40	116.66	119.60
2	B	1002	NAD	C4A-C5A-N7A	-2.38	107.86	110.58
2	D	1004	NAD	O2A-PA-O3	2.31	113.51	107.27
2	A	1001	NAD	N3A-C4A-N9A	2.28	131.05	127.17
2	F	1006	NAD	C4A-C5A-N7A	-2.27	107.98	110.58
2	D	1004	NAD	C6N-N1N-C2N	-2.27	119.95	121.88
2	D	1004	NAD	C4A-C5A-N7A	-2.27	107.99	110.58
2	F	1006	NAD	O2N-PN-O1N	2.19	122.65	112.44
2	B	1002	NAD	C5A-C4A-N9A	2.19	108.20	105.81
2	D	1004	NAD	N3A-C4A-N9A	2.19	130.89	127.17
2	E	1005	NAD	C3N-C7N-N7N	2.11	120.34	117.74
2	F	1006	NAD	O7N-C7N-C3N	-2.10	117.02	119.60
2	C	1003	NAD	C4A-N9A-C8A	2.06	107.91	105.74
2	F	1006	NAD	O3-PA-O1A	-2.06	104.52	110.70
2	B	1002	NAD	C2A-N1A-C6A	2.05	122.10	118.73
2	B	1002	NAD	O3B-C3B-C4B	-2.03	105.26	111.08
2	E	1005	NAD	C2N-C3N-C4N	2.01	120.60	118.26

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1002	NAD	C5D-O5D-PN-O3
2	B	1002	NAD	C5D-O5D-PN-O1N
2	B	1002	NAD	C5D-O5D-PN-O2N
2	C	1003	NAD	C5D-O5D-PN-O3
2	C	1003	NAD	C5D-O5D-PN-O1N
2	C	1003	NAD	C5D-O5D-PN-O2N
2	D	1004	NAD	C5D-O5D-PN-O3
2	D	1004	NAD	C5D-O5D-PN-O1N
2	D	1004	NAD	C5D-O5D-PN-O2N
2	E	1005	NAD	C5D-O5D-PN-O3
2	E	1005	NAD	C5D-O5D-PN-O1N
2	E	1005	NAD	C5D-O5D-PN-O2N
2	F	1006	NAD	C5D-O5D-PN-O2N
2	B	1002	NAD	PA-O3-PN-O2N

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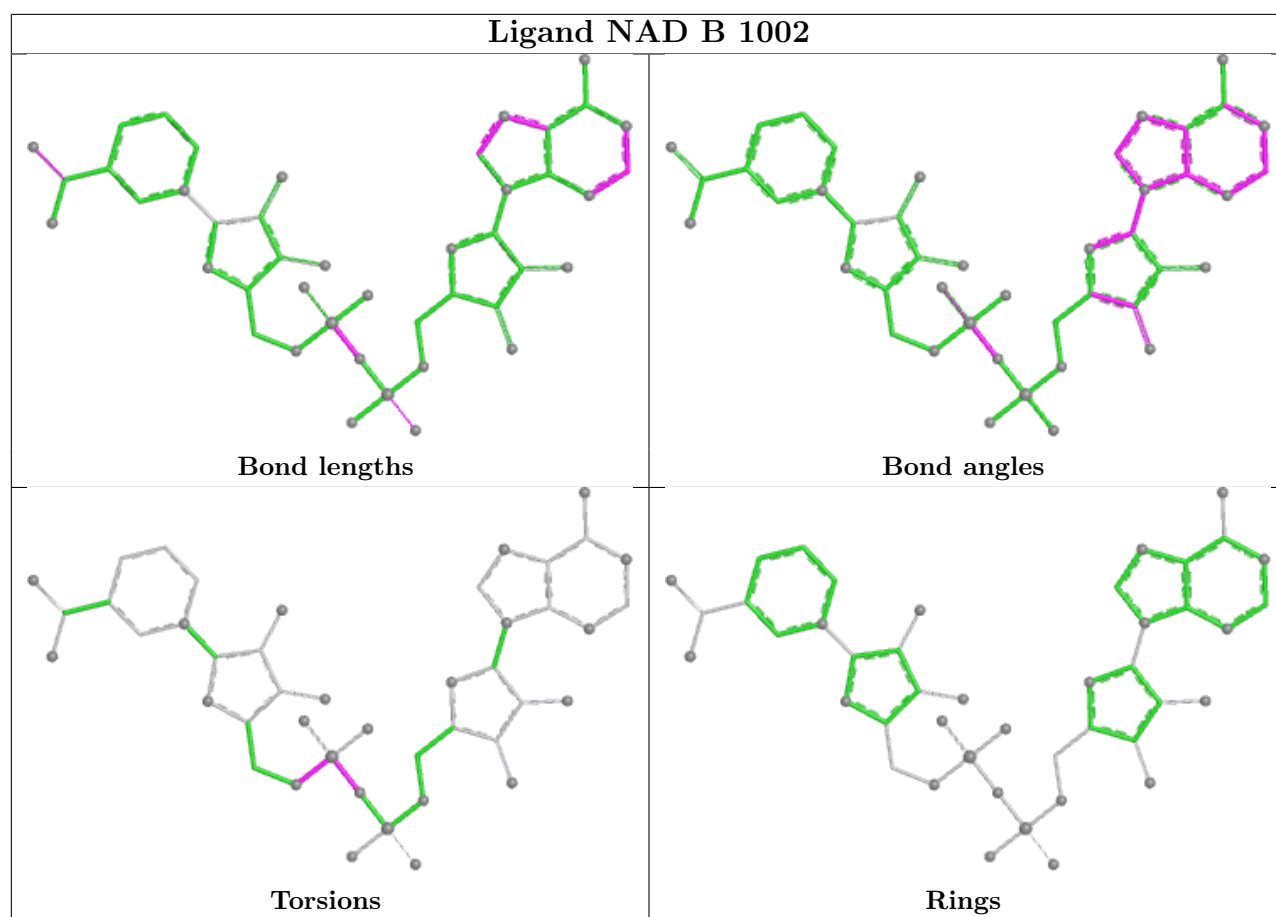
Mol	Chain	Res	Type	Atoms
2	E	1005	NAD	PA-O3-PN-O2N
2	A	1001	NAD	PA-O3-PN-O1N
2	B	1002	NAD	PA-O3-PN-O1N
2	E	1005	NAD	PA-O3-PN-O1N
2	F	1006	NAD	PA-O3-PN-O1N
2	C	1003	NAD	PA-O3-PN-O2N
2	D	1004	NAD	PA-O3-PN-O1N
2	C	1003	NAD	O4B-C4B-C5B-O5B

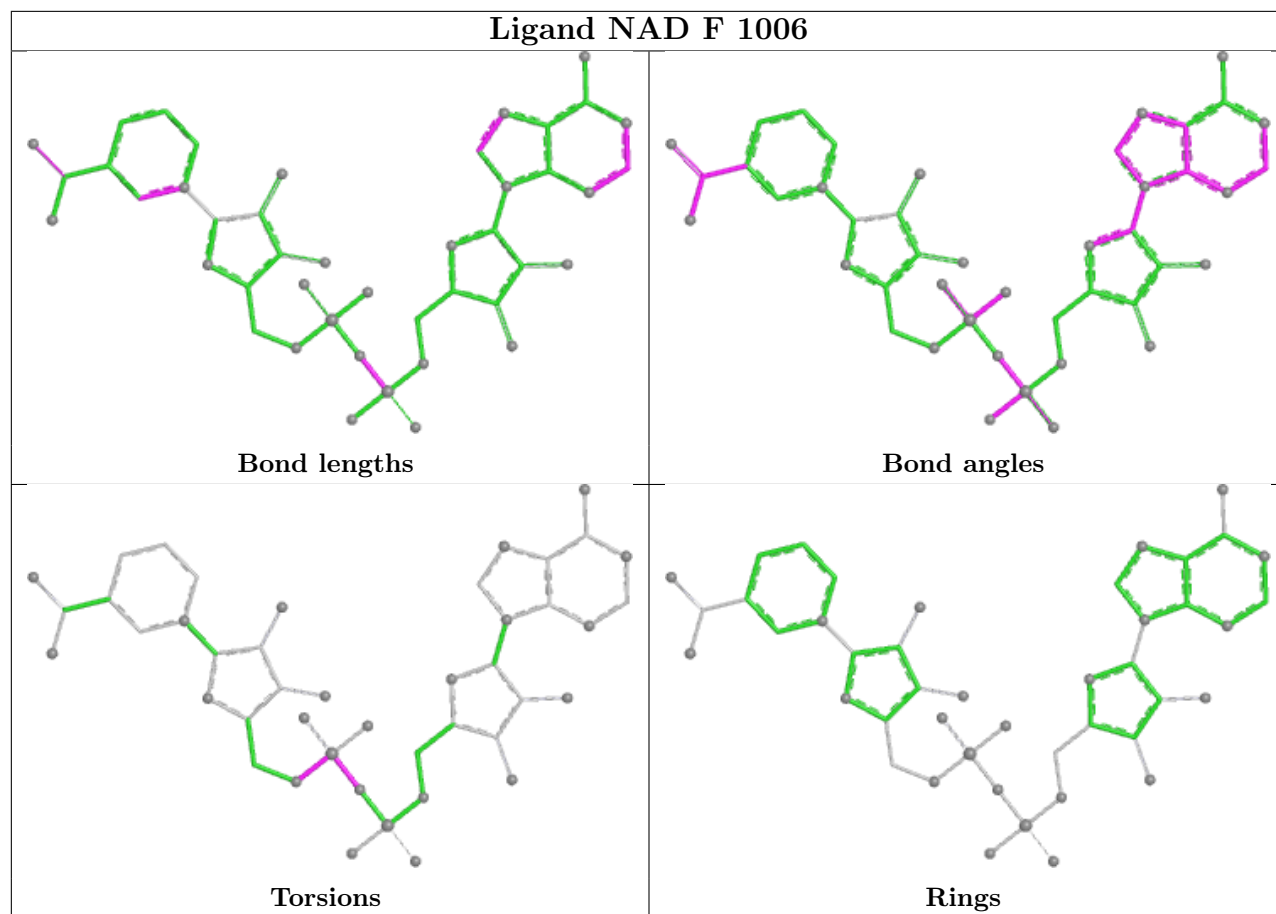
There are no ring outliers.

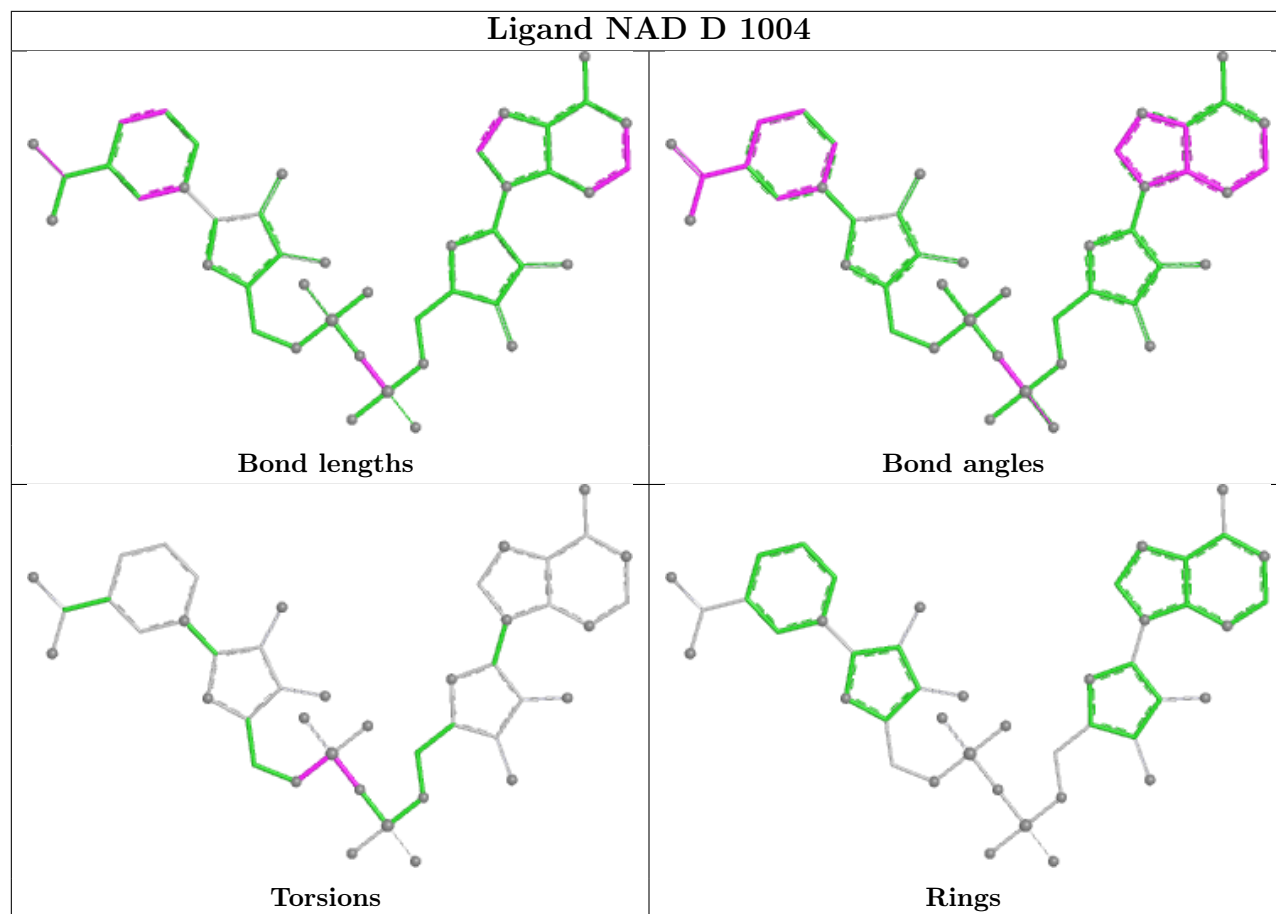
2 monomers are involved in 2 short contacts:

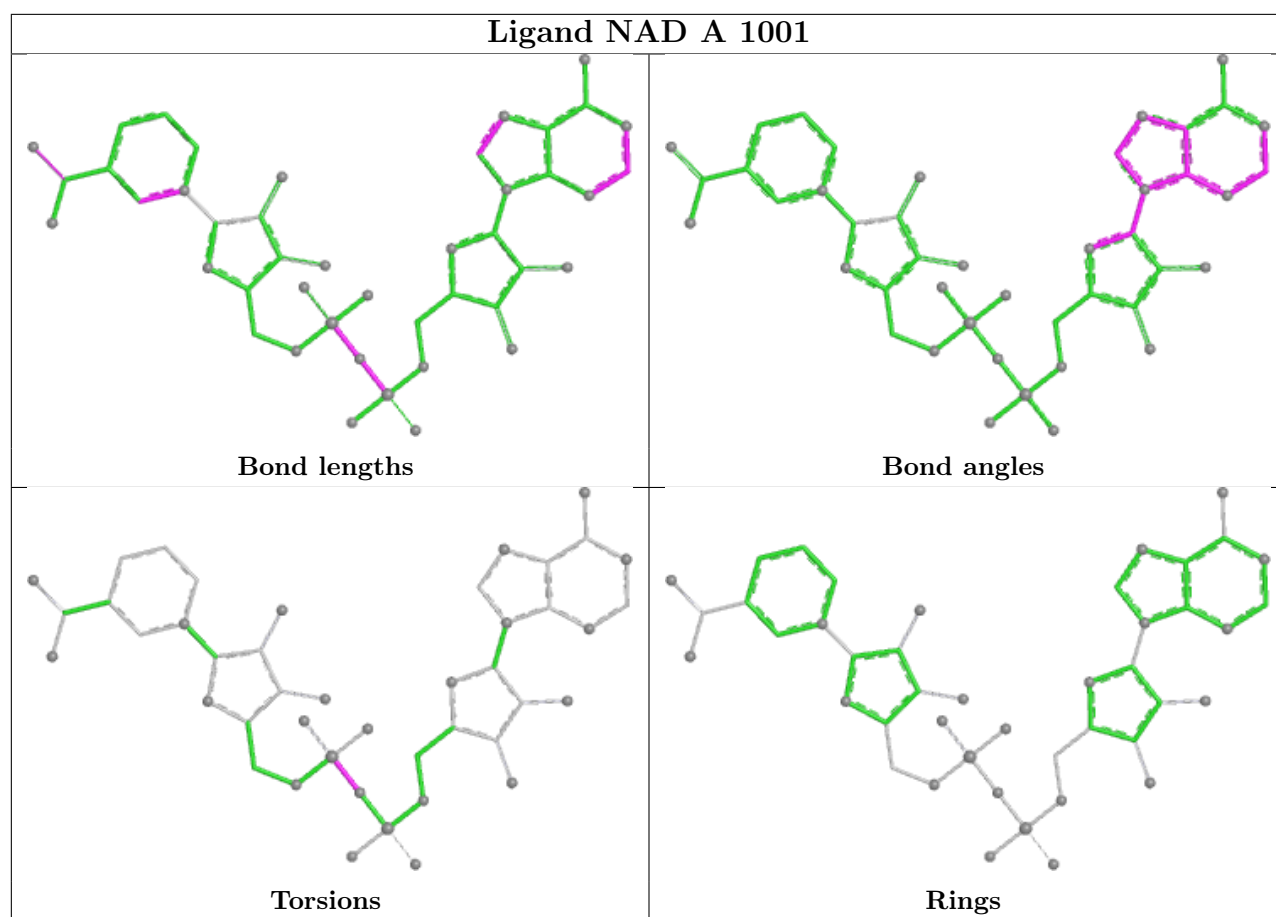
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1002	NAD	1	0
2	E	1005	NAD	1	0

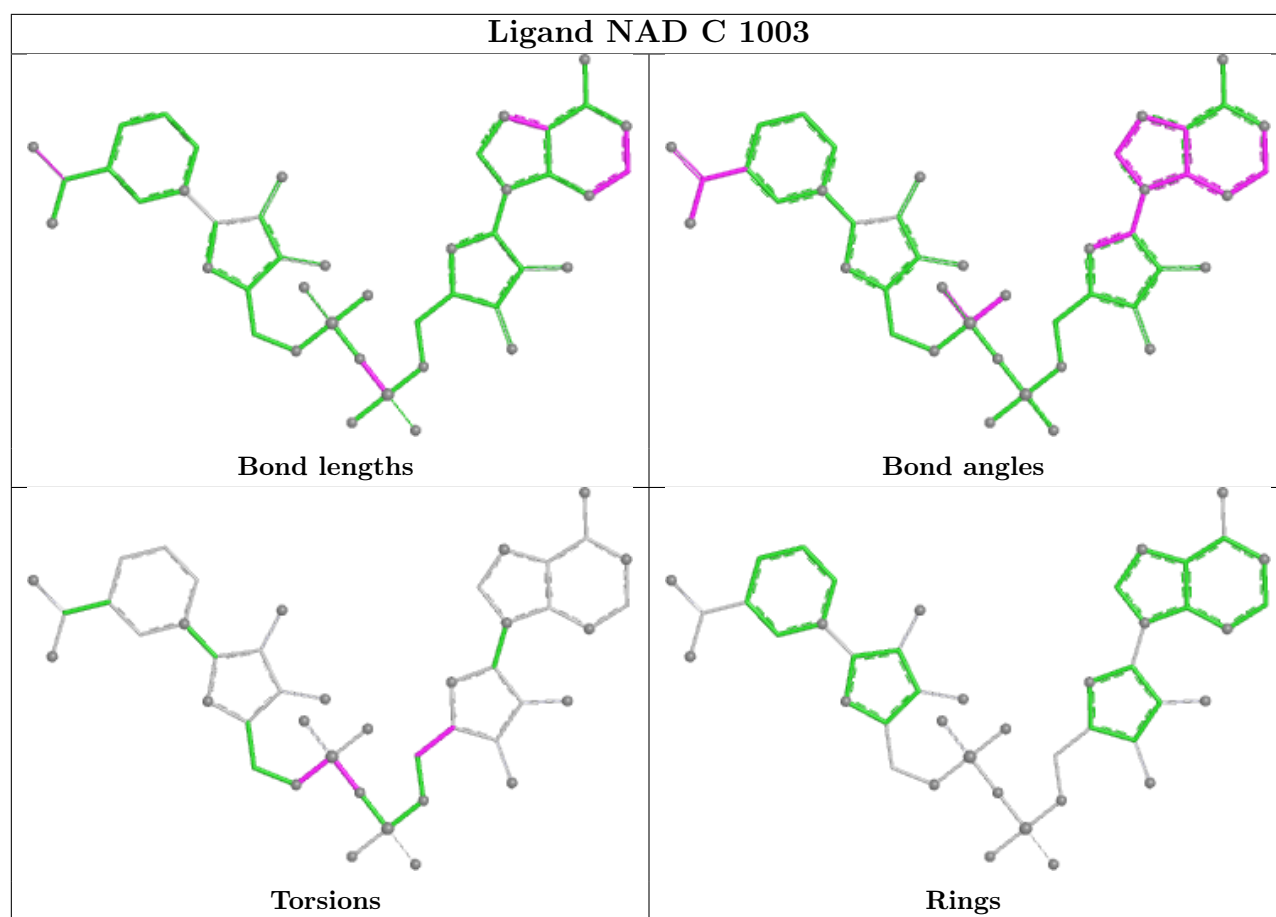
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

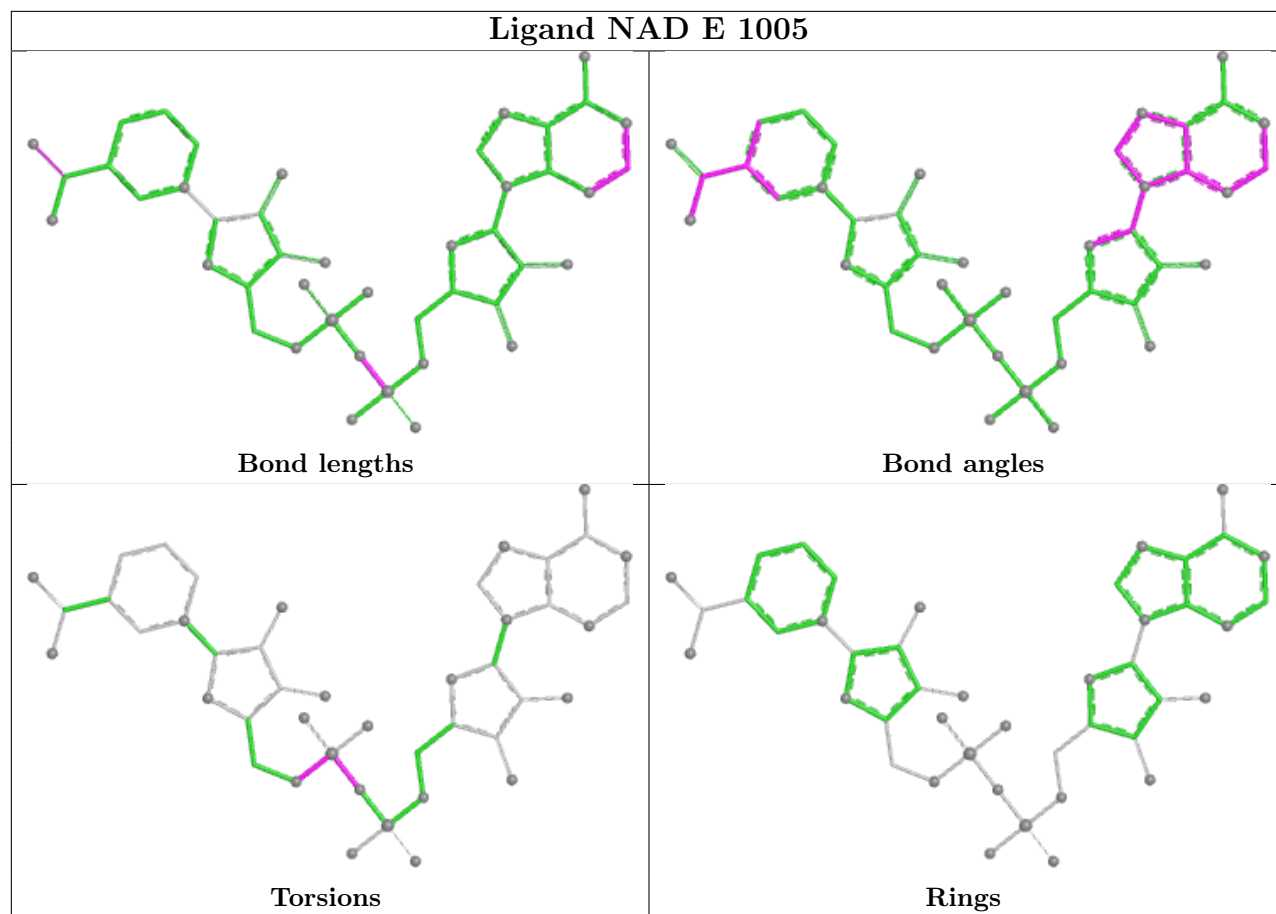












5.7 Other polymers [i](#)

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	302/327 (92%)	0.42	2 (0%) 84 82	17, 39, 47, 54	2 (0%)
1	B	302/327 (92%)	0.40	2 (0%) 84 82	20, 39, 47, 54	1 (0%)
1	C	302/327 (92%)	0.48	0 100 100	24, 39, 46, 54	1 (0%)
1	D	302/327 (92%)	0.39	7 (2%) 61 58	20, 38, 47, 54	4 (1%)
1	E	302/327 (92%)	0.86	26 (8%) 16 14	19, 39, 47, 54	1 (0%)
1	F	302/327 (92%)	0.67	5 (1%) 69 66	20, 39, 46, 54	2 (0%)
All	All	1812/1962 (92%)	0.54	42 (2%) 61 58	17, 39, 47, 54	11 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	71	ALA	3.7
1	E	252	THR	3.1
1	E	70	VAL	3.0
1	E	264	MET	3.0
1	E	265	THR	2.9
1	E	263	PRO	2.9
1	E	302	ILE	2.9
1	D	304	SER	2.9
1	E	262	HIS	2.9
1	E	273	TRP	2.9
1	E	256	ILE	2.8
1	E	298	VAL	2.7
1	F	53	LEU	2.5
1	E	257	VAL	2.5
1	D	94	ASP	2.4
1	E	303	ASP	2.4
1	E	282	ALA	2.4
1	E	276	ILE	2.4
1	D	208	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	304	SER	2.4
1	D	299	LEU	2.3
1	E	5	LEU	2.3
1	E	269	VAL	2.3
1	E	162	ALA	2.3
1	E	253	LEU	2.3
1	F	224	TRP	2.3
1	B	51	GLY	2.2
1	E	51	GLY	2.2
1	B	202	GLN	2.2
1	E	268	ALA	2.2
1	D	53	LEU	2.1
1	A	79	GLY	2.1
1	F	101	GLY	2.1
1	E	270	LYS	2.1
1	A	94	ASP	2.1
1	D	257	VAL	2.1
1	E	204	VAL	2.1
1	F	3	SER	2.1
1	E	293	GLY	2.0
1	D	286	GLN	2.0
1	F	5	LEU	2.0
1	E	18	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.

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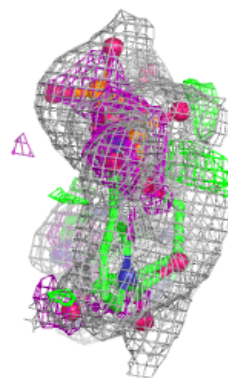
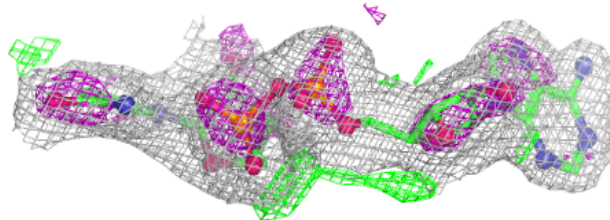
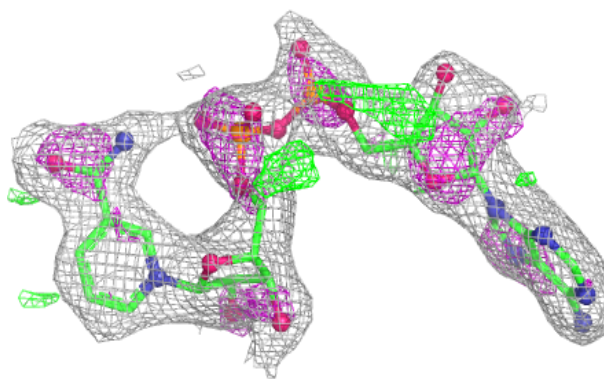
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAD	F	1006	44/44	0.91	0.10	28,34,42,45	0
2	NAD	C	1003	44/44	0.92	0.10	21,27,31,32	0
2	NAD	D	1004	44/44	0.92	0.09	25,30,32,34	0
2	NAD	E	1005	44/44	0.92	0.10	31,37,40,41	0
2	NAD	B	1002	44/44	0.92	0.10	20,28,31,33	0
2	NAD	A	1001	44/44	0.93	0.09	22,27,33,34	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.

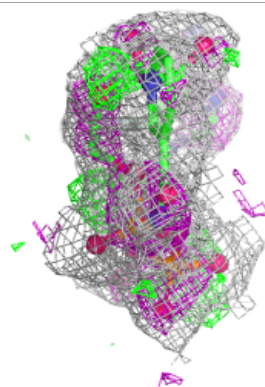
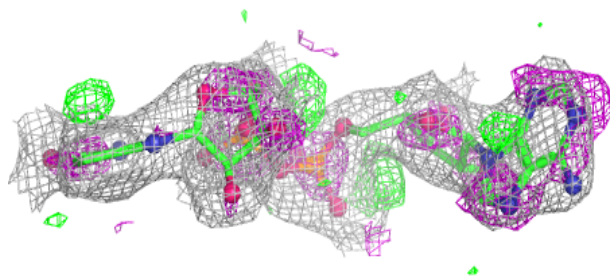
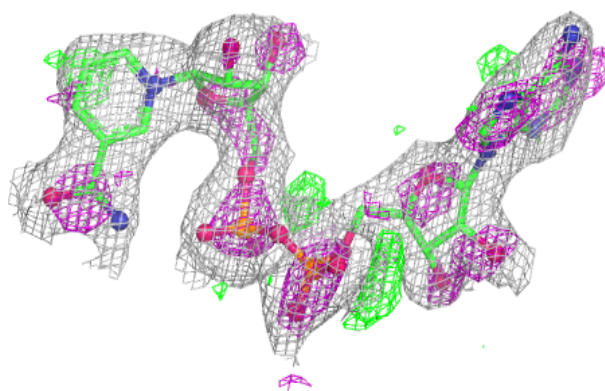
Electron density around NAD F 1006:

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

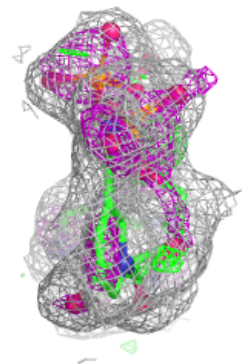
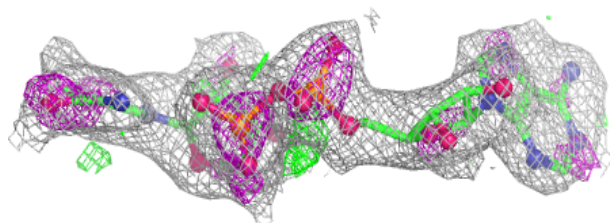
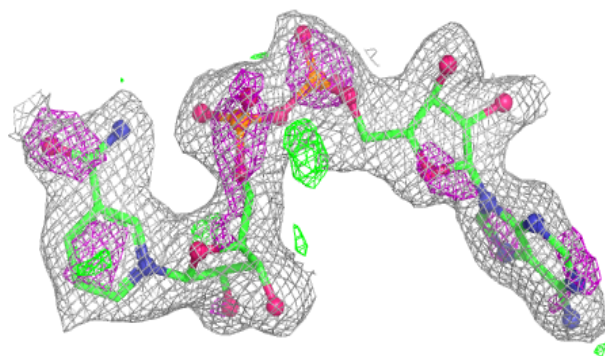


Electron density around NAD C 1003:

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

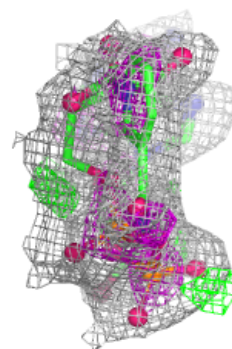
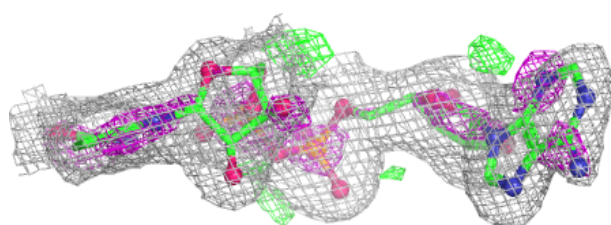
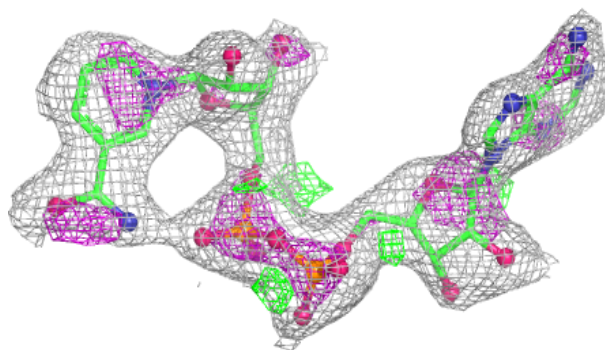
**Electron density around NAD D 1004:**

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

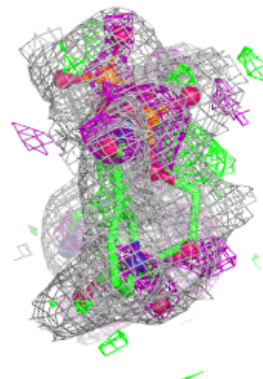
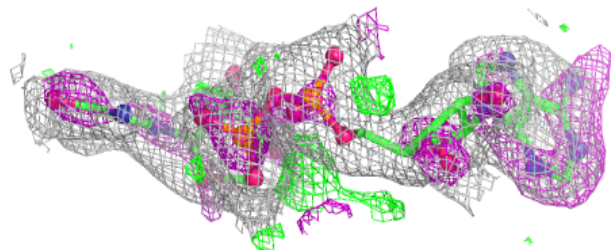
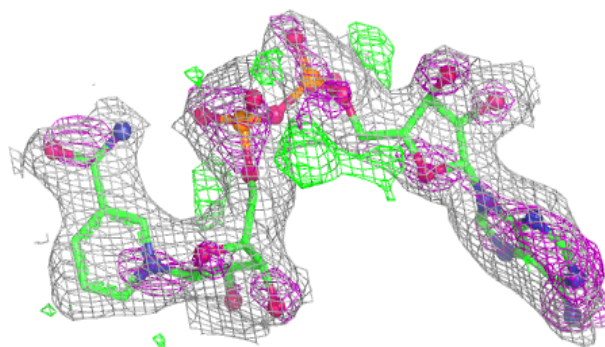


Electron density around NAD E 1005:

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

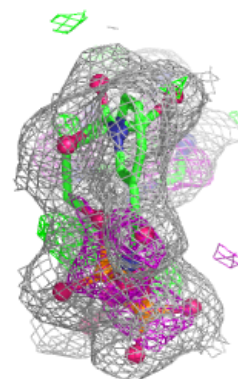
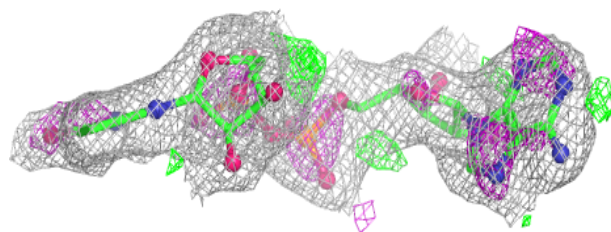
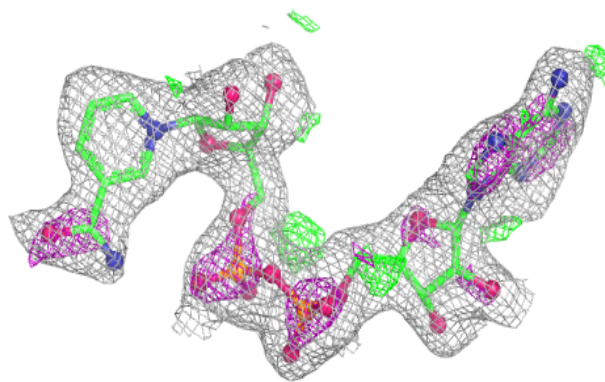
**Electron density around NAD B 1002:**

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



Electron density around NAD A 1001:

$2mF_o - DF_c$ (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.