



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:21 AM UTC

PDB ID : 8BR4 / pdb_00008br4
Title : Structure of GAPDH from Mycobacterium tuberculosis
Authors : Kumar, A.; Karthikeyan, S.
Deposited on : 2022-11-22
Resolution : 3.29 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

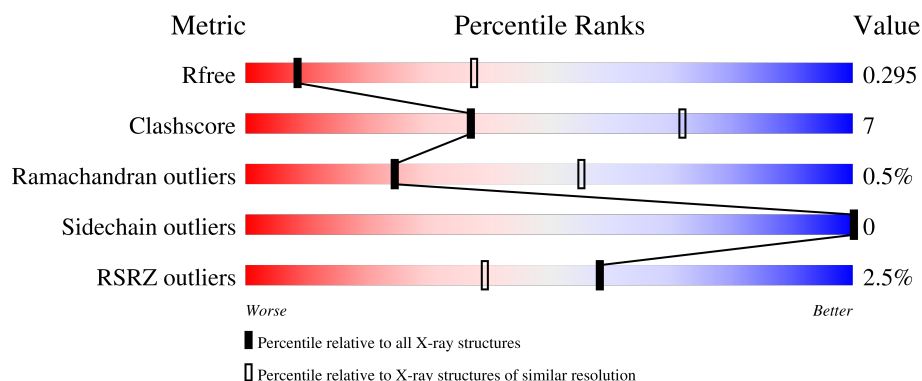
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.29 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	357	0% 82% 13% 5%
1	B	357	3% 78% 17% 5%
1	C	357	2% 82% 13% 5%
1	D	357	2% 78% 16% 5%
1	E	357	3% 78% 17% 5%

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Mol	Chain	Length	Quality of chain
1	F	357	2% 80% 14% 5%
1	G	357	3% 80% 15% 5%
1	H	357	2% 81% 13% 5%
1	I	357	% 82% 12% 5%
1	J	357	2% 81% 14% 5%
1	K	357	4% 80% 15% 5%
1	L	357	3% 81% 14% 5%
1	M	357	2% 77% 18% 5%
1	N	357	3% 79% 16% 5%
1	O	357	4% 79% 16% 5%
1	P	357	% 82% 13% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 38389 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 Glyceraldehyde-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	0	0
			2363	1464	411	483	5			
1	B	338	Total	C	N	O	S	0	0	0
			2364	1467	412	480	5			
1	C	339	Total	C	N	O	S	0	0	0
			2344	1452	410	477	5			
1	D	338	Total	C	N	O	S	0	0	0
			2347	1452	408	482	5			
1	E	339	Total	C	N	O	S	0	0	0
			2368	1468	411	484	5			
1	F	338	Total	C	N	O	S	0	0	0
			2370	1473	412	480	5			
1	G	339	Total	C	N	O	S	0	0	0
			2340	1448	409	478	5			
1	H	338	Total	C	N	O	S	0	0	0
			2359	1459	411	484	5			
1	I	339	Total	C	N	O	S	0	0	0
			2366	1467	411	483	5			
1	J	338	Total	C	N	O	S	0	0	0
			2371	1472	412	482	5			
1	K	339	Total	C	N	O	S	0	0	0
			2336	1446	408	477	5			
1	L	338	Total	C	N	O	S	0	0	0
			2344	1451	407	481	5			
1	M	339	Total	C	N	O	S	0	0	0
			2363	1464	411	483	5			
1	N	338	Total	C	N	O	S	0	0	0
			2364	1467	412	480	5			
1	O	339	Total	C	N	O	S	0	0	0
			2331	1441	408	477	5			
1	P	338	Total	C	N	O	S	0	0	0
			2347	1452	408	482	5			

There are 288 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	340	LEU	-	expression tag	UNP P9WN82
A	341	VAL	-	expression tag	UNP P9WN82
A	342	PRO	-	expression tag	UNP P9WN82
A	343	ARG	-	expression tag	UNP P9WN82
A	344	GLY	-	expression tag	UNP P9WN82
A	345	SER	-	expression tag	UNP P9WN82
A	346	GLY	-	expression tag	UNP P9WN82
A	347	GLY	-	expression tag	UNP P9WN82
A	348	GLY	-	expression tag	UNP P9WN82
A	349	GLY	-	expression tag	UNP P9WN82
A	350	HIS	-	expression tag	UNP P9WN82
A	351	HIS	-	expression tag	UNP P9WN82
A	352	HIS	-	expression tag	UNP P9WN82
A	353	HIS	-	expression tag	UNP P9WN82
A	354	HIS	-	expression tag	UNP P9WN82
A	355	HIS	-	expression tag	UNP P9WN82
A	356	HIS	-	expression tag	UNP P9WN82
A	357	HIS	-	expression tag	UNP P9WN82
B	340	LEU	-	expression tag	UNP P9WN82
B	341	VAL	-	expression tag	UNP P9WN82
B	342	PRO	-	expression tag	UNP P9WN82
B	343	ARG	-	expression tag	UNP P9WN82
B	344	GLY	-	expression tag	UNP P9WN82
B	345	SER	-	expression tag	UNP P9WN82
B	346	GLY	-	expression tag	UNP P9WN82
B	347	GLY	-	expression tag	UNP P9WN82
B	348	GLY	-	expression tag	UNP P9WN82
B	349	GLY	-	expression tag	UNP P9WN82
B	350	HIS	-	expression tag	UNP P9WN82
B	351	HIS	-	expression tag	UNP P9WN82
B	352	HIS	-	expression tag	UNP P9WN82
B	353	HIS	-	expression tag	UNP P9WN82
B	354	HIS	-	expression tag	UNP P9WN82
B	355	HIS	-	expression tag	UNP P9WN82
B	356	HIS	-	expression tag	UNP P9WN82
B	357	HIS	-	expression tag	UNP P9WN82
C	340	LEU	-	expression tag	UNP P9WN82
C	341	VAL	-	expression tag	UNP P9WN82
C	342	PRO	-	expression tag	UNP P9WN82
C	343	ARG	-	expression tag	UNP P9WN82
C	344	GLY	-	expression tag	UNP P9WN82
C	345	SER	-	expression tag	UNP P9WN82

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Chain	Residue	Modelled	Actual	Comment	Reference
C	346	GLY	-	expression tag	UNP P9WN82
C	347	GLY	-	expression tag	UNP P9WN82
C	348	GLY	-	expression tag	UNP P9WN82
C	349	GLY	-	expression tag	UNP P9WN82
C	350	HIS	-	expression tag	UNP P9WN82
C	351	HIS	-	expression tag	UNP P9WN82
C	352	HIS	-	expression tag	UNP P9WN82
C	353	HIS	-	expression tag	UNP P9WN82
C	354	HIS	-	expression tag	UNP P9WN82
C	355	HIS	-	expression tag	UNP P9WN82
C	356	HIS	-	expression tag	UNP P9WN82
C	357	HIS	-	expression tag	UNP P9WN82
D	340	LEU	-	expression tag	UNP P9WN82
D	341	VAL	-	expression tag	UNP P9WN82
D	342	PRO	-	expression tag	UNP P9WN82
D	343	ARG	-	expression tag	UNP P9WN82
D	344	GLY	-	expression tag	UNP P9WN82
D	345	SER	-	expression tag	UNP P9WN82
D	346	GLY	-	expression tag	UNP P9WN82
D	347	GLY	-	expression tag	UNP P9WN82
D	348	GLY	-	expression tag	UNP P9WN82
D	349	GLY	-	expression tag	UNP P9WN82
D	350	HIS	-	expression tag	UNP P9WN82
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D	354	HIS	-	expression tag	UNP P9WN82
D	355	HIS	-	expression tag	UNP P9WN82
D	356	HIS	-	expression tag	UNP P9WN82
D	357	HIS	-	expression tag	UNP P9WN82
E	340	LEU	-	expression tag	UNP P9WN82
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E	342	PRO	-	expression tag	UNP P9WN82
E	343	ARG	-	expression tag	UNP P9WN82
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E	346	GLY	-	expression tag	UNP P9WN82
E	347	GLY	-	expression tag	UNP P9WN82
E	348	GLY	-	expression tag	UNP P9WN82
E	349	GLY	-	expression tag	UNP P9WN82
E	350	HIS	-	expression tag	UNP P9WN82
E	351	HIS	-	expression tag	UNP P9WN82

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Chain	Residue	Modelled	Actual	Comment	Reference
E	352	HIS	-	expression tag	UNP P9WN82
E	353	HIS	-	expression tag	UNP P9WN82
E	354	HIS	-	expression tag	UNP P9WN82
E	355	HIS	-	expression tag	UNP P9WN82
E	356	HIS	-	expression tag	UNP P9WN82
E	357	HIS	-	expression tag	UNP P9WN82
F	340	LEU	-	expression tag	UNP P9WN82
F	341	VAL	-	expression tag	UNP P9WN82
F	342	PRO	-	expression tag	UNP P9WN82
F	343	ARG	-	expression tag	UNP P9WN82
F	344	GLY	-	expression tag	UNP P9WN82
F	345	SER	-	expression tag	UNP P9WN82
F	346	GLY	-	expression tag	UNP P9WN82
F	347	GLY	-	expression tag	UNP P9WN82
F	348	GLY	-	expression tag	UNP P9WN82
F	349	GLY	-	expression tag	UNP P9WN82
F	350	HIS	-	expression tag	UNP P9WN82
F	351	HIS	-	expression tag	UNP P9WN82
F	352	HIS	-	expression tag	UNP P9WN82
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F	354	HIS	-	expression tag	UNP P9WN82
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F	356	HIS	-	expression tag	UNP P9WN82
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G	342	PRO	-	expression tag	UNP P9WN82
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G	344	GLY	-	expression tag	UNP P9WN82
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G	346	GLY	-	expression tag	UNP P9WN82
G	347	GLY	-	expression tag	UNP P9WN82
G	348	GLY	-	expression tag	UNP P9WN82
G	349	GLY	-	expression tag	UNP P9WN82
G	350	HIS	-	expression tag	UNP P9WN82
G	351	HIS	-	expression tag	UNP P9WN82
G	352	HIS	-	expression tag	UNP P9WN82
G	353	HIS	-	expression tag	UNP P9WN82
G	354	HIS	-	expression tag	UNP P9WN82
G	355	HIS	-	expression tag	UNP P9WN82
G	356	HIS	-	expression tag	UNP P9WN82
G	357	HIS	-	expression tag	UNP P9WN82

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Chain	Residue	Modelled	Actual	Comment	Reference
H	340	LEU	-	expression tag	UNP P9WN82
H	341	VAL	-	expression tag	UNP P9WN82
H	342	PRO	-	expression tag	UNP P9WN82
H	343	ARG	-	expression tag	UNP P9WN82
H	344	GLY	-	expression tag	UNP P9WN82
H	345	SER	-	expression tag	UNP P9WN82
H	346	GLY	-	expression tag	UNP P9WN82
H	347	GLY	-	expression tag	UNP P9WN82
H	348	GLY	-	expression tag	UNP P9WN82
H	349	GLY	-	expression tag	UNP P9WN82
H	350	HIS	-	expression tag	UNP P9WN82
H	351	HIS	-	expression tag	UNP P9WN82
H	352	HIS	-	expression tag	UNP P9WN82
H	353	HIS	-	expression tag	UNP P9WN82
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H	355	HIS	-	expression tag	UNP P9WN82
H	356	HIS	-	expression tag	UNP P9WN82
H	357	HIS	-	expression tag	UNP P9WN82
I	340	LEU	-	expression tag	UNP P9WN82
I	341	VAL	-	expression tag	UNP P9WN82
I	342	PRO	-	expression tag	UNP P9WN82
I	343	ARG	-	expression tag	UNP P9WN82
I	344	GLY	-	expression tag	UNP P9WN82
I	345	SER	-	expression tag	UNP P9WN82
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I	347	GLY	-	expression tag	UNP P9WN82
I	348	GLY	-	expression tag	UNP P9WN82
I	349	GLY	-	expression tag	UNP P9WN82
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I	351	HIS	-	expression tag	UNP P9WN82
I	352	HIS	-	expression tag	UNP P9WN82
I	353	HIS	-	expression tag	UNP P9WN82
I	354	HIS	-	expression tag	UNP P9WN82
I	355	HIS	-	expression tag	UNP P9WN82
I	356	HIS	-	expression tag	UNP P9WN82
I	357	HIS	-	expression tag	UNP P9WN82
J	340	LEU	-	expression tag	UNP P9WN82
J	341	VAL	-	expression tag	UNP P9WN82
J	342	PRO	-	expression tag	UNP P9WN82
J	343	ARG	-	expression tag	UNP P9WN82
J	344	GLY	-	expression tag	UNP P9WN82
J	345	SER	-	expression tag	UNP P9WN82

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Chain	Residue	Modelled	Actual	Comment	Reference
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J	348	GLY	-	expression tag	UNP P9WN82
J	349	GLY	-	expression tag	UNP P9WN82
J	350	HIS	-	expression tag	UNP P9WN82
J	351	HIS	-	expression tag	UNP P9WN82
J	352	HIS	-	expression tag	UNP P9WN82
J	353	HIS	-	expression tag	UNP P9WN82
J	354	HIS	-	expression tag	UNP P9WN82
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J	356	HIS	-	expression tag	UNP P9WN82
J	357	HIS	-	expression tag	UNP P9WN82
K	340	LEU	-	expression tag	UNP P9WN82
K	341	VAL	-	expression tag	UNP P9WN82
K	342	PRO	-	expression tag	UNP P9WN82
K	343	ARG	-	expression tag	UNP P9WN82
K	344	GLY	-	expression tag	UNP P9WN82
K	345	SER	-	expression tag	UNP P9WN82
K	346	GLY	-	expression tag	UNP P9WN82
K	347	GLY	-	expression tag	UNP P9WN82
K	348	GLY	-	expression tag	UNP P9WN82
K	349	GLY	-	expression tag	UNP P9WN82
K	350	HIS	-	expression tag	UNP P9WN82
K	351	HIS	-	expression tag	UNP P9WN82
K	352	HIS	-	expression tag	UNP P9WN82
K	353	HIS	-	expression tag	UNP P9WN82
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K	356	HIS	-	expression tag	UNP P9WN82
K	357	HIS	-	expression tag	UNP P9WN82
L	340	LEU	-	expression tag	UNP P9WN82
L	341	VAL	-	expression tag	UNP P9WN82
L	342	PRO	-	expression tag	UNP P9WN82
L	343	ARG	-	expression tag	UNP P9WN82
L	344	GLY	-	expression tag	UNP P9WN82
L	345	SER	-	expression tag	UNP P9WN82
L	346	GLY	-	expression tag	UNP P9WN82
L	347	GLY	-	expression tag	UNP P9WN82
L	348	GLY	-	expression tag	UNP P9WN82
L	349	GLY	-	expression tag	UNP P9WN82
L	350	HIS	-	expression tag	UNP P9WN82
L	351	HIS	-	expression tag	UNP P9WN82

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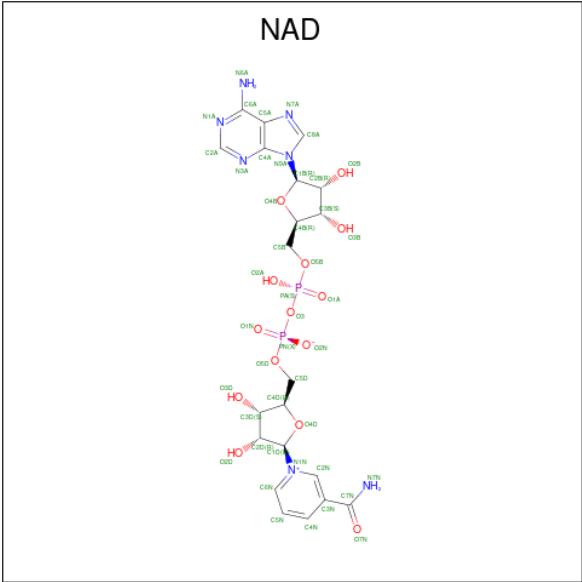
Chain	Residue	Modelled	Actual	Comment	Reference
L	352	HIS	-	expression tag	UNP P9WN82
L	353	HIS	-	expression tag	UNP P9WN82
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L	356	HIS	-	expression tag	UNP P9WN82
L	357	HIS	-	expression tag	UNP P9WN82
M	340	LEU	-	expression tag	UNP P9WN82
M	341	VAL	-	expression tag	UNP P9WN82
M	342	PRO	-	expression tag	UNP P9WN82
M	343	ARG	-	expression tag	UNP P9WN82
M	344	GLY	-	expression tag	UNP P9WN82
M	345	SER	-	expression tag	UNP P9WN82
M	346	GLY	-	expression tag	UNP P9WN82
M	347	GLY	-	expression tag	UNP P9WN82
M	348	GLY	-	expression tag	UNP P9WN82
M	349	GLY	-	expression tag	UNP P9WN82
M	350	HIS	-	expression tag	UNP P9WN82
M	351	HIS	-	expression tag	UNP P9WN82
M	352	HIS	-	expression tag	UNP P9WN82
M	353	HIS	-	expression tag	UNP P9WN82
M	354	HIS	-	expression tag	UNP P9WN82
M	355	HIS	-	expression tag	UNP P9WN82
M	356	HIS	-	expression tag	UNP P9WN82
M	357	HIS	-	expression tag	UNP P9WN82
N	340	LEU	-	expression tag	UNP P9WN82
N	341	VAL	-	expression tag	UNP P9WN82
N	342	PRO	-	expression tag	UNP P9WN82
N	343	ARG	-	expression tag	UNP P9WN82
N	344	GLY	-	expression tag	UNP P9WN82
N	345	SER	-	expression tag	UNP P9WN82
N	346	GLY	-	expression tag	UNP P9WN82
N	347	GLY	-	expression tag	UNP P9WN82
N	348	GLY	-	expression tag	UNP P9WN82
N	349	GLY	-	expression tag	UNP P9WN82
N	350	HIS	-	expression tag	UNP P9WN82
N	351	HIS	-	expression tag	UNP P9WN82
N	352	HIS	-	expression tag	UNP P9WN82
N	353	HIS	-	expression tag	UNP P9WN82
N	354	HIS	-	expression tag	UNP P9WN82
N	355	HIS	-	expression tag	UNP P9WN82
N	356	HIS	-	expression tag	UNP P9WN82
N	357	HIS	-	expression tag	UNP P9WN82

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Chain	Residue	Modelled	Actual	Comment	Reference
O	340	LEU	-	expression tag	UNP P9WN82
O	341	VAL	-	expression tag	UNP P9WN82
O	342	PRO	-	expression tag	UNP P9WN82
O	343	ARG	-	expression tag	UNP P9WN82
O	344	GLY	-	expression tag	UNP P9WN82
O	345	SER	-	expression tag	UNP P9WN82
O	346	GLY	-	expression tag	UNP P9WN82
O	347	GLY	-	expression tag	UNP P9WN82
O	348	GLY	-	expression tag	UNP P9WN82
O	349	GLY	-	expression tag	UNP P9WN82
O	350	HIS	-	expression tag	UNP P9WN82
O	351	HIS	-	expression tag	UNP P9WN82
O	352	HIS	-	expression tag	UNP P9WN82
O	353	HIS	-	expression tag	UNP P9WN82
O	354	HIS	-	expression tag	UNP P9WN82
O	355	HIS	-	expression tag	UNP P9WN82
O	356	HIS	-	expression tag	UNP P9WN82
O	357	HIS	-	expression tag	UNP P9WN82
P	340	LEU	-	expression tag	UNP P9WN82
P	341	VAL	-	expression tag	UNP P9WN82
P	342	PRO	-	expression tag	UNP P9WN82
P	343	ARG	-	expression tag	UNP P9WN82
P	344	GLY	-	expression tag	UNP P9WN82
P	345	SER	-	expression tag	UNP P9WN82
P	346	GLY	-	expression tag	UNP P9WN82
P	347	GLY	-	expression tag	UNP P9WN82
P	348	GLY	-	expression tag	UNP P9WN82
P	349	GLY	-	expression tag	UNP P9WN82
P	350	HIS	-	expression tag	UNP P9WN82
P	351	HIS	-	expression tag	UNP P9WN82
P	352	HIS	-	expression tag	UNP P9WN82
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P	354	HIS	-	expression tag	UNP P9WN82
P	355	HIS	-	expression tag	UNP P9WN82
P	356	HIS	-	expression tag	UNP P9WN82
P	357	HIS	-	expression tag	UNP P9WN82

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).



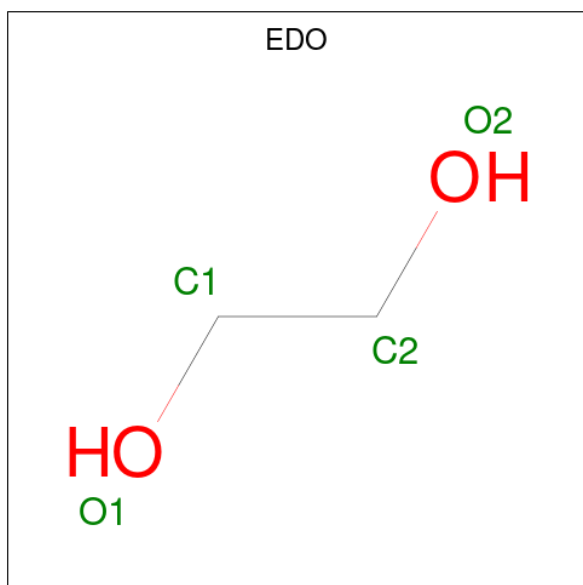
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	H	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	I	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	J	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	K	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	L	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	M	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	N	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).

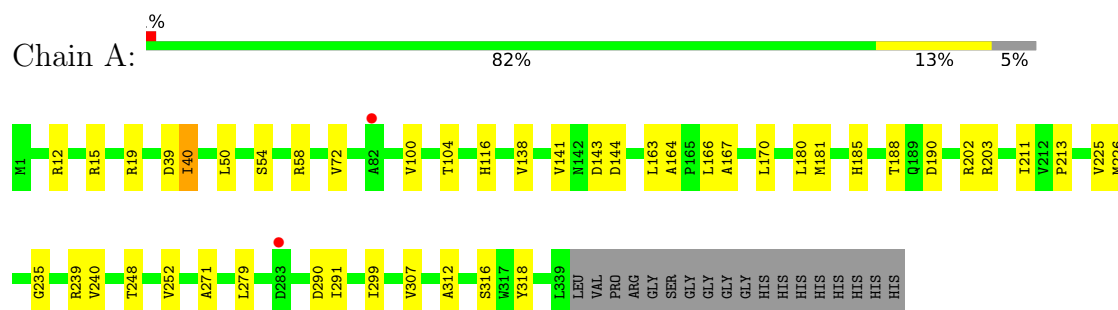


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

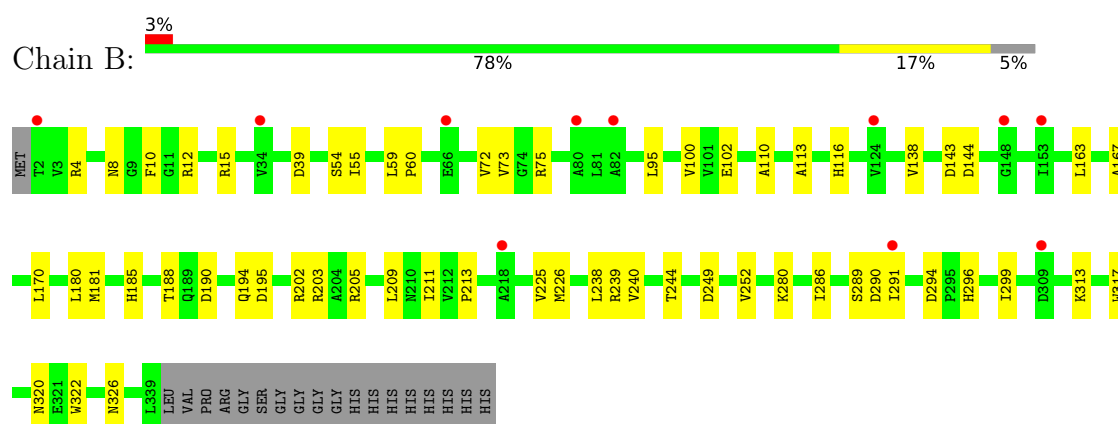
3 Residue-property plots [i](#)

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.

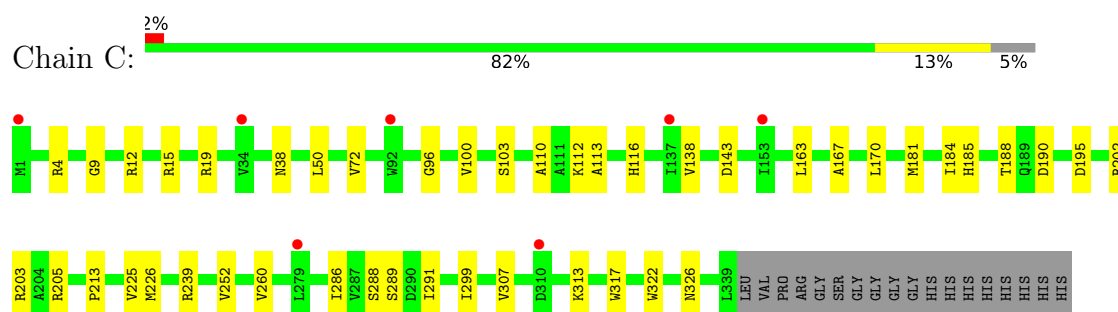
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



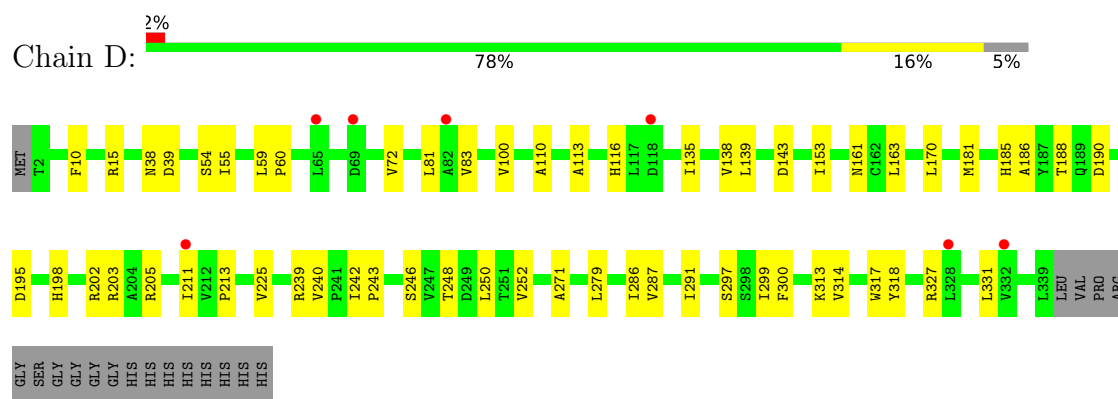
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



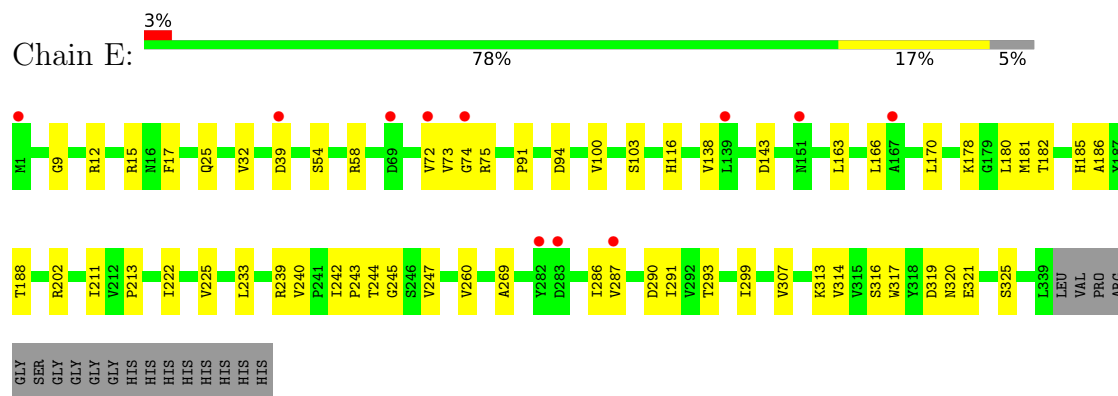
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



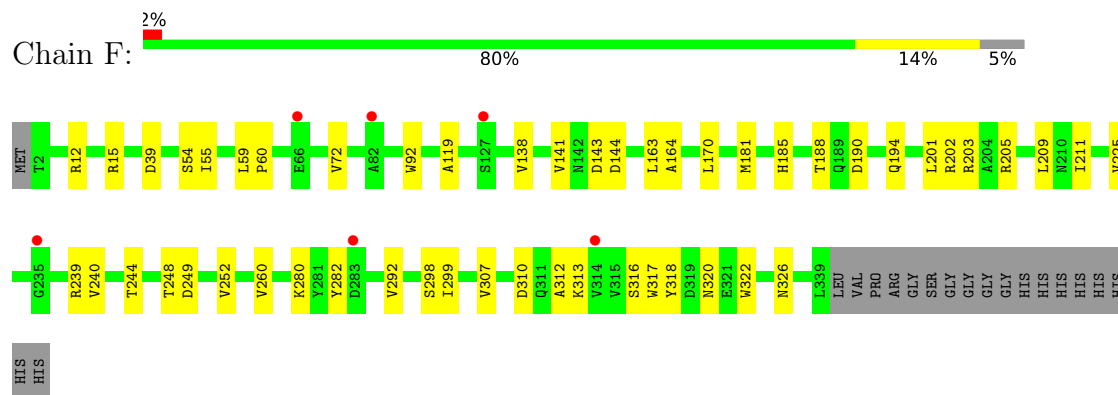
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



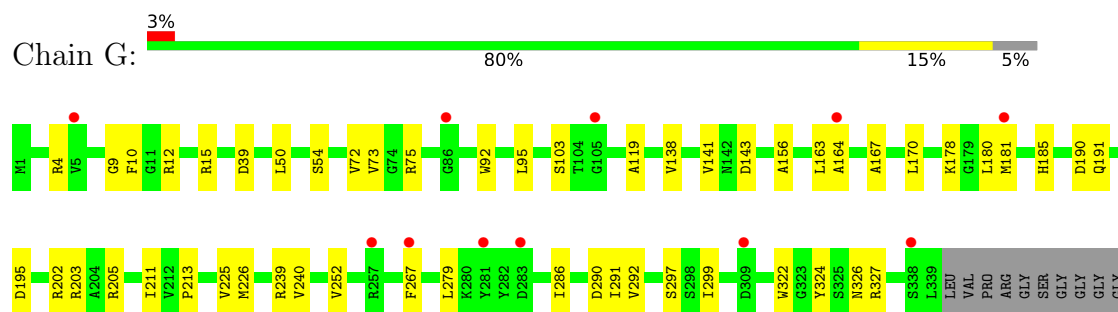
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



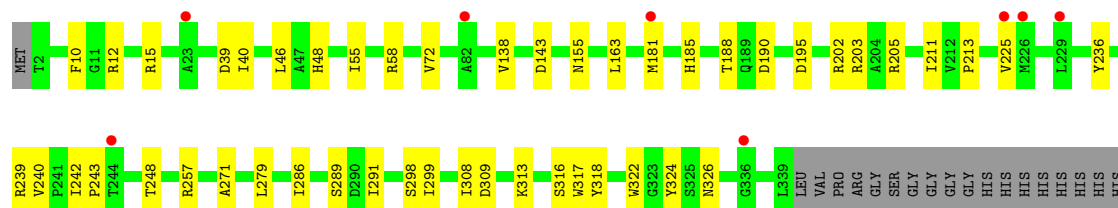
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



HIS
HIS
HIS
HIS
HIS
HIS
HIS

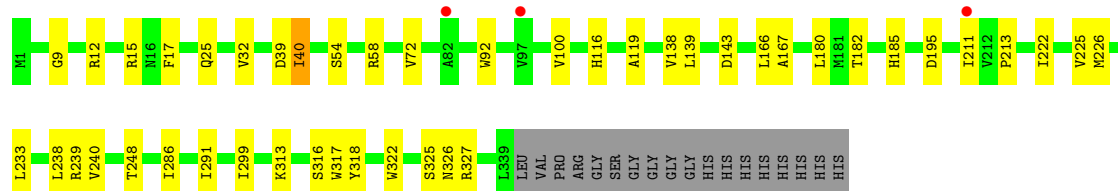
• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain H: 



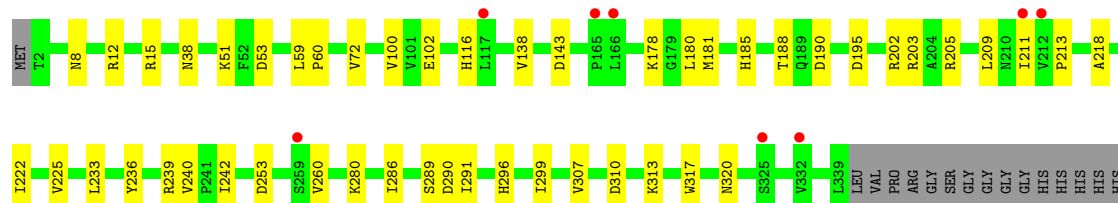
• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain I: 



• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

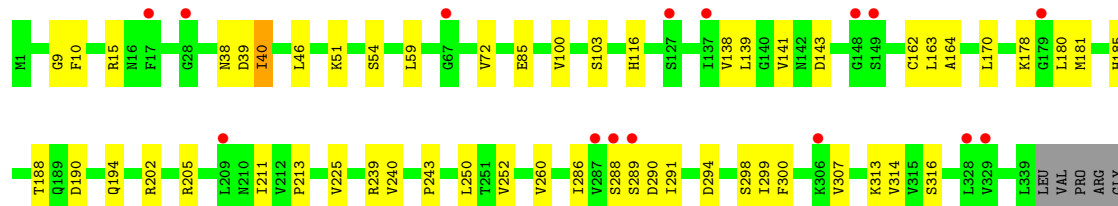
Chain J: 



HIS
HIS
HIS

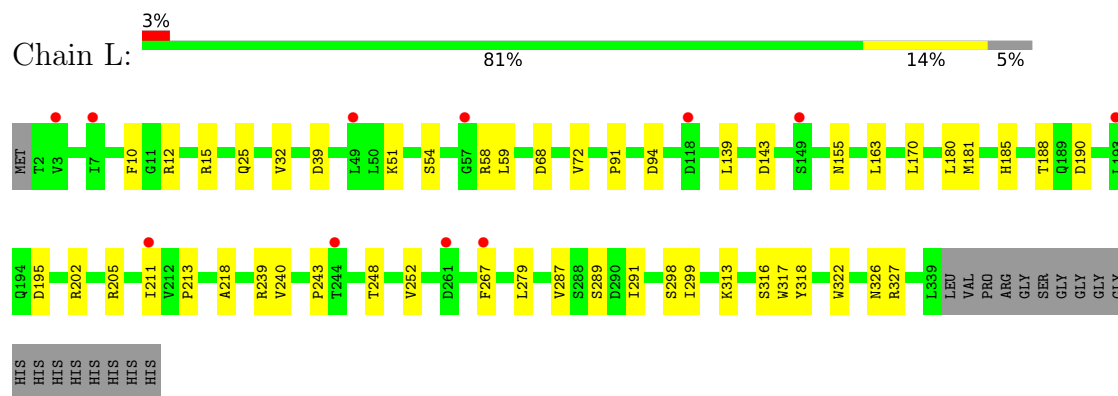
• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

Chain K: 

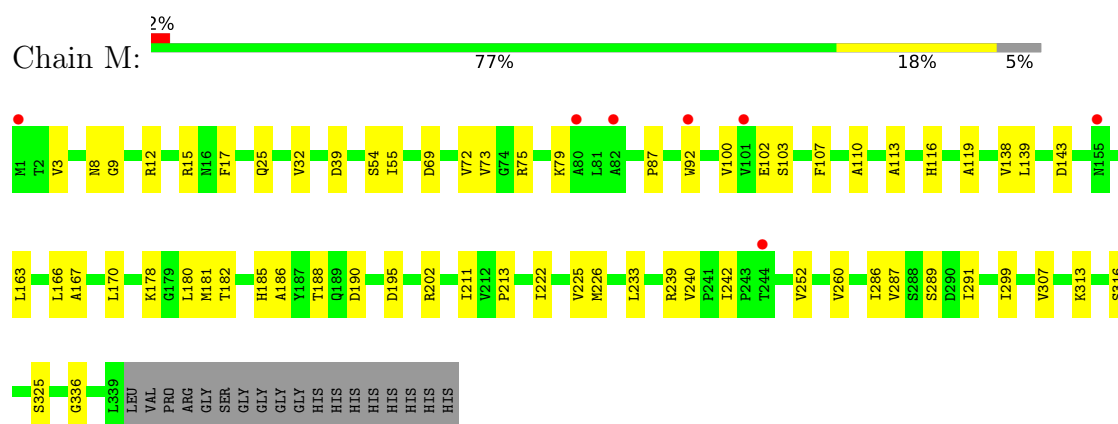


SER
GLY
GLY
GLY
GLY
HIS
HIS
HIS
HIS
HIS
HIS

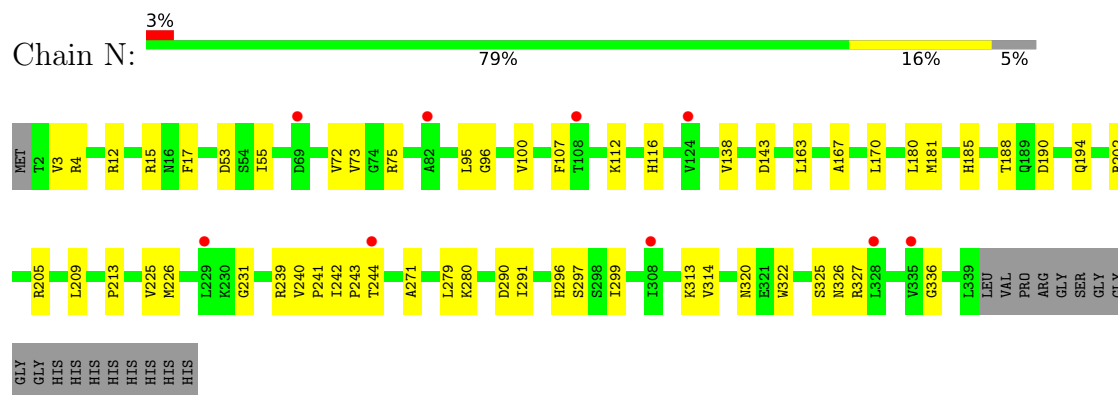
• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



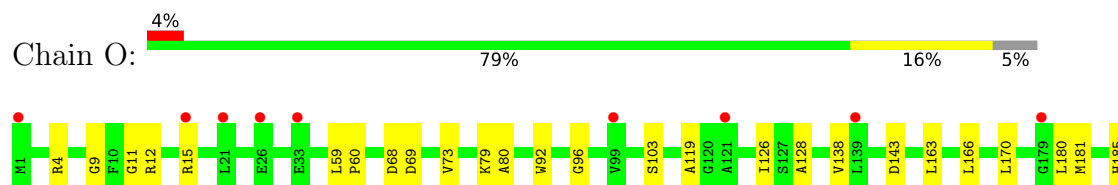
• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

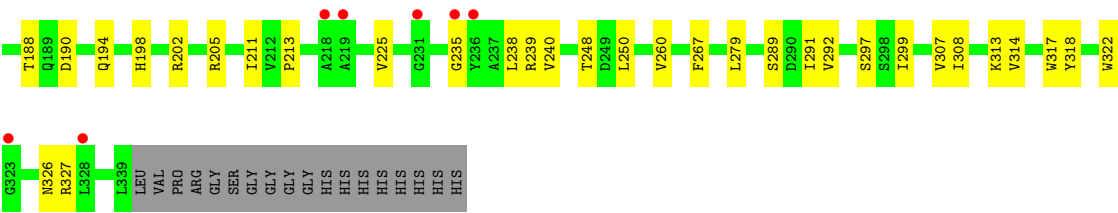


• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase

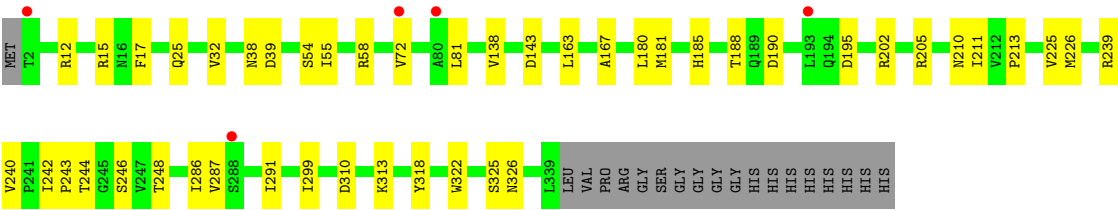
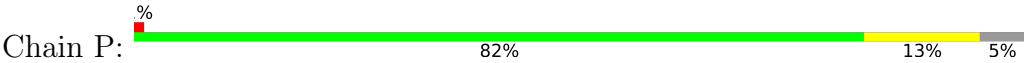


• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase





● Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	140.89Å 139.17Å 141.48Å 90.00° 108.52° 90.00°	Depositor
Resolution (Å)	47.15 – 3.29 47.15 – 3.29	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.15-3.29) 99.6 (47.15-3.29)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.255 , 0.295 0.254 , 0.295	Depositor DCC
R_{free} test set	3966 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	63.3	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 66.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	38389	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.13% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAD, EDO

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	0.07	0/2400	0.21	0/3289
1	B	0.07	0/2401	0.22	0/3290
1	C	0.08	0/2381	0.23	0/3263
1	D	0.08	0/2384	0.22	0/3268
1	E	0.09	0/2405	0.26	0/3296
1	F	0.07	0/2407	0.22	0/3298
1	G	0.07	0/2377	0.22	0/3260
1	H	0.09	0/2396	0.22	0/3283
1	I	0.07	0/2403	0.22	0/3293
1	J	0.07	0/2408	0.22	0/3299
1	K	0.07	0/2373	0.22	0/3255
1	L	0.08	0/2381	0.23	0/3264
1	M	0.08	0/2400	0.23	0/3289
1	N	0.07	0/2401	0.23	0/3290
1	O	0.07	0/2368	0.22	0/3248
1	P	0.07	0/2384	0.22	0/3268
All	All	0.07	0/38269	0.23	0/52453

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2363	0	2133	34	0
1	B	2364	0	2156	42	0
1	C	2344	0	2105	30	0
1	D	2347	0	2101	42	0
1	E	2368	0	2147	41	0
1	F	2370	0	2174	37	0
1	G	2340	0	2089	32	0
1	H	2359	0	2123	38	0
1	I	2366	0	2142	35	0
1	J	2371	0	2169	38	0
1	K	2336	0	2083	41	0
1	L	2344	0	2097	39	0
1	M	2363	0	2133	46	0
1	N	2364	0	2156	38	0
1	O	2331	0	2067	42	0
1	P	2347	0	2101	39	0
2	A	44	0	26	4	0
2	B	44	0	26	2	0
2	C	44	0	26	1	0
2	D	44	0	26	3	0
2	E	44	0	26	3	0
2	F	44	0	26	2	0
2	G	44	0	26	0	0
2	H	44	0	26	0	0
2	I	44	0	26	5	0
2	J	44	0	26	2	0
2	K	44	0	26	2	0
2	L	44	0	26	2	0
2	M	44	0	26	2	0
2	N	44	0	26	1	0
2	O	44	0	26	2	0
2	P	44	0	26	2	0
3	A	4	0	6	0	0
3	D	4	0	6	0	0
All	All	38389	0	34404	513	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 7.

All (513) 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:N:209:LEU:HD21	1:P:243:PRO:HG3	1.57	0.86
1:H:291:ILE:HG23	1:H:299:ILE:HD12	1.70	0.73
1:I:291:ILE:HG23	1:I:299:ILE:HD12	1.72	0.71
1:J:310:ASP:OD1	1:K:178:LYS:NZ	2.23	0.71
1:E:290:ASP:OD1	1:H:205:ARG:NH2	2.24	0.71
1:O:185:HIS:HB3	1:O:239:ARG:HD3	1.73	0.70
1:G:297:SER:HG	1:G:327:ARG:HH11	1.37	0.69
1:D:38:ASN:HA	1:D:81:LEU:HA	1.75	0.66
1:E:243:PRO:O	1:E:244:THR:HG22	1.94	0.66
1:N:290:ASP:OD1	1:O:205:ARG:NH2	2.28	0.66
1:B:290:ASP:OD1	1:C:205:ARG:NH2	2.28	0.66
1:F:299:ILE:HB	1:F:317:TRP:HB2	1.78	0.65
1:L:291:ILE:HG23	1:L:299:ILE:HD12	1.78	0.65
1:A:290:ASP:OD1	1:D:205:ARG:NH2	2.30	0.64
1:O:211:ILE:HG22	1:O:240:VAL:HA	1.78	0.64
1:J:222:ILE:HG21	1:J:233:LEU:HD12	1.79	0.64
1:E:182:THR:OG1	1:H:313:LYS:NZ	2.27	0.63
1:P:25:GLN:NE2	1:P:32:VAL:O	2.31	0.63
1:E:58:ARG:NH2	1:F:292:VAL:O	2.32	0.63
1:M:291:ILE:HG23	1:M:299:ILE:HD12	1.79	0.63
1:N:55:ILE:HD13	1:N:244:THR:HG23	1.80	0.63
1:D:291:ILE:HG23	1:D:299:ILE:HD12	1.80	0.62
1:L:185:HIS:HB3	1:L:239:ARG:HD3	1.81	0.62
1:I:182:THR:OG1	1:L:313:LYS:NZ	2.27	0.61
1:E:185:HIS:HB3	1:E:239:ARG:HD3	1.83	0.61
1:K:185:HIS:HB3	1:K:239:ARG:HD3	1.81	0.61
1:M:182:THR:OG1	1:P:313:LYS:NZ	2.29	0.61
1:J:260:VAL:HG23	1:J:307:VAL:HG12	1.81	0.61
1:I:185:HIS:HB3	1:I:239:ARG:HD3	1.83	0.60
1:O:12:ARG:HH11	1:O:15:ARG:HH21	1.49	0.60
1:J:290:ASP:OD1	1:K:205:ARG:NH2	2.28	0.60
1:O:292:VAL:O	1:P:58:ARG:NH2	2.33	0.60
1:I:54:SER:HA	1:J:289:SER:HB3	1.84	0.60
1:L:139:LEU:HD12	1:L:327:ARG:HD3	1.84	0.60
1:I:286:ILE:O	1:L:202:ARG:NH1	2.35	0.60
1:J:205:ARG:NH2	1:K:290:ASP:OD1	2.35	0.60
1:L:10:PHE:N	1:L:39:ASP:OD1	2.35	0.60
1:G:291:ILE:HG23	1:G:299:ILE:HD12	1.82	0.59
1:O:297:SER:OG	1:O:327:ARG:NH1	2.28	0.59
1:D:185:HIS:HB3	1:D:239:ARG:HD3	1.83	0.59
1:E:54:SER:OG	1:G:205:ARG:NH1	2.36	0.59
1:A:202:ARG:NH1	1:D:286:ILE:O	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:10:PHE:N	1:G:39:ASP:OD1	2.36	0.59
1:E:163:LEU:HD22	1:E:181:MET:HE3	1.85	0.59
1:F:320:ASN:HD22	2:F:401:NAD:H72N	1.50	0.59
1:H:185:HIS:HB3	1:H:239:ARG:HD3	1.84	0.58
1:C:185:HIS:HB3	1:C:239:ARG:HD3	1.85	0.58
1:D:135:ILE:HG23	1:D:153:ILE:HG23	1.86	0.58
1:O:12:ARG:N	2:O:401:NAD:O2N	2.37	0.58
1:M:8:ASN:ND2	1:M:102:GLU:OE1	2.36	0.58
1:M:178:LYS:NZ	1:P:310:ASP:OD1	2.32	0.58
1:F:163:LEU:HD22	1:F:181:MET:HE3	1.86	0.58
1:A:12:ARG:HH11	1:A:15:ARG:HH21	1.51	0.58
1:A:291:ILE:HG23	1:A:299:ILE:HD12	1.86	0.58
1:M:163:LEU:HD22	1:M:181:MET:HE3	1.86	0.58
1:N:202:ARG:HD2	1:N:213:PRO:HG2	1.86	0.58
1:C:291:ILE:HG23	1:C:299:ILE:HD12	1.84	0.57
1:K:289:SER:HB3	1:L:54:SER:HA	1.85	0.57
1:F:209:LEU:HD11	1:H:243:PRO:HB3	1.87	0.57
1:P:163:LEU:HD22	1:P:181:MET:HE3	1.85	0.57
1:J:202:ARG:HD2	1:J:213:PRO:HG2	1.86	0.57
1:O:260:VAL:HG23	1:O:307:VAL:HG12	1.87	0.57
1:G:297:SER:OG	1:G:327:ARG:NH1	2.32	0.57
1:I:54:SER:OG	1:K:205:ARG:NH1	2.36	0.57
1:E:222:ILE:HG21	1:E:233:LEU:HD12	1.87	0.57
1:J:185:HIS:HB3	1:J:239:ARG:HD3	1.87	0.57
1:B:194:GLN:HE22	1:D:55:ILE:HD11	1.69	0.57
1:M:195:ASP:O	1:O:15:ARG:NH2	2.30	0.57
1:N:15:ARG:NH2	1:P:195:ASP:O	2.35	0.57
1:J:320:ASN:HD22	2:J:401:NAD:H72N	1.52	0.56
1:M:39:ASP:OD1	2:M:401:NAD:H1B	2.04	0.56
1:P:291:ILE:HG23	1:P:299:ILE:HD12	1.87	0.56
1:E:138:VAL:HG23	1:E:225:VAL:HG11	1.86	0.56
1:K:39:ASP:OD2	2:K:401:NAD:O3B	2.22	0.56
1:N:194:GLN:HE22	1:P:55:ILE:HD11	1.69	0.56
1:O:291:ILE:HG23	1:O:299:ILE:HD12	1.86	0.56
1:C:195:ASP:OD2	1:C:205:ARG:NH1	2.39	0.56
1:J:195:ASP:O	1:L:15:ARG:NH2	2.33	0.56
1:M:25:GLN:NE2	1:M:32:VAL:O	2.39	0.56
1:O:181:MET:HG3	1:O:235:GLY:HA3	1.87	0.56
1:L:322:TRP:O	1:L:326:ASN:ND2	2.35	0.56
1:M:188:THR:OG1	1:M:190:ASP:OD1	2.21	0.56
1:A:211:ILE:HG22	1:A:240:VAL:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:ARG:HH11	1:C:15:ARG:HH21	1.54	0.56
1:K:39:ASP:OD1	1:K:40:ILE:N	2.36	0.56
1:L:202:ARG:HD2	1:L:213:PRO:HG2	1.88	0.56
1:I:138:VAL:HG23	1:I:225:VAL:HG11	1.87	0.56
1:J:209:LEU:HD11	1:L:243:PRO:HB3	1.86	0.56
1:I:39:ASP:OD1	1:I:40:ILE:N	2.39	0.55
1:J:53:ASP:C	1:L:205:ARG:HH22	2.14	0.55
1:J:286:ILE:O	1:K:202:ARG:NH1	2.39	0.55
1:M:286:ILE:O	1:P:202:ARG:NH1	2.38	0.55
1:F:185:HIS:HB3	1:F:239:ARG:HD3	1.87	0.55
1:F:55:ILE:HD13	1:F:244:THR:HG23	1.88	0.55
1:J:188:THR:OG1	1:J:190:ASP:OD1	2.23	0.55
1:B:138:VAL:HG23	1:B:225:VAL:HG11	1.89	0.55
1:P:202:ARG:HD2	1:P:213:PRO:HG2	1.89	0.55
1:N:185:HIS:HB3	1:N:239:ARG:HD3	1.89	0.55
1:O:188:THR:OG1	1:O:190:ASP:OD1	2.22	0.55
1:C:190:ASP:OD2	1:C:203:ARG:NH1	2.40	0.55
1:L:299:ILE:HB	1:L:317:TRP:HB2	1.88	0.55
1:B:291:ILE:HG23	1:B:299:ILE:HD12	1.89	0.54
1:K:163:LEU:HD22	1:K:181:MET:HE3	1.89	0.54
1:N:4:ARG:NH1	1:N:95:LEU:O	2.41	0.54
1:F:194:GLN:HE22	1:H:55:ILE:HD11	1.72	0.54
1:N:180:LEU:HD21	1:O:308:ILE:HG13	1.88	0.54
1:M:211:ILE:HG22	1:M:240:VAL:HA	1.90	0.54
1:O:138:VAL:HG23	1:O:225:VAL:HG11	1.89	0.54
1:B:54:SER:OG	1:D:205:ARG:NH1	2.41	0.54
1:C:138:VAL:HG23	1:C:225:VAL:HG11	1.89	0.54
1:D:188:THR:OG1	1:D:190:ASP:OD1	2.19	0.54
1:H:211:ILE:HG22	1:H:240:VAL:HA	1.88	0.54
1:J:202:ARG:NH1	1:K:286:ILE:O	2.38	0.54
1:K:211:ILE:HG22	1:K:240:VAL:HA	1.89	0.53
1:M:213:PRO:HD2	1:P:287:VAL:HG12	1.90	0.53
1:B:249:ASP:OD1	1:B:313:LYS:NZ	2.37	0.53
1:O:170:LEU:HD11	1:O:314:VAL:HG21	1.90	0.53
1:P:188:THR:OG1	1:P:190:ASP:OD1	2.19	0.53
1:B:39:ASP:OD1	2:B:401:NAD:H1B	2.08	0.53
1:J:205:ARG:NH1	1:L:54:SER:OG	2.41	0.53
1:K:170:LEU:HD13	1:K:252:VAL:HG11	1.91	0.53
1:L:12:ARG:HH11	1:L:15:ARG:HH21	1.56	0.53
1:J:181:MET:HE1	1:J:218:ALA:HB3	1.89	0.53
1:L:211:ILE:HG22	1:L:240:VAL:HA	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:185:HIS:HB3	1:M:239:ARG:HD3	1.90	0.53
1:O:322:TRP:O	1:O:326:ASN:ND2	2.37	0.53
1:J:15:ARG:NH2	1:L:195:ASP:O	2.37	0.53
1:G:12:ARG:HH11	1:G:15:ARG:HH21	1.57	0.53
1:I:180:LEU:HB3	1:L:313:LYS:HE3	1.91	0.52
1:J:190:ASP:OD2	1:J:203:ARG:NH1	2.42	0.52
1:B:205:ARG:NH1	1:D:54:SER:OG	2.43	0.52
1:E:180:LEU:HB3	1:H:313:LYS:HE3	1.91	0.52
1:F:188:THR:OG1	1:F:190:ASP:OD1	2.23	0.52
1:F:205:ARG:NH2	1:G:290:ASP:OD1	2.42	0.52
1:M:9:GLY:HA2	2:M:401:NAD:H8A	1.91	0.52
1:A:213:PRO:HD2	1:D:287:VAL:HG12	1.91	0.52
1:G:185:HIS:HB3	1:G:239:ARG:HD3	1.91	0.51
1:I:39:ASP:HA	2:I:401:NAD:C8A	2.41	0.51
1:K:141:VAL:HG21	1:K:164:ALA:HB1	1.92	0.51
1:P:39:ASP:HA	2:P:401:NAD:H8A	1.91	0.51
1:E:15:ARG:NH2	1:G:195:ASP:O	2.37	0.51
1:E:202:ARG:HD2	1:E:213:PRO:HG2	1.91	0.51
1:H:299:ILE:HB	1:H:317:TRP:HB2	1.92	0.51
1:F:249:ASP:OD1	1:F:313:LYS:NZ	2.41	0.51
1:M:180:LEU:HB3	1:P:313:LYS:HE3	1.92	0.51
1:A:202:ARG:HD2	1:A:213:PRO:HG2	1.91	0.51
1:O:202:ARG:HD2	1:O:213:PRO:HG2	1.92	0.51
1:F:260:VAL:HG23	1:F:307:VAL:HG12	1.92	0.51
1:G:202:ARG:HD3	1:G:213:PRO:HG2	1.92	0.51
1:O:69:ASP:HA	1:O:79:LYS:H	1.76	0.51
1:B:4:ARG:NH1	1:B:95:LEU:O	2.44	0.51
1:B:211:ILE:HG22	1:B:240:VAL:HA	1.91	0.51
1:B:322:TRP:O	1:B:326:ASN:ND2	2.34	0.51
1:H:195:ASP:OD2	1:H:205:ARG:NH1	2.44	0.51
1:K:188:THR:OG1	1:K:190:ASP:OD1	2.22	0.51
1:A:39:ASP:HA	2:A:401:NAD:C8A	2.40	0.51
1:B:202:ARG:NH1	1:C:286:ILE:O	2.42	0.51
1:K:288:SER:O	1:K:291:ILE:HG22	2.11	0.51
1:I:139:LEU:HD12	1:I:327:ARG:HD3	1.93	0.51
1:A:54:SER:HA	1:B:289:SER:HB3	1.93	0.51
1:F:190:ASP:OD2	1:F:203:ARG:NH1	2.44	0.51
1:D:202:ARG:HD2	1:D:213:PRO:HG2	1.93	0.50
1:E:25:GLN:NE2	1:E:32:VAL:O	2.44	0.50
1:I:166:LEU:HD11	1:I:316:SER:HB3	1.93	0.50
1:K:163:LEU:HD11	1:K:250:LEU:HD22	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:260:VAL:HG23	1:M:307:VAL:HG12	1.94	0.50
1:A:166:LEU:HD11	1:A:316:SER:HB3	1.94	0.50
1:D:161:ASN:O	1:D:297:SER:OG	2.26	0.50
1:N:205:ARG:NH1	1:P:54:SER:OG	2.44	0.50
1:A:138:VAL:HG23	1:A:225:VAL:HG11	1.93	0.50
1:B:286:ILE:O	1:C:202:ARG:NH1	2.44	0.50
1:G:170:LEU:HD13	1:G:252:VAL:HG11	1.93	0.50
1:I:195:ASP:O	1:K:15:ARG:NH2	2.36	0.50
1:G:54:SER:HA	1:H:289:SER:HB3	1.93	0.50
1:J:291:ILE:HD11	1:J:299:ILE:HG21	1.93	0.50
1:E:260:VAL:HG23	1:E:307:VAL:HG12	1.94	0.50
1:M:9:GLY:O	1:M:103:SER:OG	2.30	0.50
1:E:180:LEU:HD21	1:H:308:ILE:HG12	1.94	0.50
1:O:299:ILE:HB	1:O:317:TRP:HB2	1.94	0.50
1:A:185:HIS:HB3	1:A:239:ARG:HD3	1.94	0.50
1:B:167:ALA:HB3	1:B:226:MET:HE1	1.93	0.50
1:I:12:ARG:HH11	1:I:15:ARG:HH21	1.58	0.50
1:M:12:ARG:HH11	1:M:15:ARG:HH21	1.58	0.50
1:E:91:PRO:HB2	1:E:94:ASP:HB2	1.94	0.50
1:F:298:SER:OG	1:F:316:SER:OG	2.22	0.50
1:N:209:LEU:CD2	1:P:243:PRO:HG3	2.36	0.50
1:P:167:ALA:HB3	1:P:226:MET:HE1	1.94	0.50
1:B:15:ARG:NH2	1:D:195:ASP:O	2.38	0.49
1:F:298:SER:HG	1:F:316:SER:HG	1.51	0.49
1:E:286:ILE:O	1:H:202:ARG:NH1	2.42	0.49
1:G:292:VAL:O	1:H:58:ARG:NH2	2.44	0.49
1:P:322:TRP:O	1:P:326:ASN:ND2	2.38	0.49
1:D:10:PHE:N	1:D:39:ASP:OD1	2.46	0.49
1:I:167:ALA:HB3	1:I:226:MET:HE1	1.95	0.49
1:A:12:ARG:HB2	2:A:401:NAD:O1N	2.13	0.49
1:A:180:LEU:HB3	1:D:313:LYS:HE2	1.93	0.49
1:B:320:ASN:HD22	2:B:401:NAD:H72N	1.61	0.49
1:G:141:VAL:HG21	1:G:164:ALA:HB1	1.93	0.49
1:M:170:LEU:HD13	1:M:252:VAL:HG11	1.94	0.49
1:G:138:VAL:HG23	1:G:225:VAL:HG11	1.93	0.49
1:M:202:ARG:HD2	1:M:213:PRO:HG2	1.94	0.49
1:K:291:ILE:HD11	1:K:299:ILE:HD12	1.95	0.49
1:L:267:PHE:HD1	1:L:279:LEU:HD21	1.77	0.49
1:M:138:VAL:HG23	1:M:225:VAL:HG11	1.93	0.49
1:N:297:SER:OG	1:N:327:ARG:NH1	2.34	0.49
1:B:12:ARG:HH11	1:B:15:ARG:HH21	1.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:ILE:HG22	1:D:240:VAL:HA	1.95	0.49
1:J:178:LYS:NZ	1:J:253:ASP:OD2	2.46	0.49
1:D:271:ALA:HB2	1:D:279:LEU:HD22	1.94	0.48
1:J:180:LEU:HB3	1:K:313:LYS:HE2	1.94	0.48
1:P:39:ASP:HA	2:P:401:NAD:C8A	2.43	0.48
1:G:73:VAL:C	1:G:75:ARG:H	2.21	0.48
1:J:8:ASN:ND2	1:J:102:GLU:OE1	2.43	0.48
1:N:138:VAL:HG23	1:N:225:VAL:HG11	1.95	0.48
1:M:110:ALA:HA	1:M:113:ALA:HB3	1.94	0.48
1:D:299:ILE:HB	1:D:317:TRP:HB2	1.96	0.48
1:O:11:GLY:HA3	2:O:401:NAD:O5B	2.13	0.48
1:A:163:LEU:HD22	1:A:181:MET:HE3	1.95	0.48
1:B:8:ASN:ND2	1:B:102:GLU:OE1	2.45	0.48
1:N:53:ASP:C	1:P:205:ARG:HH22	2.22	0.48
1:F:138:VAL:HG23	1:F:225:VAL:HG11	1.95	0.48
1:L:51:LYS:HD2	1:L:59:LEU:HB3	1.94	0.48
1:A:19:ARG:NH2	1:A:50:LEU:O	2.47	0.48
1:K:9:GLY:O	1:K:103:SER:OG	2.31	0.48
1:A:39:ASP:HA	2:A:401:NAD:H8A	1.96	0.48
1:G:4:ARG:NH1	1:G:95:LEU:O	2.47	0.48
1:G:163:LEU:HD22	1:G:181:MET:SD	2.54	0.48
1:K:162:CYS:HA	1:K:298:SER:HB2	1.95	0.48
1:K:138:VAL:HG23	1:K:225:VAL:HG11	1.96	0.47
1:L:163:LEU:HD22	1:L:181:MET:SD	2.54	0.47
1:L:188:THR:OG1	1:L:190:ASP:OD1	2.27	0.47
1:N:280:LYS:HB2	1:N:296:HIS:CD2	2.48	0.47
1:P:138:VAL:HG23	1:P:225:VAL:HG11	1.96	0.47
1:E:247:VAL:HB	1:E:317:TRP:CE3	2.49	0.47
1:H:257:ARG:HH21	1:N:231:GLY:HA3	1.80	0.47
1:N:73:VAL:C	1:N:75:ARG:H	2.23	0.47
1:N:170:LEU:HD11	1:N:314:VAL:HG21	1.96	0.47
1:N:188:THR:OG1	1:N:239:ARG:NH1	2.48	0.47
1:G:191:GLN:OE1	1:G:239:ARG:NH1	2.46	0.47
1:I:39:ASP:HA	2:I:401:NAD:H8A	1.97	0.47
1:J:317:TRP:HZ2	1:K:213:PRO:HG3	1.80	0.47
1:B:163:LEU:HD22	1:B:181:MET:HE3	1.96	0.47
1:D:138:VAL:HG23	1:D:225:VAL:HG11	1.97	0.47
1:J:12:ARG:HH11	1:J:15:ARG:HH21	1.62	0.47
1:K:139:LEU:HA	1:K:143:ASP:HB3	1.97	0.47
1:M:166:LEU:HD11	1:M:316:SER:HB3	1.96	0.47
1:M:222:ILE:HG21	1:M:233:LEU:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:ARG:HD2	1:B:213:PRO:HG2	1.95	0.47
1:H:322:TRP:O	1:H:326:ASN:ND2	2.42	0.47
1:D:190:ASP:O	1:D:198:HIS:NE2	2.38	0.47
1:E:100:VAL:HG11	1:E:116:HIS:ND1	2.29	0.47
1:E:211:ILE:HG23	1:H:242:ILE:HD11	1.96	0.47
1:K:38:ASN:ND2	2:K:401:NAD:H2A	2.30	0.47
1:P:17:PHE:HA	1:P:325:SER:HB3	1.97	0.47
1:A:54:SER:N	1:C:205:ARG:HH12	2.13	0.47
1:B:185:HIS:HB3	1:B:239:ARG:HD3	1.96	0.47
1:B:195:ASP:O	1:D:15:ARG:NH2	2.39	0.47
1:E:166:LEU:HD11	1:E:316:SER:HB3	1.97	0.47
1:E:240:VAL:O	1:E:242:ILE:N	2.48	0.47
1:G:267:PHE:HD1	1:G:279:LEU:HD21	1.80	0.46
1:A:144:ASP:OD1	1:A:144:ASP:N	2.47	0.46
1:E:12:ARG:NH2	1:E:321:GLU:OE1	2.45	0.46
1:F:12:ARG:HH11	1:F:15:ARG:HH21	1.64	0.46
1:O:190:ASP:O	1:O:198:HIS:NE2	2.45	0.46
1:B:299:ILE:HB	1:B:317:TRP:HB2	1.97	0.46
1:D:163:LEU:HD11	1:D:250:LEU:HD22	1.97	0.46
1:N:12:ARG:HH11	1:N:15:ARG:HH21	1.64	0.46
1:P:185:HIS:HB3	1:P:239:ARG:HD3	1.98	0.46
1:C:163:LEU:HD22	1:C:181:MET:SD	2.56	0.46
1:E:73:VAL:C	1:E:75:ARG:H	2.23	0.46
1:C:188:THR:OG1	1:C:190:ASP:OD1	2.27	0.46
1:F:201:LEU:HB3	1:H:48:HIS:CE1	2.50	0.46
1:L:298:SER:OG	1:L:316:SER:OG	2.22	0.46
1:P:248:THR:HG23	1:P:318:TYR:CE1	2.50	0.46
1:H:138:VAL:HG23	1:H:225:VAL:HG11	1.97	0.46
1:J:100:VAL:HG11	1:J:116:HIS:ND1	2.30	0.46
1:N:167:ALA:HB3	1:N:226:MET:HE1	1.98	0.46
1:O:163:LEU:HD11	1:O:250:LEU:HD22	1.97	0.46
1:L:170:LEU:HD13	1:L:252:VAL:HG11	1.98	0.46
1:C:188:THR:OG1	1:C:239:ARG:NH1	2.46	0.46
1:G:15:ARG:HD2	1:G:50:LEU:HA	1.98	0.46
1:M:202:ARG:NH1	1:P:286:ILE:O	2.47	0.46
1:N:322:TRP:O	1:N:326:ASN:ND2	2.39	0.46
1:N:3:VAL:HG21	1:N:336:GLY:HA3	1.98	0.46
1:O:92:TRP:HD1	1:O:119:ALA:HB3	1.81	0.46
1:E:39:ASP:HA	2:E:401:NAD:C8A	2.46	0.45
1:E:188:THR:OG1	1:E:239:ARG:NH1	2.48	0.45
1:H:271:ALA:HB2	1:H:279:LEU:HD22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:167:ALA:HB3	1:M:226:MET:HE1	1.96	0.45
1:N:291:ILE:HD12	1:N:299:ILE:HD12	1.98	0.45
1:D:163:LEU:HD22	1:D:181:MET:HE3	1.99	0.45
1:E:211:ILE:HG22	1:E:240:VAL:HA	1.99	0.45
1:I:185:HIS:N	1:I:238:LEU:O	2.37	0.45
1:M:287:VAL:HG12	1:P:213:PRO:HD2	1.99	0.45
1:A:167:ALA:HB3	1:A:226:MET:HE1	1.98	0.45
1:I:12:ARG:HB2	2:I:401:NAD:O1N	2.17	0.45
1:J:138:VAL:HG23	1:J:225:VAL:HG11	1.98	0.45
1:A:39:ASP:OD1	1:A:40:ILE:N	2.48	0.45
1:D:170:LEU:HD13	1:D:252:VAL:HG11	1.99	0.45
1:F:202:ARG:NH1	1:G:286:ILE:O	2.48	0.45
1:I:25:GLN:NE2	1:I:32:VAL:O	2.49	0.45
1:J:211:ILE:HG22	1:J:240:VAL:HA	1.99	0.45
1:L:91:PRO:HB2	1:L:94:ASP:HB2	1.99	0.45
1:M:12:ARG:HG3	1:O:194:GLN:HG2	1.99	0.45
1:N:163:LEU:HD22	1:N:181:MET:SD	2.57	0.45
1:H:39:ASP:OD1	1:H:40:ILE:N	2.50	0.45
1:A:100:VAL:HG11	1:A:116:HIS:ND1	2.32	0.45
1:J:236:TYR:HE1	1:K:313:LYS:HD3	1.81	0.45
1:J:313:LYS:HE3	1:K:180:LEU:HB3	1.98	0.45
1:L:25:GLN:NE2	1:L:32:VAL:O	2.49	0.45
1:N:271:ALA:HB2	1:N:279:LEU:HD22	1.99	0.45
1:O:213:PRO:HB3	1:O:238:LEU:HD21	1.98	0.45
1:G:167:ALA:HB3	1:G:226:MET:HE1	1.99	0.45
1:K:260:VAL:HG23	1:K:307:VAL:HG12	1.99	0.45
1:L:39:ASP:HA	2:L:401:NAD:H8A	1.98	0.45
1:B:144:ASP:OD1	1:B:144:ASP:N	2.50	0.44
1:F:54:SER:N	1:H:205:ARG:HH12	2.14	0.44
1:F:280:LYS:HE3	1:F:282:TYR:HE2	1.82	0.44
1:B:280:LYS:HB2	1:B:296:HIS:ND1	2.33	0.44
1:D:100:VAL:HG11	1:D:116:HIS:ND1	2.32	0.44
1:F:322:TRP:O	1:F:326:ASN:ND2	2.38	0.44
1:N:17:PHE:HA	1:N:325:SER:HB3	1.98	0.44
1:G:211:ILE:HG23	1:G:240:VAL:HG12	2.00	0.44
1:A:170:LEU:HD13	1:A:252:VAL:HG11	1.99	0.44
1:D:190:ASP:OD2	1:D:203:ARG:NH1	2.51	0.44
1:F:170:LEU:HD13	1:F:252:VAL:HG11	2.00	0.44
1:M:69:ASP:HA	1:M:79:LYS:H	1.82	0.44
1:C:170:LEU:HD13	1:C:252:VAL:HG11	2.00	0.44
1:F:141:VAL:HG21	1:F:164:ALA:HB1	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:190:ASP:OD1	1:H:203:ARG:NH1	2.50	0.44
1:F:248:THR:HG23	1:F:318:TYR:CE1	2.53	0.44
1:I:143:ASP:N	1:I:143:ASP:OD1	2.51	0.44
1:I:322:TRP:O	1:I:326:ASN:ND2	2.39	0.44
1:N:240:VAL:O	1:N:242:ILE:N	2.50	0.44
1:K:298:SER:OG	1:K:316:SER:OG	2.30	0.44
1:O:289:SER:HB3	1:P:54:SER:HA	1.99	0.44
1:A:213:PRO:HG3	1:D:317:TRP:HZ2	1.83	0.44
1:B:73:VAL:C	1:B:75:ARG:H	2.26	0.44
1:E:320:ASN:O	2:E:401:NAD:H4N	2.18	0.44
1:A:188:THR:OG1	1:A:190:ASP:OD1	2.23	0.44
1:D:39:ASP:HA	2:D:401:NAD:C8A	2.48	0.44
1:D:110:ALA:HA	1:D:113:ALA:HB3	1.99	0.44
1:E:291:ILE:HG23	1:E:299:ILE:HD12	2.00	0.44
1:M:313:LYS:HE2	1:P:180:LEU:HB3	2.00	0.44
1:N:188:THR:OG1	1:N:190:ASP:OD1	2.27	0.44
1:O:4:ARG:HD3	1:O:96:GLY:O	2.18	0.44
1:O:267:PHE:HD1	1:O:279:LEU:HD21	1.81	0.44
1:A:143:ASP:OD1	1:A:143:ASP:N	2.51	0.43
1:C:167:ALA:HB3	1:C:226:MET:HE1	1.99	0.43
1:F:59:LEU:HD12	1:F:60:PRO:HD2	1.99	0.43
1:M:166:LEU:HD21	1:M:316:SER:HB2	1.99	0.43
1:B:100:VAL:HG11	1:B:116:HIS:ND1	2.33	0.43
1:C:4:ARG:HD3	1:C:96:GLY:O	2.18	0.43
1:F:143:ASP:OD1	1:F:143:ASP:N	2.51	0.43
1:I:213:PRO:HG3	1:L:317:TRP:HZ2	1.82	0.43
1:K:54:SER:HA	1:L:289:SER:HB3	2.00	0.43
1:B:170:LEU:HD13	1:B:252:VAL:HG11	2.00	0.43
1:D:246:SER:HB2	1:D:318:TYR:CZ	2.53	0.43
1:E:291:ILE:HD12	1:E:299:ILE:HD12	2.00	0.43
1:F:211:ILE:HG22	1:F:240:VAL:HG12	2.00	0.43
1:H:10:PHE:HD2	1:H:46:LEU:HD22	1.83	0.43
1:K:100:VAL:HG11	1:K:116:HIS:ND1	2.33	0.43
1:L:143:ASP:N	1:L:143:ASP:OD1	2.52	0.43
1:F:39:ASP:OD2	2:F:401:NAD:O3B	2.36	0.43
1:G:322:TRP:O	1:G:326:ASN:ND2	2.44	0.43
1:I:58:ARG:NE	1:J:290:ASP:O	2.52	0.43
1:N:242:ILE:HD11	1:O:211:ILE:HG23	2.01	0.43
1:E:170:LEU:HD11	1:E:314:VAL:HG21	1.99	0.43
1:F:211:ILE:HG22	1:F:240:VAL:HA	1.99	0.43
1:G:143:ASP:OD1	1:G:143:ASP:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:240:VAL:O	1:H:242:ILE:N	2.51	0.43
1:K:143:ASP:OD1	1:K:143:ASP:N	2.51	0.43
1:B:209:LEU:HD11	1:D:243:PRO:HB3	2.00	0.43
1:C:288:SER:HB3	1:C:317:TRP:CZ3	2.53	0.43
1:E:287:VAL:HG12	1:H:213:PRO:HD2	2.01	0.43
1:K:300:PHE:CE1	1:K:314:VAL:HG13	2.53	0.43
1:L:155:ASN:ND2	1:L:155:ASN:O	2.51	0.43
1:M:139:LEU:HA	1:M:143:ASP:HB3	2.00	0.43
1:P:143:ASP:OD1	1:P:143:ASP:N	2.51	0.43
1:C:9:GLY:O	1:C:103:SER:OG	2.37	0.43
1:F:310:ASP:H	1:G:178:LYS:HE3	1.83	0.43
1:I:166:LEU:HD21	1:I:316:SER:HB2	2.01	0.43
1:J:143:ASP:OD1	1:J:143:ASP:N	2.52	0.43
1:L:181:MET:HE1	1:L:218:ALA:HB3	2.01	0.43
1:O:248:THR:HG23	1:O:318:TYR:CE1	2.54	0.43
1:A:141:VAL:HG21	1:A:164:ALA:HB1	2.00	0.43
1:C:19:ARG:NH2	1:C:50:LEU:O	2.52	0.43
1:F:280:LYS:HB3	1:F:299:ILE:HD13	2.01	0.43
1:I:291:ILE:HG21	1:I:317:TRP:HB3	2.01	0.43
1:K:170:LEU:HD11	1:K:314:VAL:HG21	2.01	0.43
1:N:107:PHE:HD1	1:N:112:LYS:HD3	1.83	0.43
1:E:178:LYS:HD2	1:H:309:ASP:HB3	2.01	0.42
1:P:38:ASN:HA	1:P:81:LEU:HA	2.01	0.42
1:C:202:ARG:HD2	1:C:213:PRO:HG2	1.99	0.42
1:I:12:ARG:HG3	1:K:194:GLN:HG2	2.00	0.42
1:L:248:THR:HG23	1:L:318:TYR:CE1	2.54	0.42
1:D:297:SER:HB2	1:D:327:ARG:HH11	1.85	0.42
1:K:294:ASP:OD2	1:L:58:ARG:NH1	2.52	0.42
1:M:143:ASP:OD1	1:M:143:ASP:N	2.52	0.42
1:N:320:ASN:HD22	2:N:401:NAD:H72N	1.68	0.42
1:O:9:GLY:O	1:O:103:SER:OG	2.31	0.42
1:B:180:LEU:HB3	1:C:313:LYS:HE2	2.00	0.42
1:F:144:ASP:OD1	1:F:144:ASP:N	2.52	0.42
1:L:188:THR:OG1	1:L:239:ARG:NH1	2.52	0.42
1:O:143:ASP:OD1	1:O:143:ASP:N	2.52	0.42
1:B:291:ILE:HD12	1:B:299:ILE:HD12	2.01	0.42
1:C:143:ASP:OD1	1:C:143:ASP:N	2.52	0.42
1:G:92:TRP:HD1	1:G:119:ALA:HB3	1.84	0.42
1:G:190:ASP:OD1	1:G:203:ARG:NH1	2.52	0.42
1:H:12:ARG:HH11	1:H:15:ARG:HH21	1.65	0.42
1:J:38:ASN:ND2	2:J:401:NAD:H2A	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ARG:NH1	1:B:294:ASP:OD2	2.53	0.42
1:C:110:ALA:HA	1:C:113:ALA:HB3	2.02	0.42
1:H:143:ASP:OD1	1:H:143:ASP:N	2.53	0.42
1:I:313:LYS:HE2	1:L:180:LEU:HB3	2.01	0.42
1:K:10:PHE:HD2	1:K:46:LEU:HD22	1.85	0.42
1:O:163:LEU:HD22	1:O:181:MET:SD	2.60	0.42
1:P:12:ARG:HH11	1:P:15:ARG:HH21	1.68	0.42
1:G:9:GLY:O	1:G:103:SER:OG	2.34	0.42
1:H:163:LEU:HD22	1:H:181:MET:SD	2.60	0.42
1:L:39:ASP:HA	2:L:401:NAD:C8A	2.49	0.42
1:M:211:ILE:HG23	1:P:242:ILE:HD12	2.02	0.42
1:N:313:LYS:HE2	1:O:180:LEU:HB3	2.02	0.42
1:B:143:ASP:OD1	1:B:143:ASP:N	2.52	0.42
1:E:293:THR:N	1:E:319:ASP:OD2	2.50	0.42
1:H:248:THR:HG23	1:H:318:TYR:CE1	2.55	0.42
1:I:248:THR:HG23	1:I:318:TYR:CE1	2.54	0.42
1:P:291:ILE:HD12	1:P:299:ILE:HD12	2.02	0.42
1:A:190:ASP:OD2	1:A:203:ARG:NH1	2.53	0.42
1:B:110:ALA:HA	1:B:113:ALA:HB3	2.01	0.42
1:B:188:THR:OG1	1:B:239:ARG:NH1	2.53	0.42
1:I:17:PHE:HA	1:I:325:SER:HB3	2.01	0.42
1:K:85:GLU:OE1	1:K:85:GLU:N	2.53	0.42
1:K:202:ARG:HD2	1:K:213:PRO:HG2	2.02	0.42
1:N:180:LEU:HB3	1:O:313:LYS:HE2	2.01	0.42
2:E:401:NAD:H2N	2:E:401:NAD:H2D	1.73	0.41
1:O:166:LEU:HA	1:O:267:PHE:HE1	1.85	0.41
1:O:188:THR:OG1	1:O:239:ARG:NH1	2.43	0.41
1:B:55:ILE:HD13	1:B:244:THR:HG23	2.02	0.41
1:C:260:VAL:HG23	1:C:307:VAL:HG12	2.03	0.41
1:E:143:ASP:OD1	1:E:143:ASP:N	2.50	0.41
1:E:313:LYS:HD3	1:H:236:TYR:HE1	1.86	0.41
1:I:100:VAL:HG11	1:I:116:HIS:ND1	2.35	0.41
1:O:213:PRO:CA	1:O:238:LEU:HD23	2.50	0.41
1:B:59:LEU:HD12	1:B:60:PRO:HD2	2.02	0.41
1:C:100:VAL:HG11	1:C:116:HIS:ND1	2.36	0.41
1:J:280:LYS:HB2	1:J:296:HIS:CD2	2.55	0.41
1:K:51:LYS:NZ	1:K:59:LEU:O	2.41	0.41
1:O:68:ASP:OD1	1:O:68:ASP:N	2.52	0.41
1:C:112:LYS:HE2	1:E:269:ALA:HB2	2.02	0.41
1:C:289:SER:HB3	1:D:54:SER:HA	2.02	0.41
1:D:139:LEU:HD21	1:D:331:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:186:ALA:HB2	1:E:245:GLY:O	2.19	0.41
1:H:188:THR:OG1	1:H:239:ARG:NH1	2.51	0.41
1:I:92:TRP:HD1	1:I:119:ALA:HB3	1.84	0.41
1:J:51:LYS:NZ	1:J:59:LEU:O	2.44	0.41
1:M:211:ILE:HG23	1:P:242:ILE:CD1	2.50	0.41
1:A:307:VAL:HG23	1:A:312:ALA:HB2	2.01	0.41
1:C:38:ASN:ND2	2:C:401:NAD:H2A	2.36	0.41
1:H:257:ARG:NH2	1:N:231:GLY:HA3	2.35	0.41
1:J:240:VAL:O	1:J:242:ILE:N	2.54	0.41
1:M:289:SER:HB3	1:P:210:ASN:HB3	2.03	0.41
1:N:143:ASP:OD1	1:N:143:ASP:N	2.51	0.41
1:O:59:LEU:HD12	1:O:60:PRO:HD2	2.02	0.41
1:A:181:MET:HG3	1:A:235:GLY:HA3	2.02	0.41
1:D:186:ALA:HA	1:D:242:ILE:HG22	2.02	0.41
1:H:155:ASN:HB3	1:H:324:TYR:OH	2.20	0.41
1:J:280:LYS:HB3	1:J:299:ILE:HG23	2.03	0.41
1:C:322:TRP:O	1:C:326:ASN:ND2	2.46	0.41
1:E:17:PHE:HA	1:E:325:SER:HB3	2.02	0.41
1:I:9:GLY:HA2	2:I:401:NAD:H8A	2.02	0.41
1:M:3:VAL:HG21	1:M:336:GLY:HA3	2.03	0.41
1:P:240:VAL:O	1:P:242:ILE:N	2.52	0.41
1:D:39:ASP:HA	2:D:401:NAD:H8A	2.03	0.41
1:D:300:PHE:HE1	1:D:314:VAL:HG13	1.86	0.41
1:F:15:ARG:NH2	1:H:195:ASP:O	2.38	0.41
1:F:313:LYS:HE2	1:G:180:LEU:HB3	2.02	0.41
1:G:156:ALA:O	1:G:324:TYR:OH	2.37	0.41
1:I:211:ILE:HG22	1:I:240:VAL:HA	2.03	0.41
2:I:401:NAD:N3A	2:I:401:NAD:H2B	2.36	0.41
1:L:68:ASP:OD1	1:L:68:ASP:N	2.53	0.41
1:M:87:PRO:HD3	1:M:107:PHE:CZ	2.55	0.41
1:M:100:VAL:HG11	1:M:116:HIS:ND1	2.36	0.41
1:N:4:ARG:HD3	1:N:96:GLY:O	2.21	0.41
1:O:126:ILE:HG22	1:O:128:ALA:H	1.84	0.41
1:A:271:ALA:HB2	1:A:279:LEU:HD22	2.03	0.41
1:B:190:ASP:OD1	1:B:203:ARG:NH1	2.54	0.41
1:M:12:ARG:HH21	1:M:55:ILE:HD12	1.85	0.41
1:M:54:SER:OG	1:O:205:ARG:NH1	2.53	0.41
1:M:73:VAL:C	1:M:75:ARG:H	2.28	0.41
1:M:186:ALA:HB1	1:M:242:ILE:O	2.21	0.41
1:N:100:VAL:HG11	1:N:116:HIS:ND1	2.36	0.41
1:P:246:SER:HB2	1:P:318:TYR:CZ	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:THR:HA	2:A:401:NAD:H52A	2.02	0.40
1:A:248:THR:HG23	1:A:318:TYR:CE1	2.56	0.40
1:D:39:ASP:OD1	2:D:401:NAD:H1B	2.22	0.40
1:D:59:LEU:HD12	1:D:60:PRO:HD2	2.02	0.40
1:D:248:THR:HG23	1:D:318:TYR:CE1	2.56	0.40
1:F:307:VAL:HG23	1:F:312:ALA:HB2	2.02	0.40
1:B:10:PHE:CE2	1:B:15:ARG:HG2	2.57	0.40
1:M:92:TRP:HD1	1:M:119:ALA:HB3	1.85	0.40
1:D:143:ASP:OD1	1:D:143:ASP:N	2.52	0.40
1:F:92:TRP:HD1	1:F:119:ALA:HB3	1.85	0.40
1:I:213:PRO:HD2	1:L:287:VAL:HG12	2.03	0.40
1:M:17:PHE:HA	1:M:325:SER:HB3	2.04	0.40
1:B:238:LEU:HD13	1:C:184:ILE:HD12	2.02	0.40
1:E:9:GLY:O	1:E:103:SER:OG	2.39	0.40
1:H:298:SER:OG	1:H:316:SER:OG	2.34	0.40
1:M:240:VAL:HG21	1:P:211:ILE:HD11	2.02	0.40
1:B:280:LYS:HB2	1:B:296:HIS:CE1	2.56	0.40
1:E:202:ARG:NH1	1:H:286:ILE:O	2.53	0.40
1:I:222:ILE:HG21	1:I:233:LEU:HD12	2.04	0.40
1:J:59:LEU:HD12	1:J:60:PRO:HD2	2.03	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	337/357 (94%)	313 (93%)	22 (6%)	2 (1%)	21	52
1	B	336/357 (94%)	309 (92%)	26 (8%)	1 (0%)	36	65
1	C	337/357 (94%)	310 (92%)	26 (8%)	1 (0%)	36	65
1	D	336/357 (94%)	311 (93%)	23 (7%)	2 (1%)	21	52

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	337/357 (94%)	313 (93%)	22 (6%)	2 (1%)	21	52
1	F	336/357 (94%)	307 (91%)	28 (8%)	1 (0%)	36	65
1	G	337/357 (94%)	310 (92%)	26 (8%)	1 (0%)	36	65
1	H	336/357 (94%)	309 (92%)	26 (8%)	1 (0%)	36	65
1	I	337/357 (94%)	311 (92%)	24 (7%)	2 (1%)	21	52
1	J	336/357 (94%)	308 (92%)	27 (8%)	1 (0%)	36	65
1	K	337/357 (94%)	315 (94%)	19 (6%)	3 (1%)	14	43
1	L	336/357 (94%)	306 (91%)	29 (9%)	1 (0%)	36	65
1	M	337/357 (94%)	310 (92%)	26 (8%)	1 (0%)	36	65
1	N	336/357 (94%)	311 (93%)	22 (6%)	3 (1%)	14	43
1	O	337/357 (94%)	313 (93%)	22 (6%)	2 (1%)	21	52
1	P	336/357 (94%)	311 (93%)	23 (7%)	2 (1%)	21	52
All	All	5384/5712 (94%)	4967 (92%)	391 (7%)	26 (0%)	24	55

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	VAL
1	C	72	VAL
1	D	72	VAL
1	E	72	VAL
1	F	72	VAL
1	G	72	VAL
1	H	72	VAL
1	I	72	VAL
1	J	72	VAL
1	K	72	VAL
1	L	72	VAL
1	M	72	VAL
1	N	72	VAL
1	P	72	VAL
1	P	244	THR
1	O	73	VAL
1	B	72	VAL
1	K	243	PRO
1	O	80	ALA
1	I	40	ILE
1	N	243	PRO

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Mol	Chain	Res	Type
1	A	40	ILE
1	D	83	VAL
1	N	241	PRO
1	E	74	GLY
1	K	40	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	220/284 (78%)	220 (100%)	0	100	100
1	B	223/284 (78%)	223 (100%)	0	100	100
1	C	215/284 (76%)	215 (100%)	0	100	100
1	D	217/284 (76%)	217 (100%)	0	100	100
1	E	222/284 (78%)	222 (100%)	0	100	100
1	F	225/284 (79%)	225 (100%)	0	100	100
1	G	214/284 (75%)	214 (100%)	0	100	100
1	H	220/284 (78%)	220 (100%)	0	100	100
1	I	221/284 (78%)	221 (100%)	0	100	100
1	J	225/284 (79%)	225 (100%)	0	100	100
1	K	213/284 (75%)	213 (100%)	0	100	100
1	L	216/284 (76%)	216 (100%)	0	100	100
1	M	220/284 (78%)	220 (100%)	0	100	100
1	N	223/284 (78%)	223 (100%)	0	100	100
1	O	211/284 (74%)	211 (100%)	0	100	100
1	P	217/284 (76%)	217 (100%)	0	100	100
All	All	3502/4544 (77%)	3502 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	155	ASN
1	A	161	ASN
1	A	264	ASN
1	C	16	ASN
1	C	43	ASN
1	D	48	HIS
1	D	194	GLN
1	E	264	ASN
1	E	296	HIS
1	F	210	ASN
1	F	320	ASN
1	G	264	ASN
1	G	296	HIS
1	H	38	ASN
1	H	43	ASN
1	I	161	ASN
1	I	264	ASN
1	J	264	ASN
1	K	38	ASN
1	K	189	GLN
1	K	264	ASN
1	M	210	ASN
1	N	189	GLN
1	N	264	ASN
1	O	38	ASN
1	O	264	ASN
1	O	311	GLN
1	P	38	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

18 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	P	401	-	46,48,48	0.65	1 (2%)	64,73,73	0.68	0
2	NAD	M	401	-	46,48,48	0.64	1 (2%)	64,73,73	0.87	4 (6%)
2	NAD	L	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.69	1 (1%)
3	EDO	D	402	-	3,3,3	0.43	0	2,2,2	0.38	0
2	NAD	N	401	-	46,48,48	0.65	1 (2%)	64,73,73	0.62	0
2	NAD	H	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.67	0
2	NAD	A	401	-	46,48,48	0.67	1 (2%)	64,73,73	0.66	1 (1%)
2	NAD	C	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.61	0
2	NAD	F	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.62	1 (1%)
2	NAD	J	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.62	0
2	NAD	B	401	-	46,48,48	0.65	1 (2%)	64,73,73	0.74	2 (3%)
2	NAD	O	401	-	46,48,48	0.64	1 (2%)	64,73,73	0.68	0
2	NAD	D	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.67	0
2	NAD	I	401	-	46,48,48	0.67	1 (2%)	64,73,73	0.66	0
2	NAD	E	401	-	46,48,48	0.67	1 (2%)	64,73,73	0.73	1 (1%)
2	NAD	K	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.60	0
2	NAD	G	401	-	46,48,48	0.66	1 (2%)	64,73,73	0.58	0
3	EDO	A	402	-	3,3,3	0.42	0	2,2,2	0.38	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
2	NAD	P	401	-	-	11/30/62/62	0/5/5/5
2	NAD	M	401	-	-	14/30/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	L	401	-	-	7/30/62/62	0/5/5/5
3	EDO	D	402	-	-	1/1/1/1	-
2	NAD	N	401	-	-	6/30/62/62	0/5/5/5
2	NAD	H	401	-	-	7/30/62/62	0/5/5/5
2	NAD	A	401	-	-	12/30/62/62	0/5/5/5
2	NAD	C	401	-	-	6/30/62/62	0/5/5/5
2	NAD	F	401	-	-	12/30/62/62	0/5/5/5
2	NAD	J	401	-	-	6/30/62/62	0/5/5/5
2	NAD	B	401	-	-	7/30/62/62	0/5/5/5
2	NAD	O	401	-	-	4/30/62/62	0/5/5/5
2	NAD	D	401	-	-	7/30/62/62	0/5/5/5
2	NAD	I	401	-	-	10/30/62/62	0/5/5/5
2	NAD	E	401	-	-	15/30/62/62	0/5/5/5
2	NAD	K	401	-	-	6/30/62/62	0/5/5/5
2	NAD	G	401	-	-	5/30/62/62	0/5/5/5
3	EDO	A	402	-	-	0/1/1/1	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAD	C2N-N1N	2.85	1.38	1.35
2	E	401	NAD	C2N-N1N	2.85	1.38	1.35
2	L	401	NAD	C2N-N1N	2.84	1.38	1.35
2	C	401	NAD	C2N-N1N	2.84	1.38	1.35
2	H	401	NAD	C2N-N1N	2.83	1.38	1.35
2	K	401	NAD	C2N-N1N	2.83	1.38	1.35
2	P	401	NAD	C2N-N1N	2.82	1.38	1.35
2	F	401	NAD	C2N-N1N	2.82	1.38	1.35
2	I	401	NAD	C2N-N1N	2.82	1.38	1.35
2	N	401	NAD	C2N-N1N	2.81	1.38	1.35
2	G	401	NAD	C2N-N1N	2.81	1.38	1.35
2	D	401	NAD	C2N-N1N	2.81	1.38	1.35
2	J	401	NAD	C2N-N1N	2.81	1.38	1.35
2	B	401	NAD	C2N-N1N	2.80	1.38	1.35
2	M	401	NAD	C2N-N1N	2.76	1.38	1.35
2	O	401	NAD	C2N-N1N	2.76	1.38	1.35

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	401	NAD	C4D-O4D-C1D	-2.47	107.66	109.92
2	M	401	NAD	C4B-O4B-C1B	-2.22	104.57	109.47
2	B	401	NAD	O4B-C1B-N9A	2.15	112.22	108.09
2	E	401	NAD	C6N-N1N-C2N	-2.14	120.06	121.88
2	M	401	NAD	O4B-C1B-C2B	-2.08	102.16	106.62
2	M	401	NAD	O4B-C1B-N9A	2.08	112.08	108.09
2	B	401	NAD	C6N-N1N-C2N	-2.06	120.12	121.88
2	A	401	NAD	C6N-N1N-C2N	-2.05	120.14	121.88
2	F	401	NAD	C6N-N1N-C2N	-2.03	120.15	121.88
2	L	401	NAD	C6N-N1N-C2N	-2.02	120.16	121.88

There are no chirality outliers.

All (136) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAD	C5D-O5D-PN-O2N
2	B	401	NAD	O4D-C1D-N1N-C2N
2	B	401	NAD	O4D-C1D-N1N-C6N
2	B	401	NAD	C2D-C1D-N1N-C2N
2	B	401	NAD	C2D-C1D-N1N-C6N
2	C	401	NAD	O4D-C1D-N1N-C2N
2	C	401	NAD	O4D-C1D-N1N-C6N
2	C	401	NAD	C2D-C1D-N1N-C2N
2	C	401	NAD	C2D-C1D-N1N-C6N
2	D	401	NAD	O4D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	C2D-C1D-N1N-C2N
2	D	401	NAD	C2D-C1D-N1N-C6N
2	E	401	NAD	C5B-O5B-PA-O1A
2	E	401	NAD	C5B-O5B-PA-O2A
2	E	401	NAD	C5B-O5B-PA-O3
2	E	401	NAD	O4D-C1D-N1N-C2N
2	E	401	NAD	O4D-C1D-N1N-C6N
2	E	401	NAD	C2D-C1D-N1N-C2N
2	E	401	NAD	C2D-C1D-N1N-C6N
2	F	401	NAD	PN-O3-PA-O5B
2	F	401	NAD	C5D-O5D-PN-O1N
2	F	401	NAD	O4D-C1D-N1N-C2N
2	F	401	NAD	O4D-C1D-N1N-C6N
2	F	401	NAD	C2D-C1D-N1N-C2N
2	F	401	NAD	C2D-C1D-N1N-C6N
2	G	401	NAD	O4D-C1D-N1N-C2N
2	G	401	NAD	O4D-C1D-N1N-C6N

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Mol	Chain	Res	Type	Atoms
2	G	401	NAD	C2D-C1D-N1N-C2N
2	G	401	NAD	C2D-C1D-N1N-C6N
2	H	401	NAD	O4D-C1D-N1N-C2N
2	H	401	NAD	O4D-C1D-N1N-C6N
2	H	401	NAD	C2D-C1D-N1N-C2N
2	H	401	NAD	C2D-C1D-N1N-C6N
2	I	401	NAD	PA-O3-PN-O5D
2	I	401	NAD	C5D-O5D-PN-O3
2	I	401	NAD	C5D-O5D-PN-O2N
2	J	401	NAD	O4D-C1D-N1N-C2N
2	J	401	NAD	O4D-C1D-N1N-C6N
2	J	401	NAD	C2D-C1D-N1N-C2N
2	J	401	NAD	C2D-C1D-N1N-C6N
2	K	401	NAD	O4D-C1D-N1N-C2N
2	K	401	NAD	O4D-C1D-N1N-C6N
2	K	401	NAD	C2D-C1D-N1N-C2N
2	K	401	NAD	C2D-C1D-N1N-C6N
2	L	401	NAD	O4D-C1D-N1N-C2N
2	L	401	NAD	O4D-C1D-N1N-C6N
2	L	401	NAD	C2D-C1D-N1N-C2N
2	L	401	NAD	C2D-C1D-N1N-C6N
2	M	401	NAD	C5B-O5B-PA-O2A
2	M	401	NAD	C5B-O5B-PA-O3
2	M	401	NAD	C5D-O5D-PN-O3
2	M	401	NAD	C5D-O5D-PN-O2N
2	M	401	NAD	O4D-C4D-C5D-O5D
2	N	401	NAD	O4D-C1D-N1N-C2N
2	N	401	NAD	O4D-C1D-N1N-C6N
2	N	401	NAD	C2D-C1D-N1N-C2N
2	N	401	NAD	C2D-C1D-N1N-C6N
2	O	401	NAD	O4D-C1D-N1N-C2N
2	O	401	NAD	O4D-C1D-N1N-C6N
2	O	401	NAD	C2D-C1D-N1N-C6N
2	P	401	NAD	C5B-O5B-PA-O1A
2	P	401	NAD	C5B-O5B-PA-O3
2	P	401	NAD	O4D-C1D-N1N-C2N
2	P	401	NAD	O4D-C1D-N1N-C6N
2	P	401	NAD	C2D-C1D-N1N-C2N
2	P	401	NAD	C2D-C1D-N1N-C6N
2	F	401	NAD	O4D-C4D-C5D-O5D
2	M	401	NAD	C3D-C4D-C5D-O5D
2	A	401	NAD	O4D-C4D-C5D-O5D

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Mol	Chain	Res	Type	Atoms
2	A	401	NAD	C3D-C4D-C5D-O5D
2	E	401	NAD	O4D-C4D-C5D-O5D
2	F	401	NAD	C3D-C4D-C5D-O5D
2	A	401	NAD	O4B-C4B-C5B-O5B
2	A	401	NAD	C3B-C4B-C5B-O5B
2	E	401	NAD	O4B-C4B-C5B-O5B
2	M	401	NAD	O4B-C4B-C5B-O5B
2	E	401	NAD	C3B-C4B-C5B-O5B
2	I	401	NAD	O4B-C4B-C5B-O5B
2	M	401	NAD	C3B-C4B-C5B-O5B
2	E	401	NAD	C3D-C4D-C5D-O5D
2	I	401	NAD	C3B-C4B-C5B-O5B
2	E	401	NAD	C2B-C1B-N9A-C4A
2	M	401	NAD	C2B-C1B-N9A-C4A
2	A	401	NAD	PA-O3-PN-O5D
2	E	401	NAD	PN-O3-PA-O5B
2	M	401	NAD	PN-O3-PA-O5B
2	C	401	NAD	C2B-C1B-N9A-C8A
2	J	401	NAD	C2B-C1B-N9A-C8A
2	K	401	NAD	C2B-C1B-N9A-C8A
2	N	401	NAD	C2B-C1B-N9A-C8A
2	F	401	NAD	PA-O3-PN-O2N
2	I	401	NAD	C4B-C5B-O5B-PA
2	E	401	NAD	C2B-C1B-N9A-C8A
2	M	401	NAD	C2B-C1B-N9A-C8A
2	P	401	NAD	C2B-C1B-N9A-C4A
3	D	402	EDO	O1-C1-C2-O2
2	A	401	NAD	C5D-O5D-PN-O3
2	A	401	NAD	C5D-O5D-PN-O1N
2	I	401	NAD	C5D-O5D-PN-O1N
2	M	401	NAD	C5D-O5D-PN-O1N
2	P	401	NAD	C5B-O5B-PA-O2A
2	A	401	NAD	C4B-C5B-O5B-PA
2	M	401	NAD	C4B-C5B-O5B-PA
2	A	401	NAD	C2B-C1B-N9A-C8A
2	F	401	NAD	C2B-C1B-N9A-C8A
2	I	401	NAD	C2B-C1B-N9A-C8A
2	L	401	NAD	C2B-C1B-N9A-C8A
2	P	401	NAD	C2B-C1B-N9A-C8A
2	A	401	NAD	C2B-C1B-N9A-C4A
2	D	401	NAD	C2B-C1B-N9A-C4A
2	I	401	NAD	C2B-C1B-N9A-C4A

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Mol	Chain	Res	Type	Atoms
2	L	401	NAD	C2B-C1B-N9A-C4A
2	D	401	NAD	C2B-C1B-N9A-C8A
2	H	401	NAD	C2B-C1B-N9A-C8A
2	O	401	NAD	C2D-C1D-N1N-C2N
2	H	401	NAD	C2B-C1B-N9A-C4A
2	P	401	NAD	O4B-C1B-N9A-C8A
2	B	401	NAD	PN-O3-PA-O1A
2	E	401	NAD	C4B-C5B-O5B-PA
2	D	401	NAD	O4B-C1B-N9A-C8A
2	L	401	NAD	O4B-C1B-N9A-C8A
2	M	401	NAD	C4D-C5D-O5D-PN
2	P	401	NAD	C4B-C5B-O5B-PA
2	A	401	NAD	O4B-C1B-N9A-C8A
2	H	401	NAD	O4B-C1B-N9A-C8A
2	I	401	NAD	O4B-C1B-N9A-C8A
2	J	401	NAD	O4B-C1B-N9A-C8A
2	B	401	NAD	PN-O3-PA-O2A
2	F	401	NAD	PA-O3-PN-O1N
2	G	401	NAD	PN-O3-PA-O2A
2	F	401	NAD	O4B-C4B-C5B-O5B
2	C	401	NAD	O4B-C1B-N9A-C8A
2	K	401	NAD	O4B-C1B-N9A-C8A
2	N	401	NAD	O4B-C1B-N9A-C8A
2	B	401	NAD	C2B-C1B-N9A-C8A

There are no ring outliers.

14 monomers are involved in 33 short contacts:

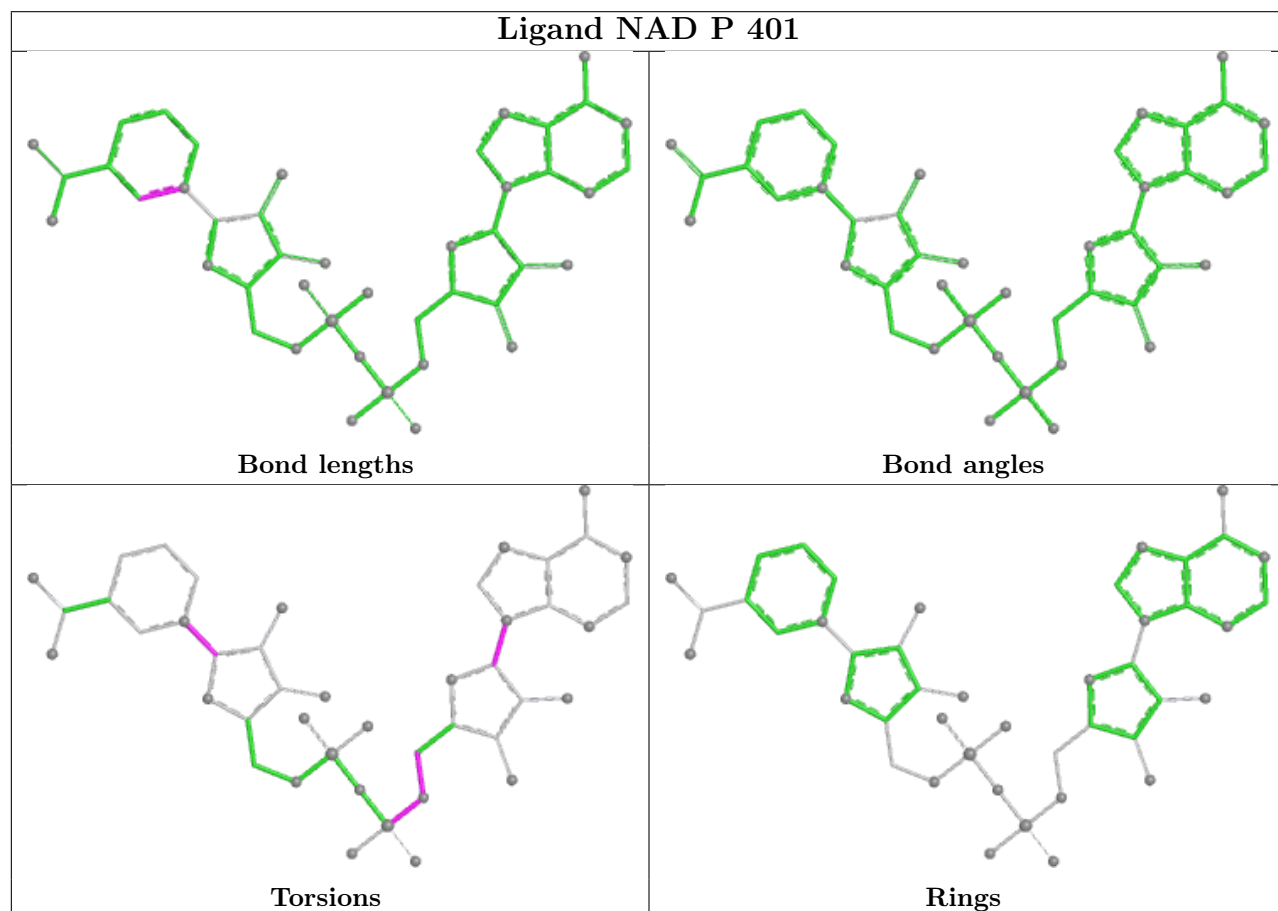
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	401	NAD	2	0
2	M	401	NAD	2	0
2	L	401	NAD	2	0
2	N	401	NAD	1	0
2	A	401	NAD	4	0
2	C	401	NAD	1	0
2	F	401	NAD	2	0
2	J	401	NAD	2	0
2	B	401	NAD	2	0
2	O	401	NAD	2	0
2	D	401	NAD	3	0
2	I	401	NAD	5	0
2	E	401	NAD	3	0

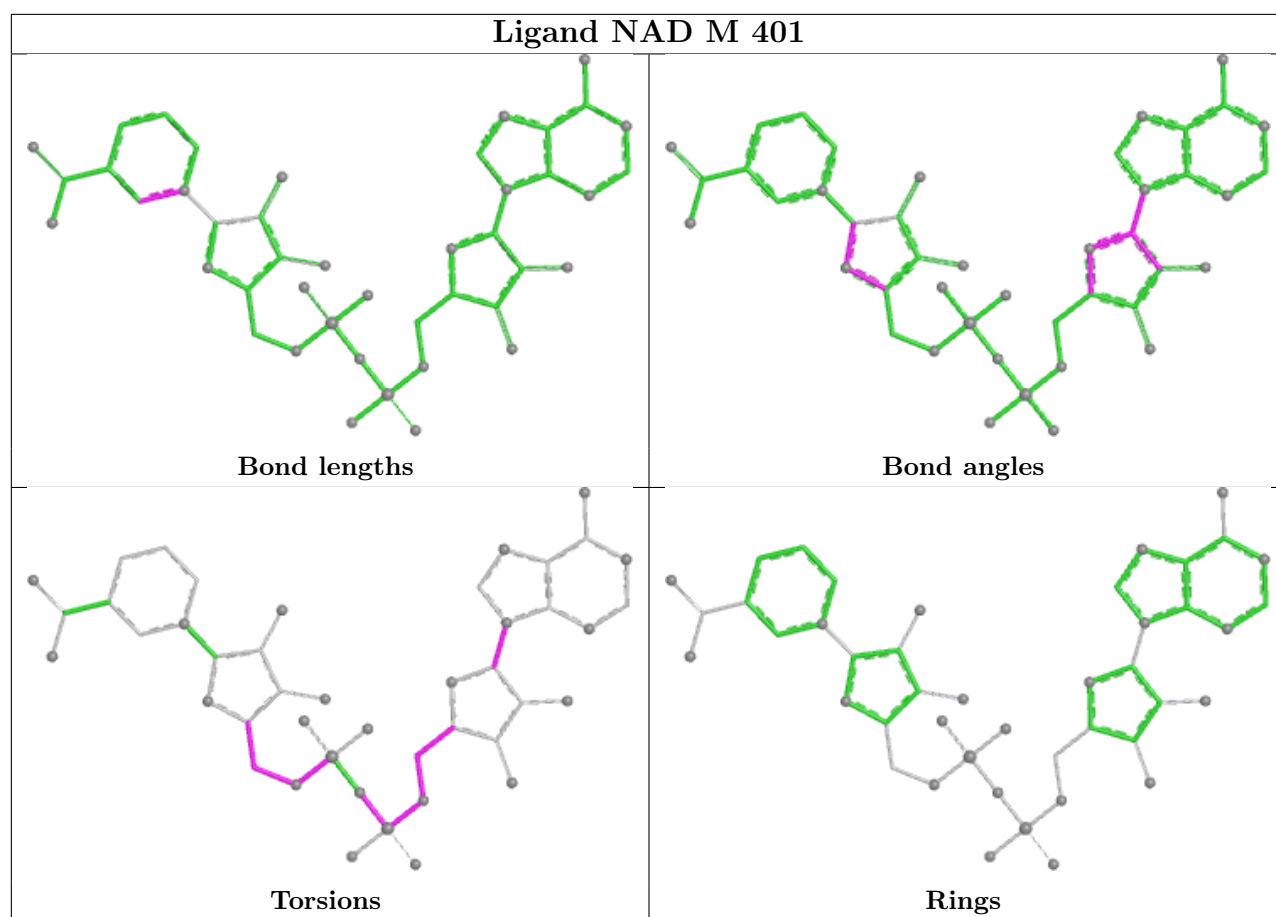
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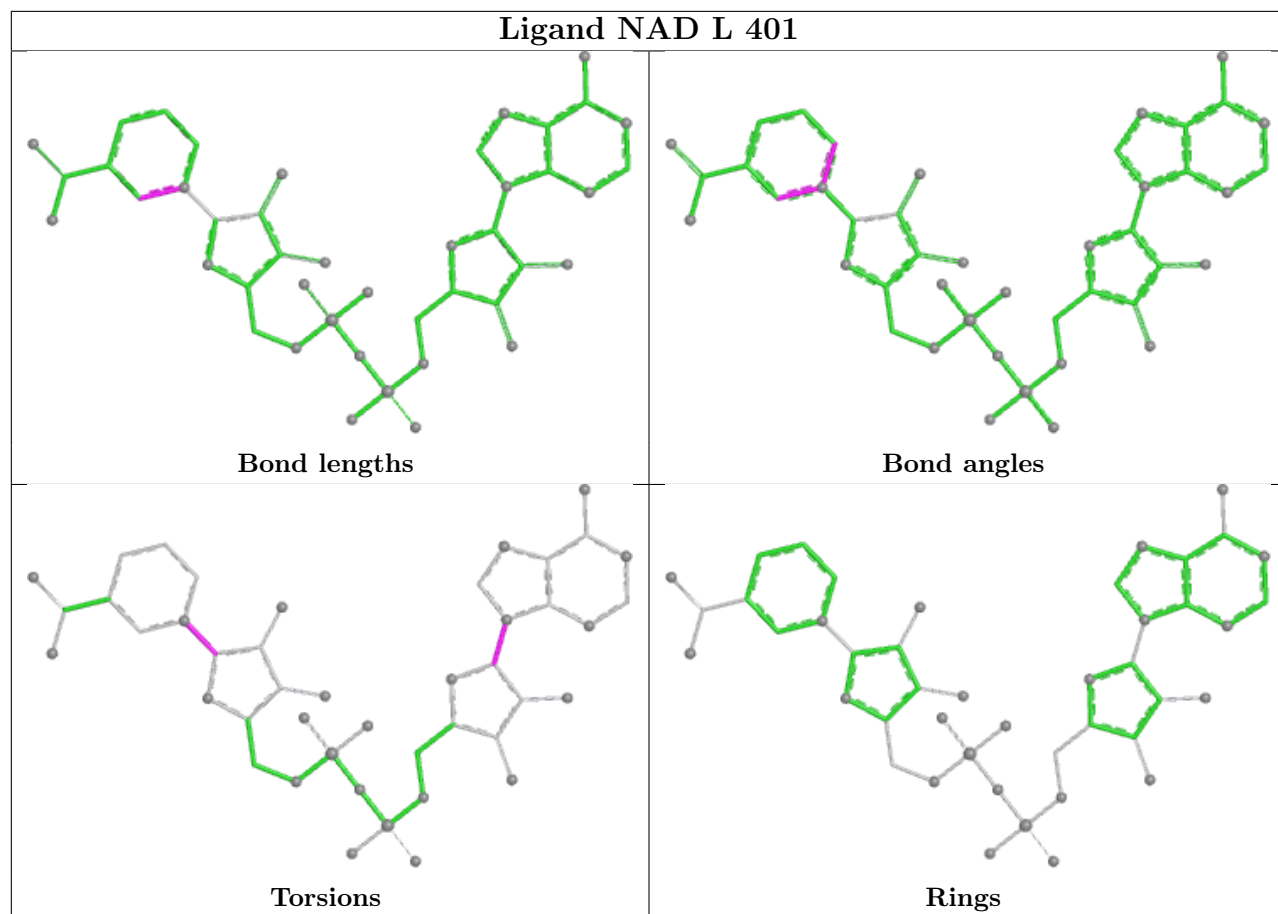
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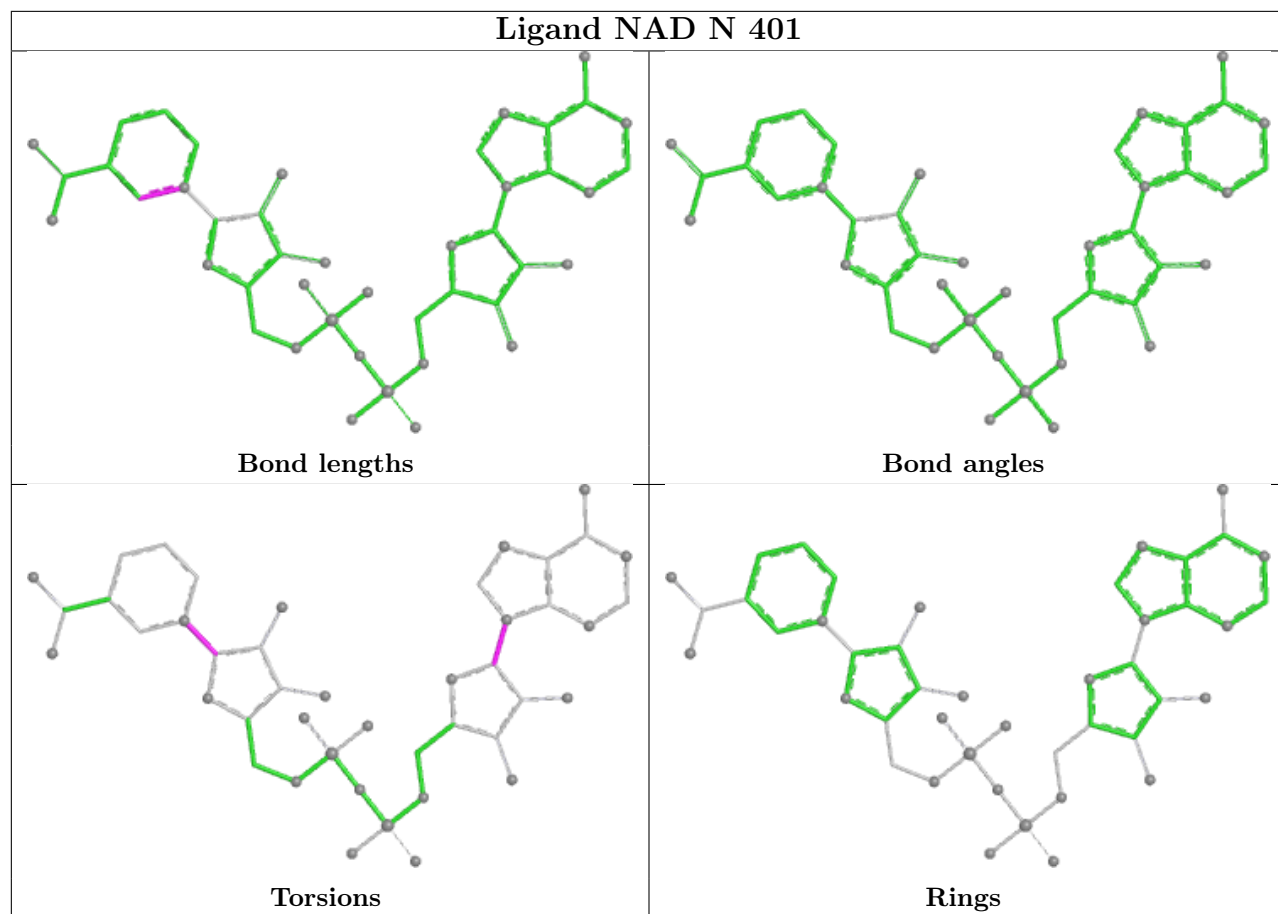
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	401	NAD	2	0

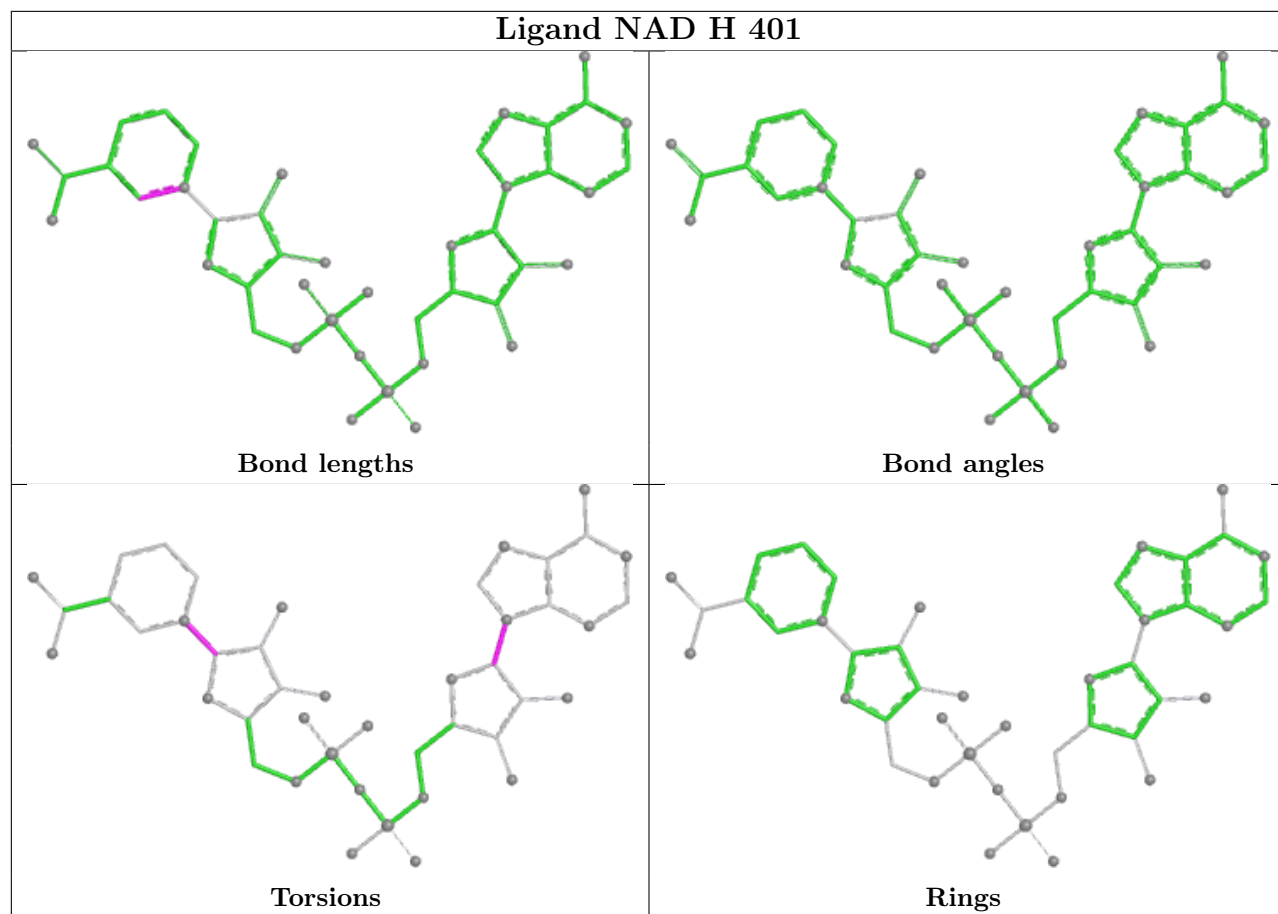
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

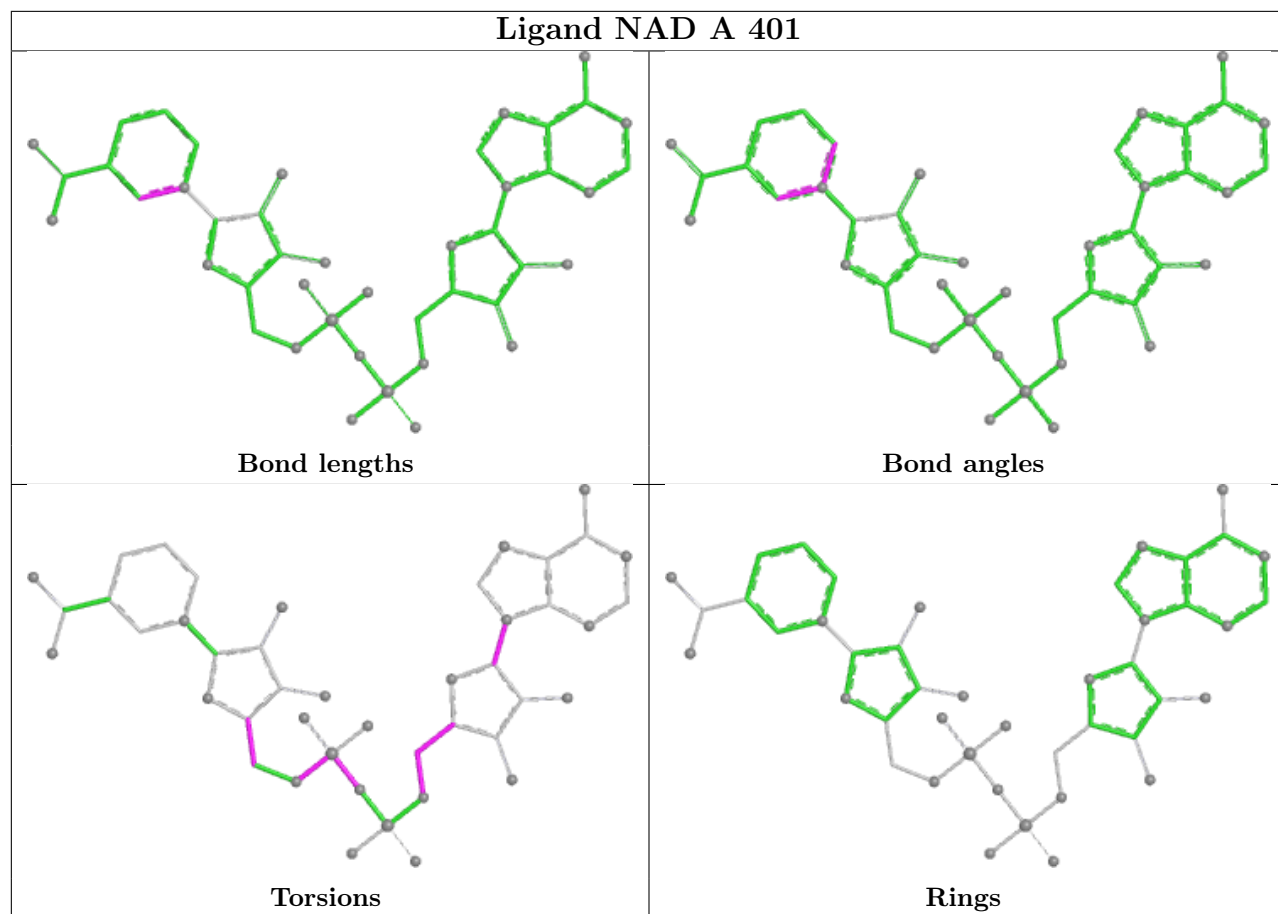


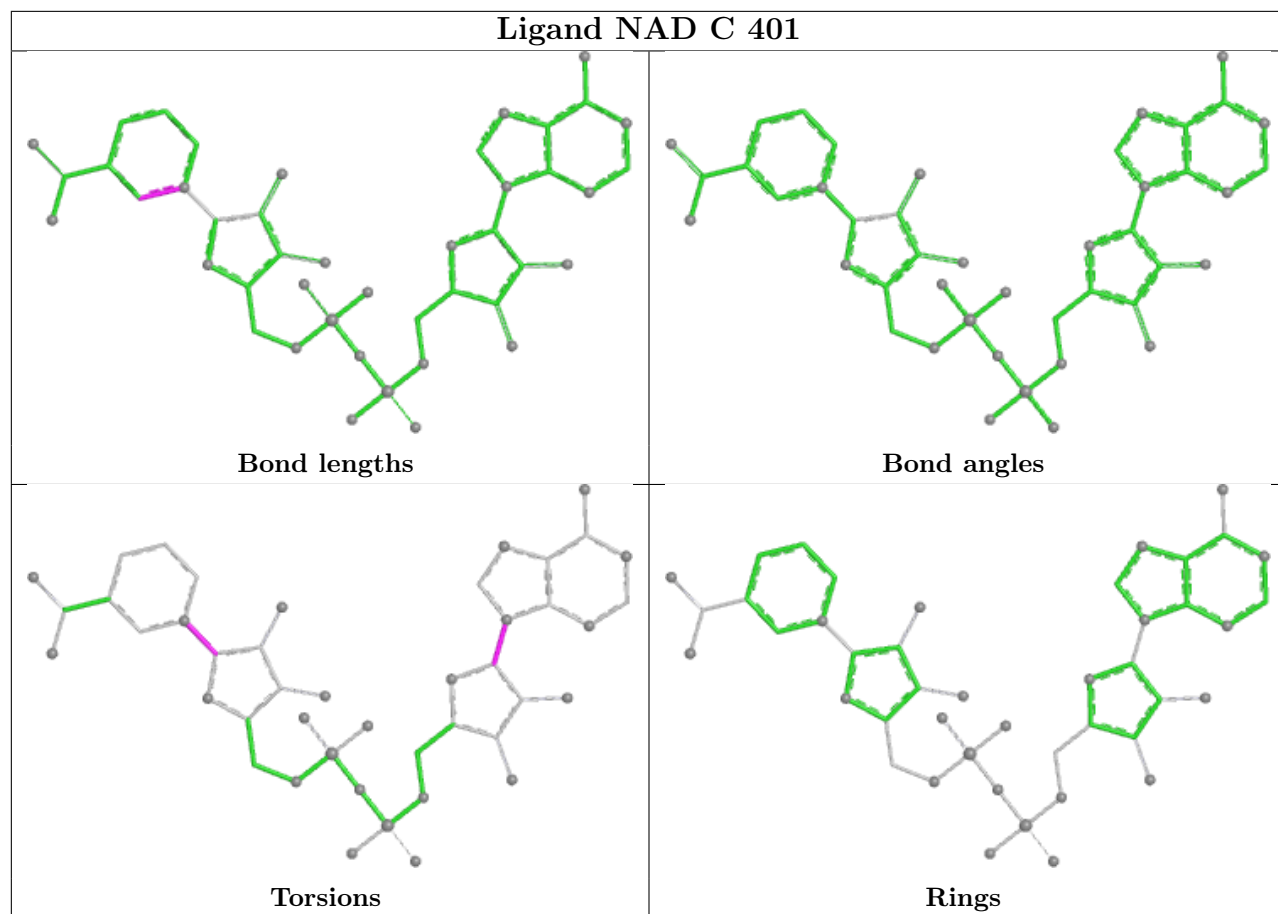


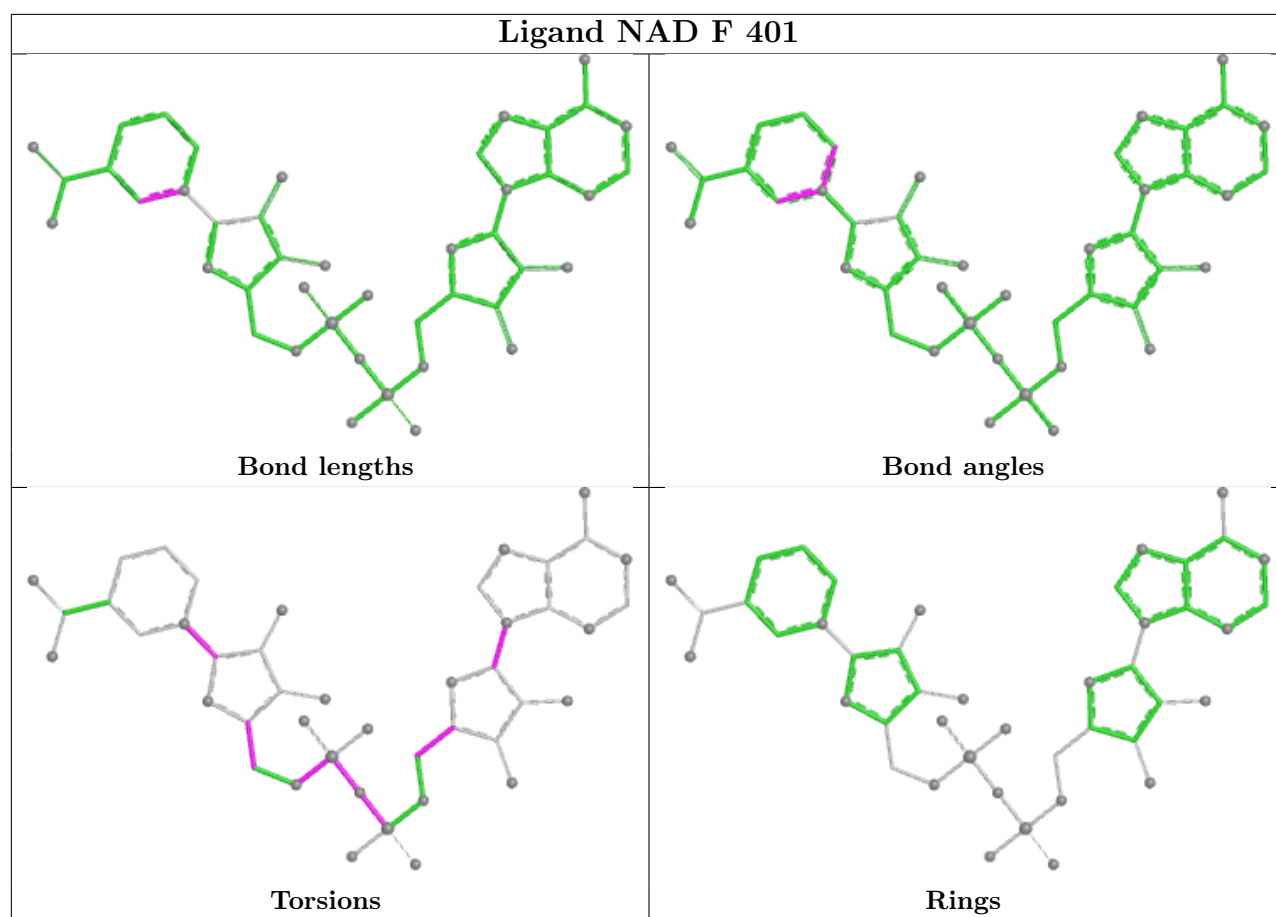


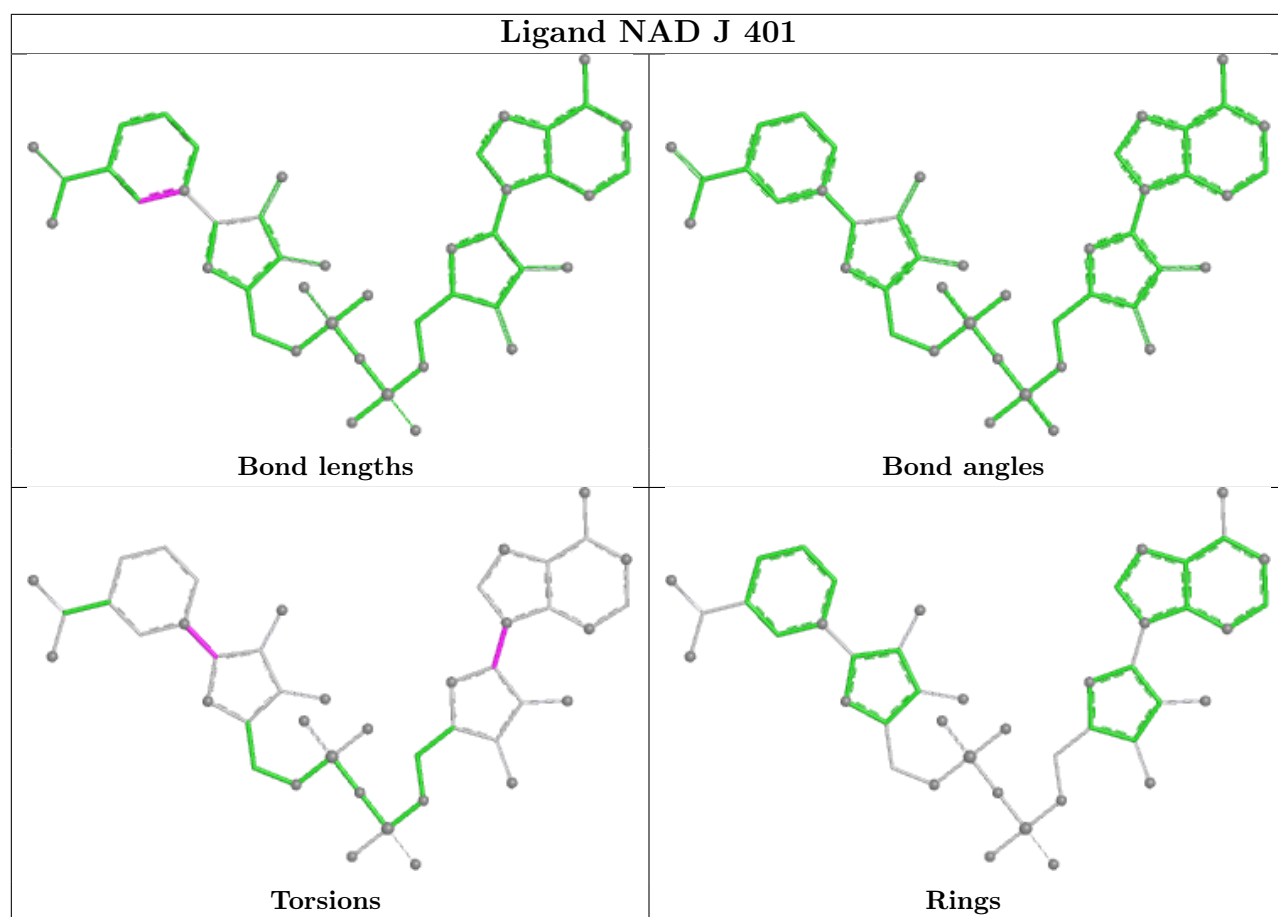


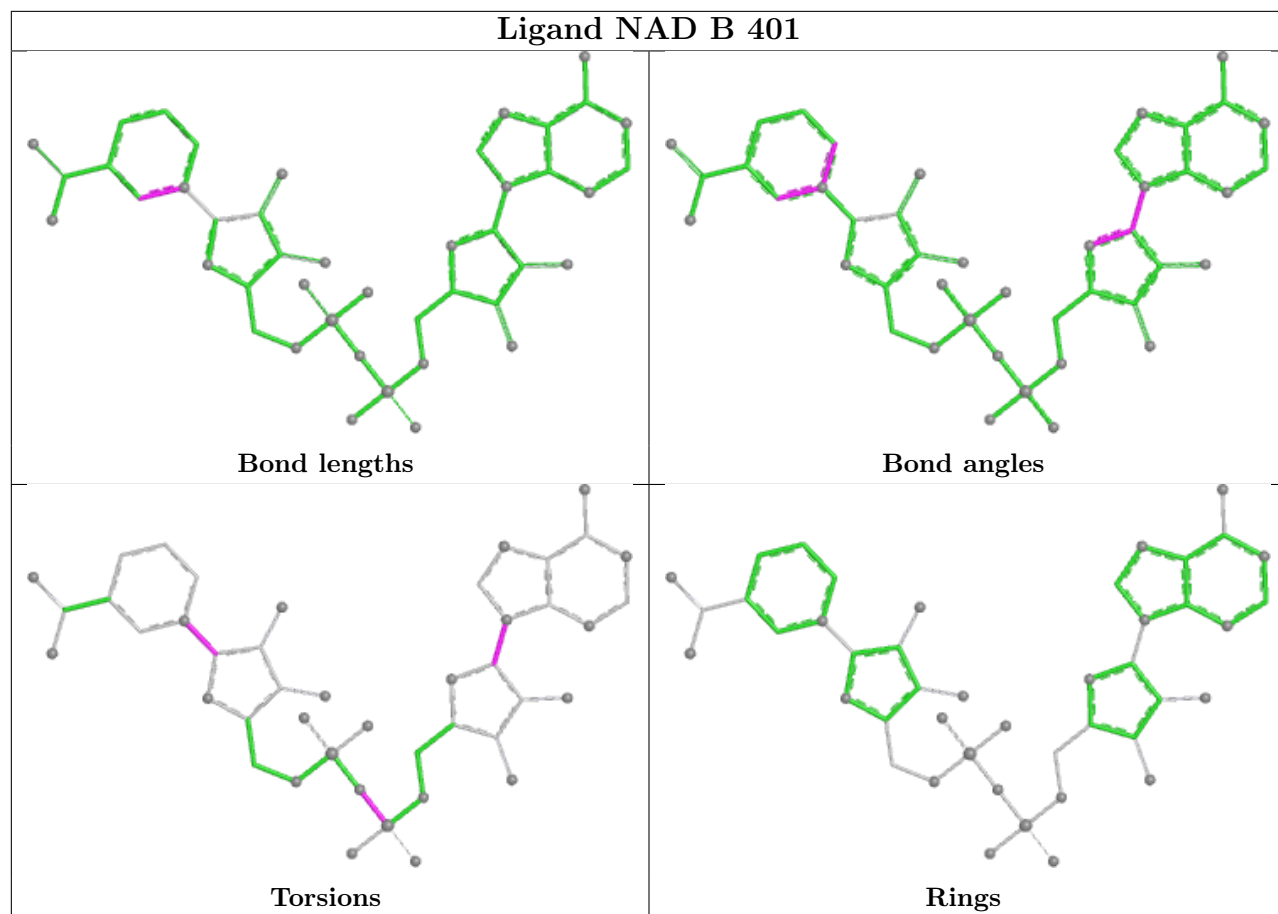


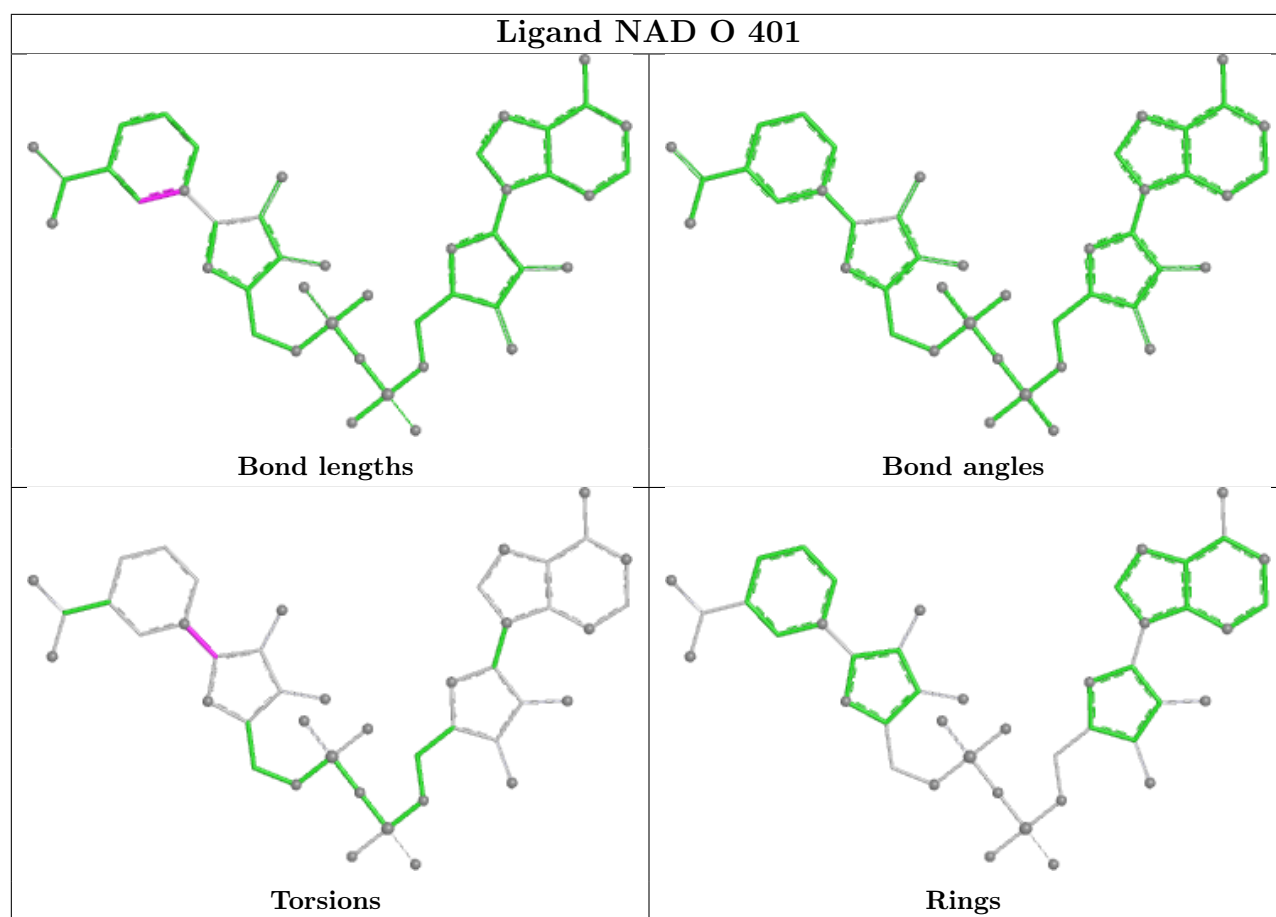


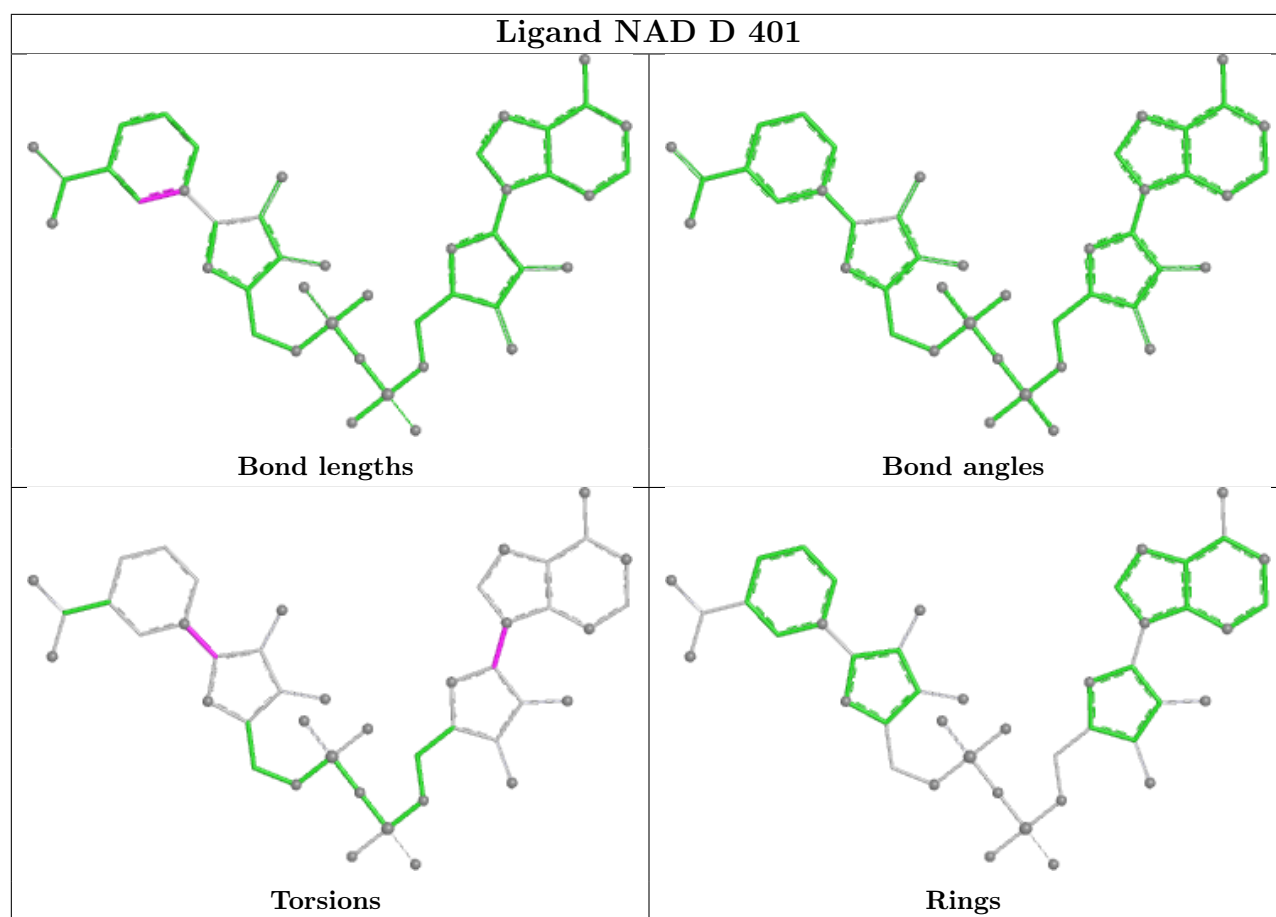


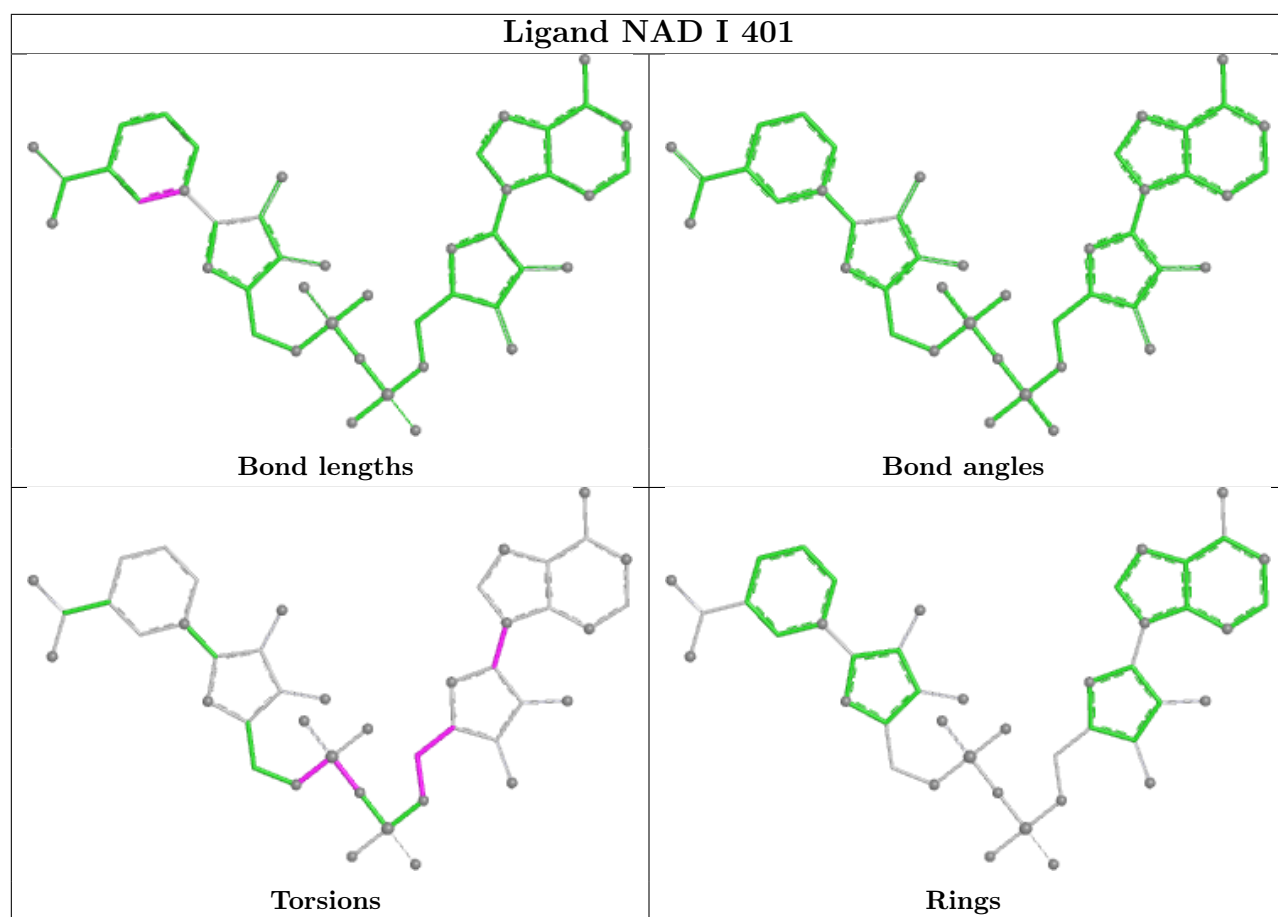


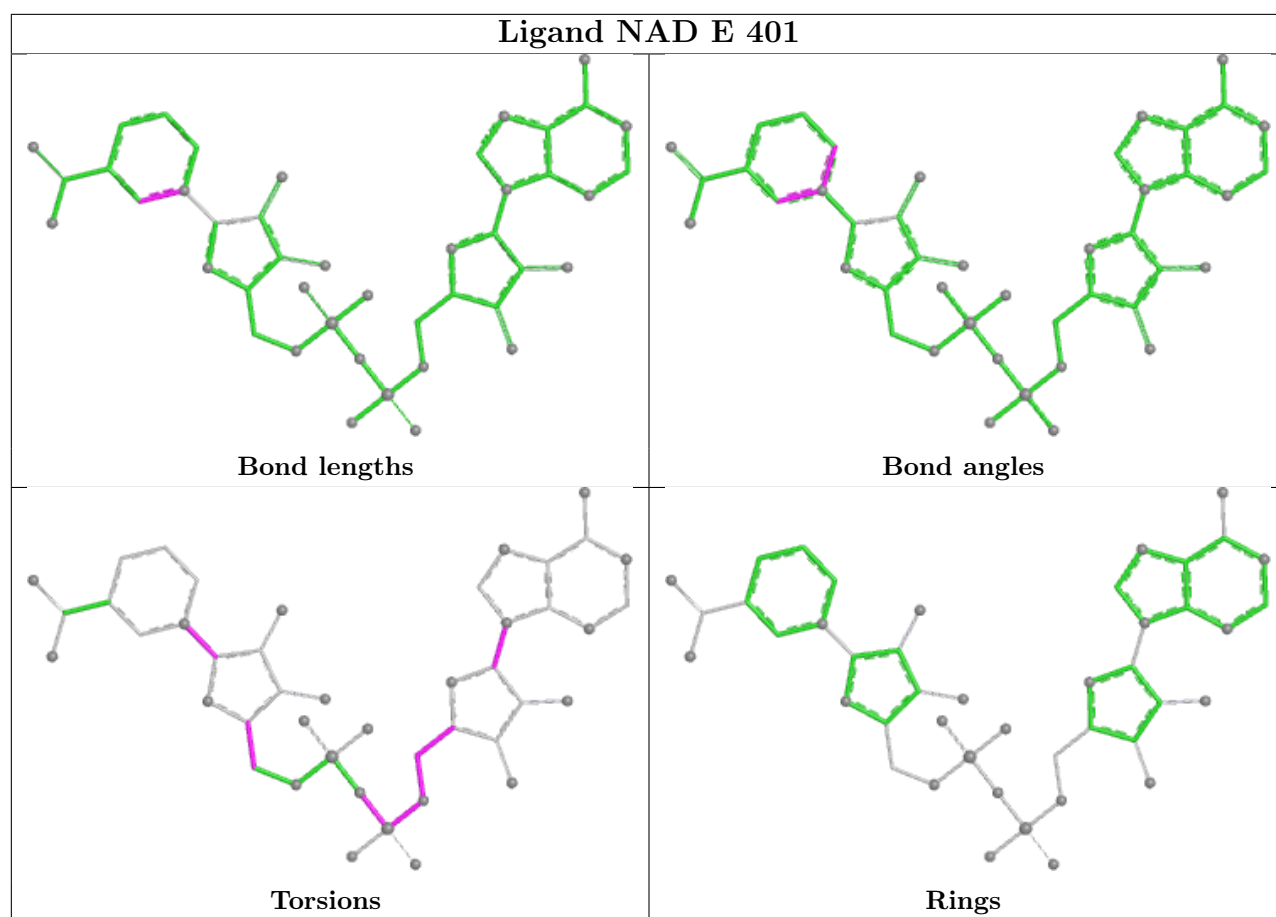


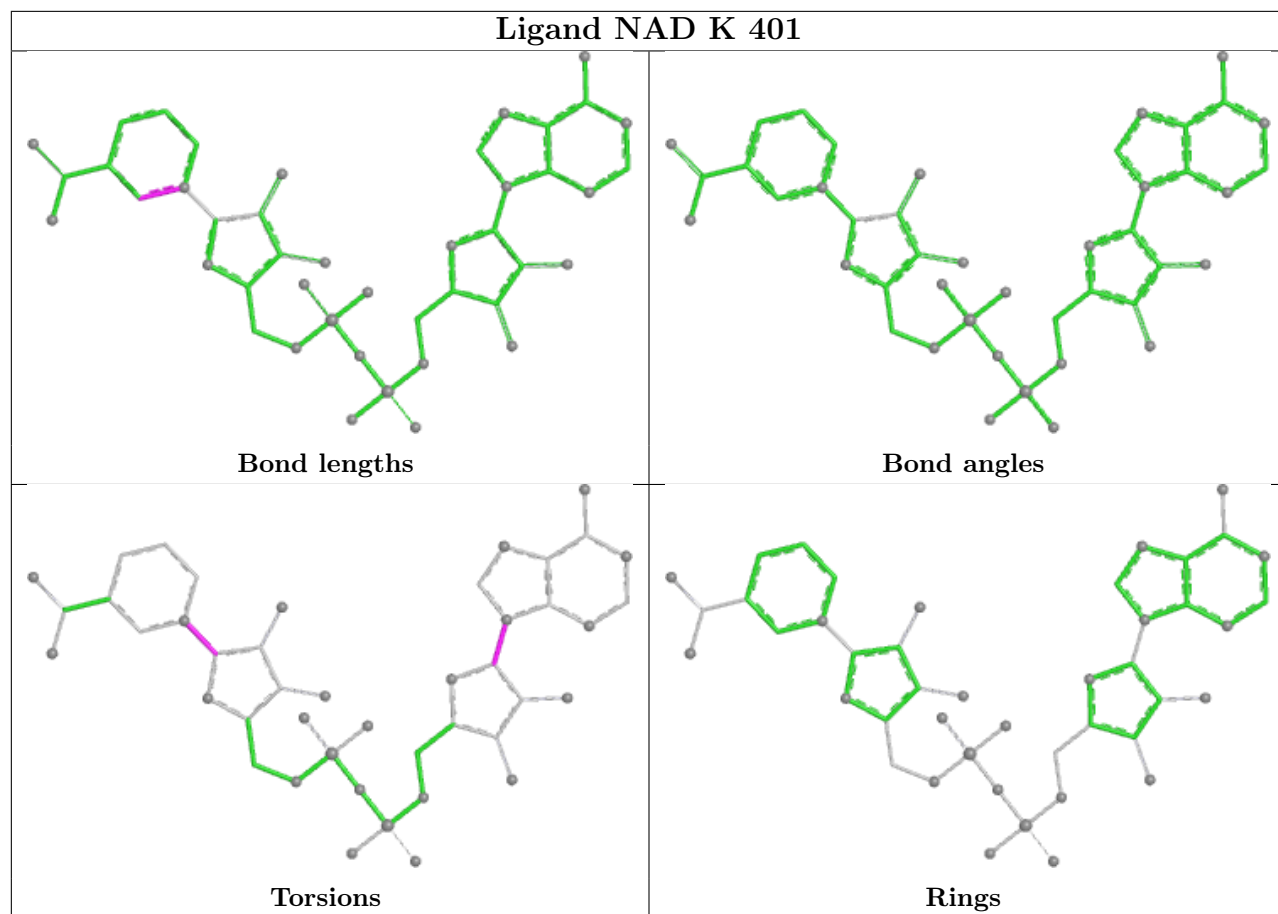


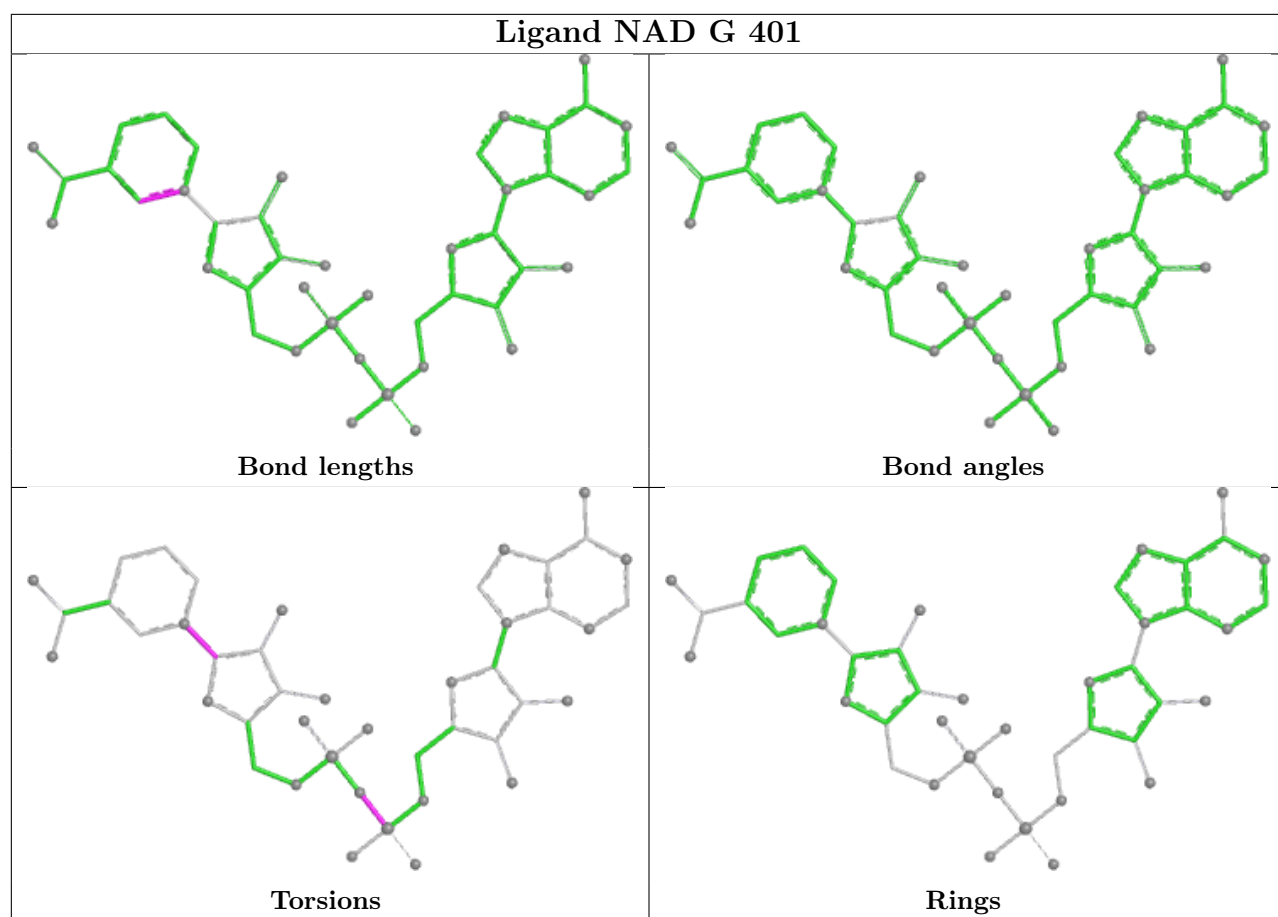












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	339/357 (94%)	0.47	2 (0%) 85 73	30, 47, 61, 77	0
1	B	338/357 (94%)	0.48	11 (3%) 49 33	33, 46, 63, 71	0
1	C	339/357 (94%)	0.49	7 (2%) 63 44	29, 47, 68, 85	0
1	D	338/357 (94%)	0.54	7 (2%) 63 44	33, 53, 67, 78	0
1	E	339/357 (94%)	0.43	11 (3%) 50 34	33, 43, 58, 69	0
1	F	338/357 (94%)	0.43	6 (1%) 67 49	30, 47, 60, 77	0
1	G	339/357 (94%)	0.50	11 (3%) 50 34	29, 47, 63, 75	0
1	H	338/357 (94%)	0.54	8 (2%) 59 40	33, 48, 63, 73	0
1	I	339/357 (94%)	0.48	3 (0%) 81 65	31, 47, 63, 75	0
1	J	338/357 (94%)	0.68	8 (2%) 59 40	39, 56, 71, 84	0
1	K	339/357 (94%)	0.65	15 (4%) 39 25	39, 56, 70, 81	0
1	L	338/357 (94%)	0.50	11 (3%) 49 33	35, 47, 62, 73	0
1	M	339/357 (94%)	0.61	7 (2%) 63 44	35, 57, 69, 85	0
1	N	338/357 (94%)	0.59	9 (2%) 56 37	36, 56, 72, 81	0
1	O	339/357 (94%)	0.77	16 (4%) 36 24	38, 66, 84, 104	0
1	P	338/357 (94%)	0.49	5 (1%) 72 53	34, 53, 67, 80	0
All	All	5416/5712 (94%)	0.54	137 (2%) 58 39	29, 50, 70, 104	0

All (137) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	82	ALA	4.2
1	I	97	VAL	4.1
1	O	235	GLY	4.0
1	C	34	VAL	3.9
1	A	82	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	N	244	THR	3.8
1	I	82	ALA	3.7
1	E	167	ALA	3.6
1	O	219	ALA	3.5
1	K	17	PHE	3.4
1	O	179	GLY	3.3
1	M	82	ALA	3.3
1	B	309	ASP	3.3
1	N	82	ALA	3.3
1	F	127	SER	3.2
1	J	165	PRO	3.2
1	D	118	ASP	3.2
1	D	82	ALA	3.2
1	J	325	SER	3.1
1	K	328	LEU	3.1
1	K	209	LEU	3.1
1	F	82	ALA	3.0
1	B	148	GLY	3.0
1	L	7	ILE	3.0
1	P	80	ALA	3.0
1	O	99	VAL	3.0
1	H	226	MET	3.0
1	O	139	LEU	2.9
1	O	33	GLU	2.9
1	K	306	LYS	2.9
1	O	121	ALA	2.9
1	F	66	GLU	2.9
1	G	309	ASP	2.9
1	E	139	LEU	2.8
1	J	166	LEU	2.8
1	O	236	TYR	2.8
1	E	1	MET	2.8
1	D	332	VAL	2.8
1	C	92	TRP	2.8
1	H	23	ALA	2.8
1	E	283	ASP	2.7
1	B	80	ALA	2.7
1	K	137	ILE	2.7
1	M	101	VAL	2.6
1	N	328	LEU	2.6
1	O	328	LEU	2.6
1	F	235	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	287	VAL	2.6
1	K	287	VAL	2.6
1	K	67	GLY	2.6
1	K	28	GLY	2.5
1	O	21	LEU	2.5
1	M	155	ASN	2.5
1	P	288	SER	2.5
1	G	181	MET	2.5
1	G	164	ALA	2.4
1	O	323	GLY	2.4
1	G	105	GLY	2.4
1	K	179	GLY	2.4
1	E	151	ASN	2.4
1	P	72	VAL	2.4
1	O	1	MET	2.4
1	C	310	ASP	2.4
1	H	336	GLY	2.4
1	H	82	ALA	2.4
1	D	328	LEU	2.4
1	B	2	THR	2.3
1	B	218	ALA	2.3
1	L	267	PHE	2.3
1	J	332	VAL	2.3
1	K	288	SER	2.3
1	N	108	THR	2.3
1	G	283	ASP	2.3
1	E	69	ASP	2.3
1	D	211	ILE	2.2
1	J	117	LEU	2.2
1	G	338	SER	2.2
1	K	289	SER	2.2
1	L	149	SER	2.2
1	D	69	ASP	2.2
1	F	283	ASP	2.2
1	L	193	LEU	2.2
1	G	257	ARG	2.2
1	L	211	ILE	2.2
1	L	49	LEU	2.2
1	N	229	LEU	2.2
1	M	92	TRP	2.2
1	J	212	VAL	2.2
1	J	259	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	K	149	SER	2.2
1	A	283	ASP	2.2
1	E	39	ASP	2.2
1	L	118	ASP	2.2
1	H	244	THR	2.1
1	M	244	THR	2.1
1	P	2	THR	2.1
1	B	66	GLU	2.1
1	M	80	ALA	2.1
1	K	329	VAL	2.1
1	C	153	ILE	2.1
1	E	282	TYR	2.1
1	M	1	MET	2.1
1	G	86	GLY	2.1
1	K	148	GLY	2.1
1	B	124	VAL	2.1
1	H	229	LEU	2.1
1	L	57	GLY	2.1
1	G	267	PHE	2.1
1	E	72	VAL	2.1
1	N	335	VAL	2.1
1	N	308	ILE	2.1
1	O	15	ARG	2.1
1	C	1	MET	2.1
1	E	74	GLY	2.1
1	O	231	GLY	2.1
1	F	314	VAL	2.1
1	H	225	VAL	2.1
1	B	153	ILE	2.1
1	H	181	MET	2.1
1	L	261	ASP	2.1
1	G	281	TYR	2.1
1	L	3	VAL	2.0
1	K	127	SER	2.0
1	B	291	ILE	2.0
1	C	137	ILE	2.0
1	J	211	ILE	2.0
1	P	193	LEU	2.0
1	L	244	THR	2.0
1	N	69	ASP	2.0
1	B	34	VAL	2.0
1	G	5	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	N	124	VAL	2.0
1	O	218	ALA	2.0
1	O	26	GLU	2.0
1	C	279	LEU	2.0
1	D	65	LEU	2.0
1	I	211	ILE	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.

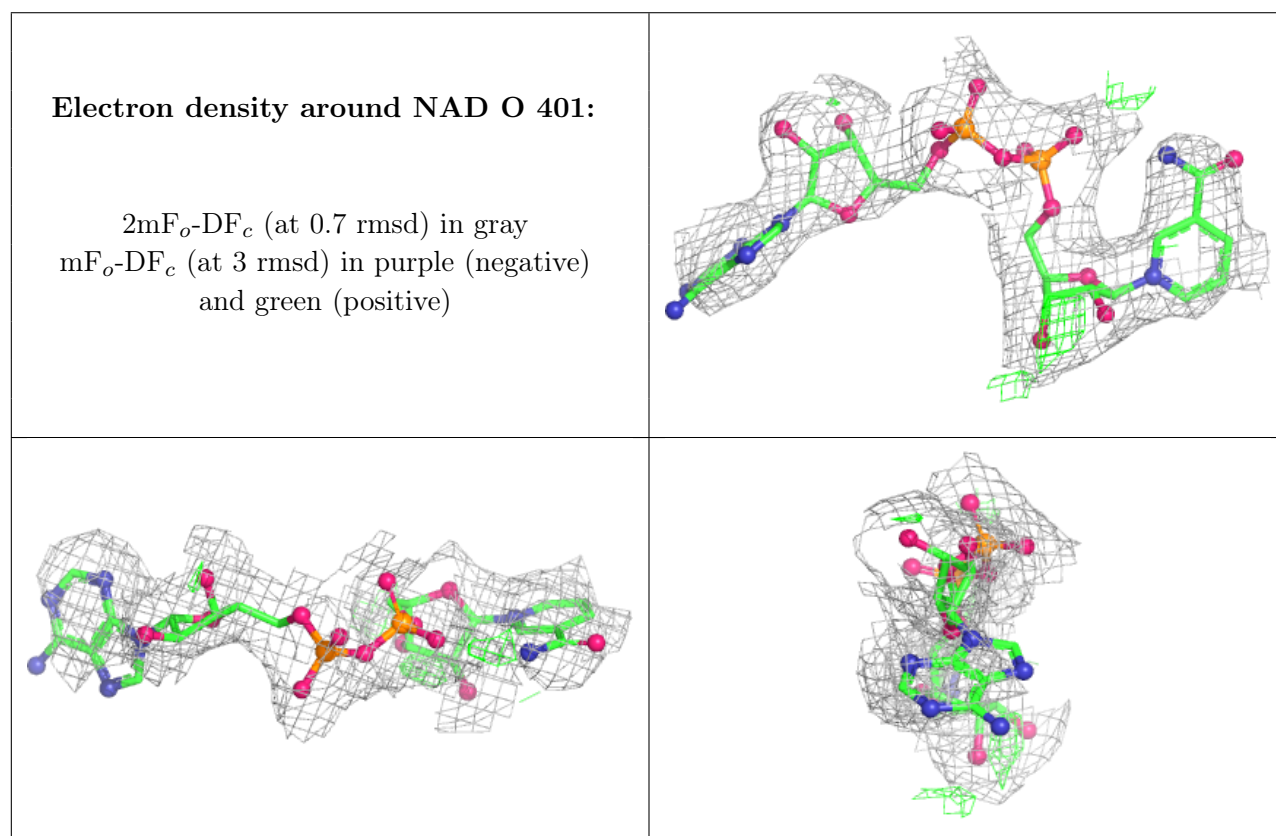
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	EDO	D	402	4/4	0.77	0.18	29,39,39,40	0
2	NAD	O	401	44/44	0.82	0.15	46,59,75,76	0
2	NAD	I	401	44/44	0.85	0.14	35,44,50,54	0
2	NAD	P	401	44/44	0.86	0.13	36,48,54,64	0
2	NAD	M	401	44/44	0.86	0.12	43,51,56,61	0
2	NAD	E	401	44/44	0.87	0.12	33,41,48,50	0
3	EDO	A	402	4/4	0.87	0.17	29,35,37,38	0
2	NAD	A	401	44/44	0.87	0.13	33,41,47,52	0
2	NAD	D	401	44/44	0.88	0.10	33,44,48,50	0
2	NAD	H	401	44/44	0.88	0.12	31,40,46,47	0
2	NAD	L	401	44/44	0.89	0.11	35,46,51,57	0
2	NAD	B	401	44/44	0.89	0.13	29,41,46,48	0
2	NAD	J	401	44/44	0.89	0.10	47,55,60,65	0
2	NAD	G	401	44/44	0.90	0.11	29,41,47,49	0
2	NAD	N	401	44/44	0.90	0.10	39,49,53,54	0
2	NAD	K	401	44/44	0.91	0.10	38,47,53,54	0
2	NAD	F	401	44/44	0.92	0.09	34,38,43,47	0

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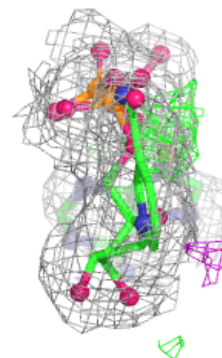
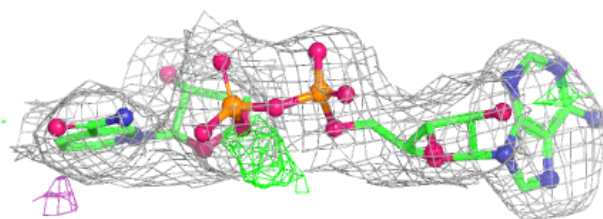
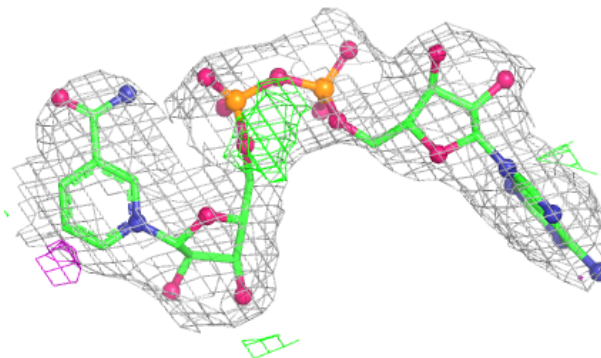
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAD	C	401	44/44	0.93	0.09	35,41,45,49	0

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.

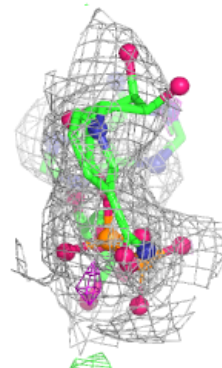
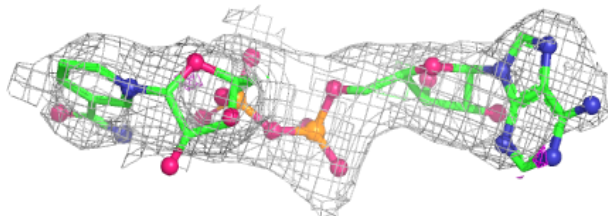
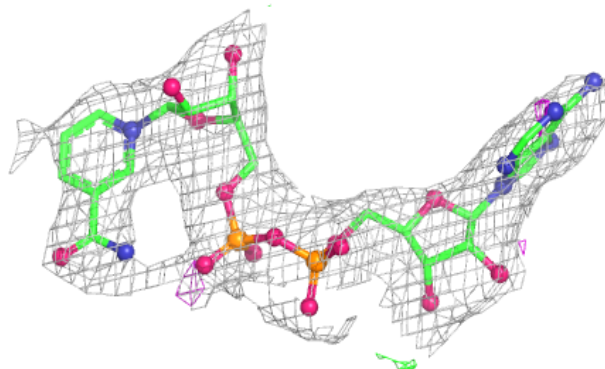


Electron density around NAD I 401:

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

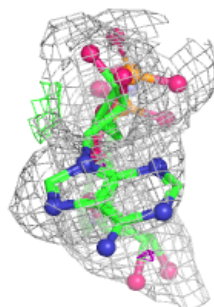
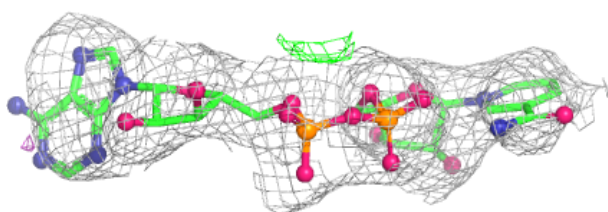
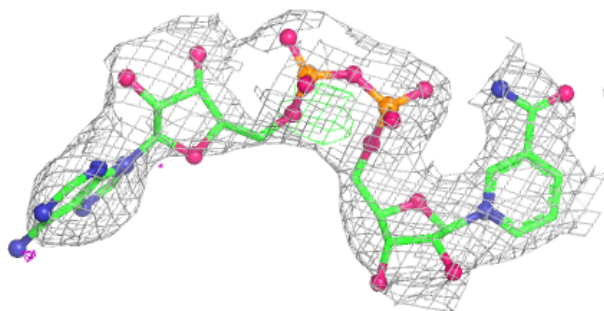
**Electron density around NAD P 401:**

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

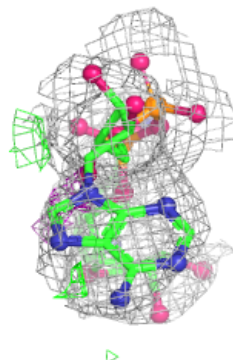
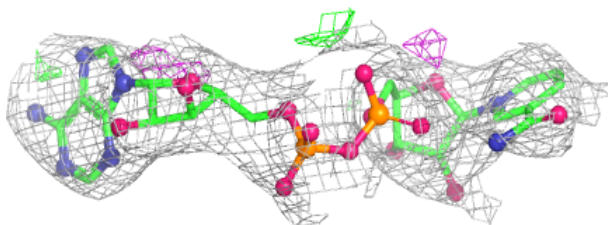
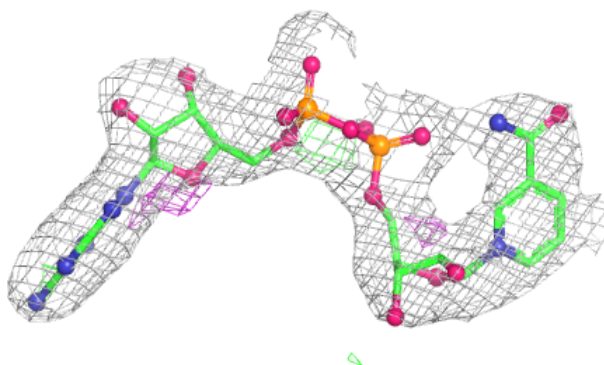


Electron density around NAD M 401:

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

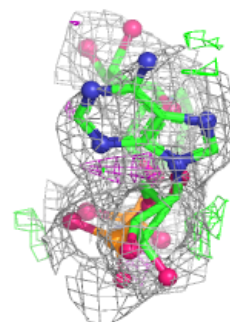
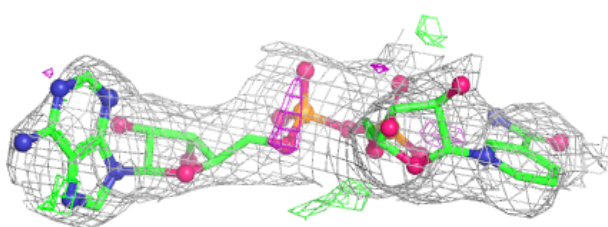
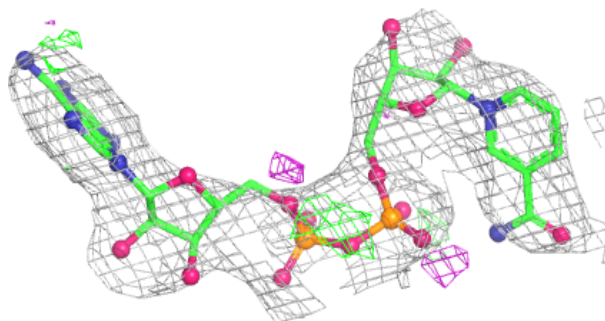
**Electron density around NAD E 401:**

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

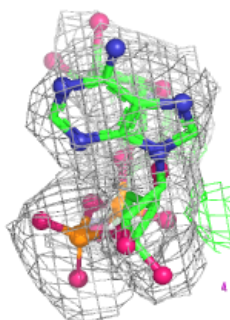
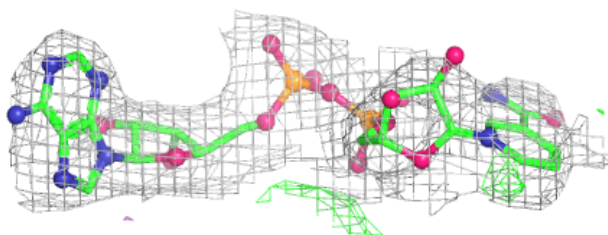
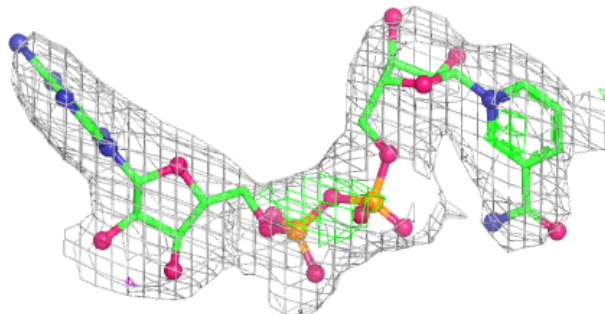


Electron density around NAD A 401:

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

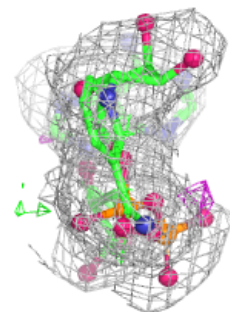
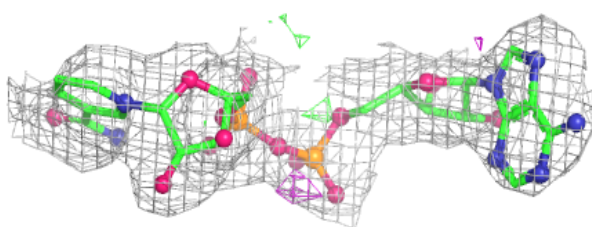
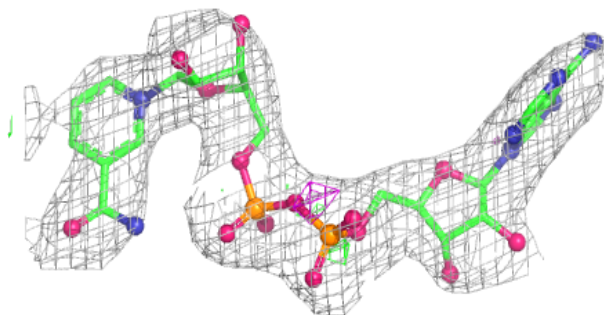
**Electron density around NAD D 401:**

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

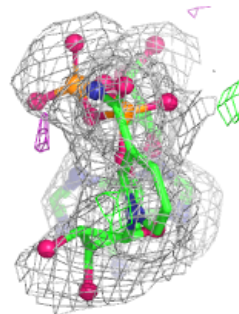
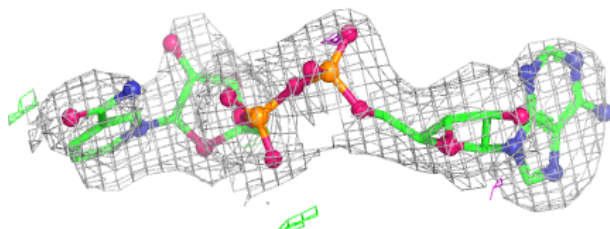
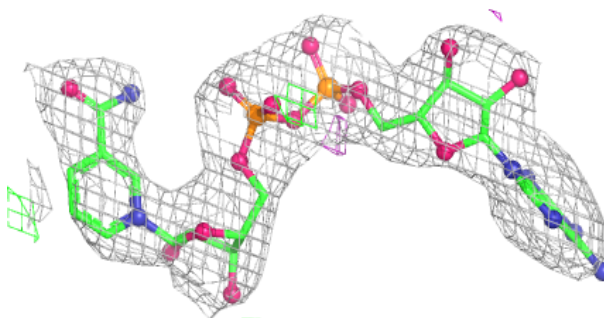


Electron density around NAD H 401:

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

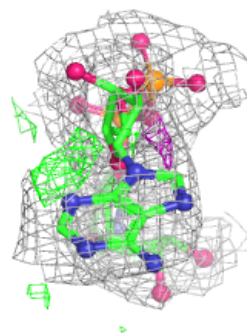
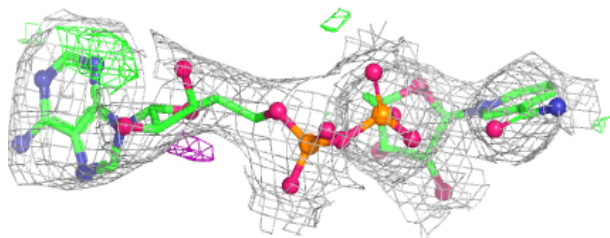
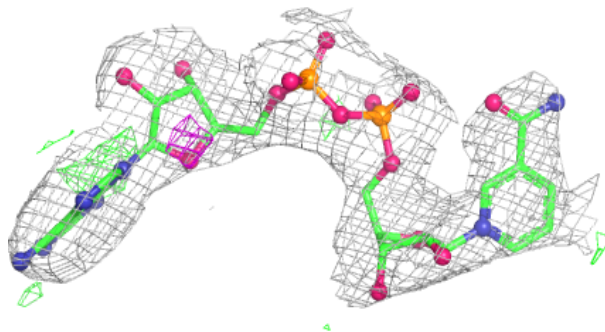
**Electron density around NAD L 401:**

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

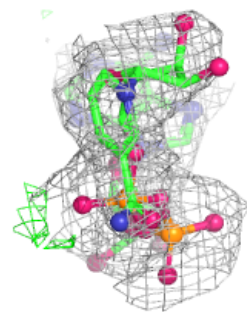
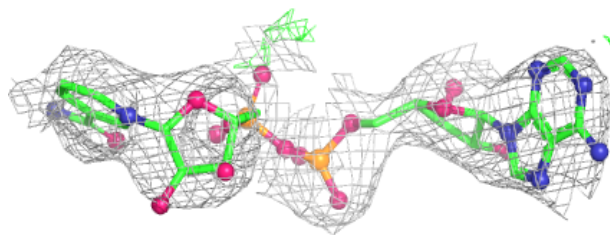
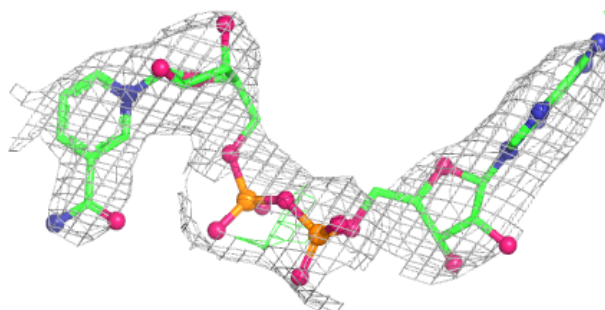


Electron density around NAD B 401:

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

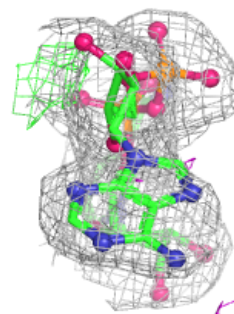
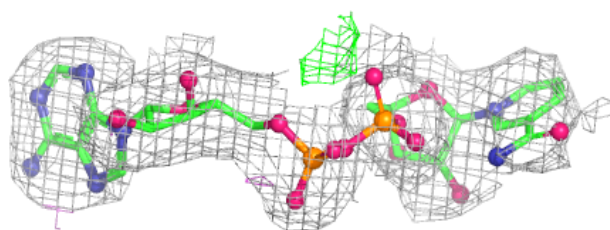
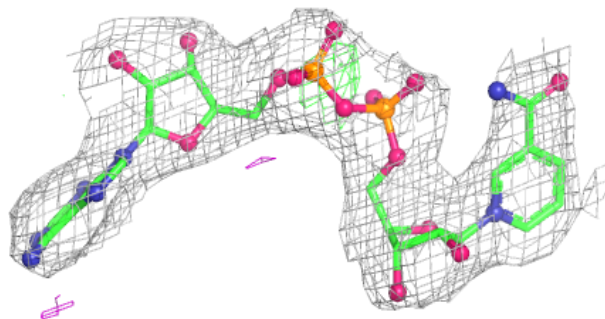
**Electron density around NAD J 401:**

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

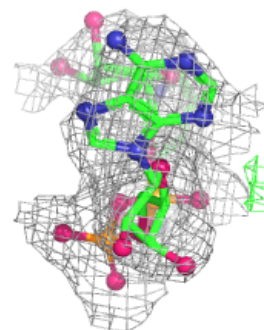
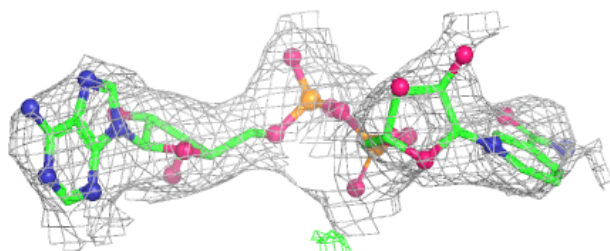
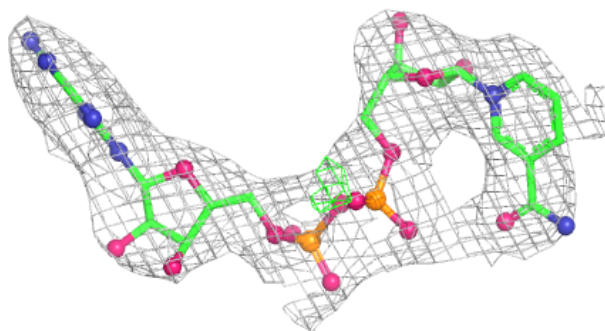


Electron density around NAD G 401:

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

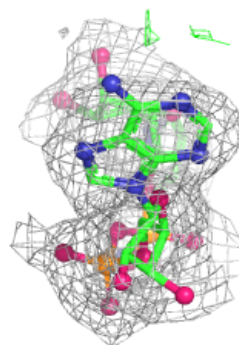
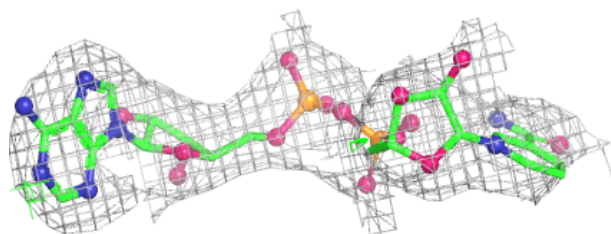
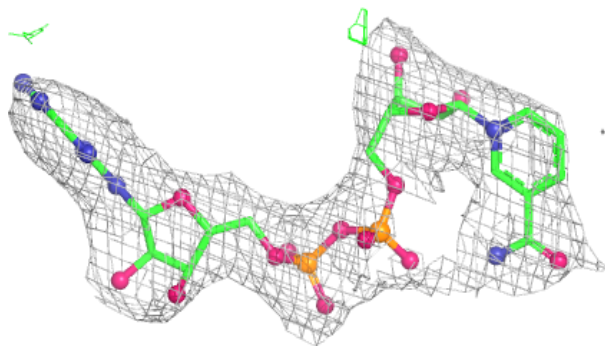
**Electron density around NAD N 401:**

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

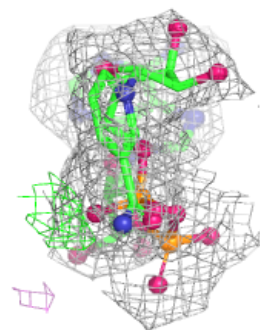
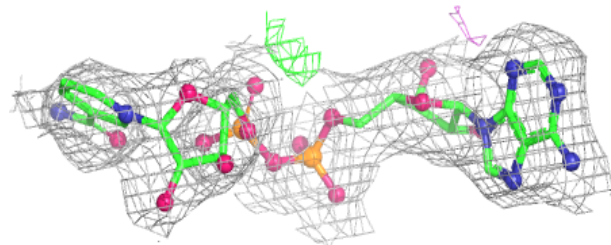
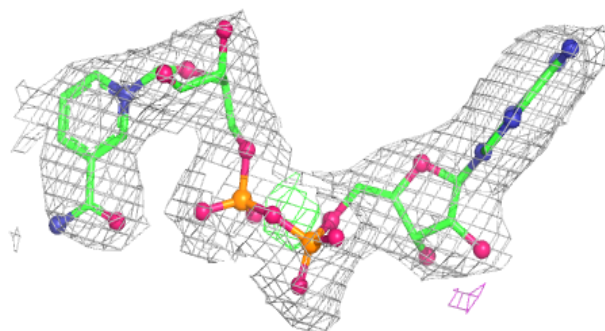


Electron density around NAD K 401:

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and green (positive)

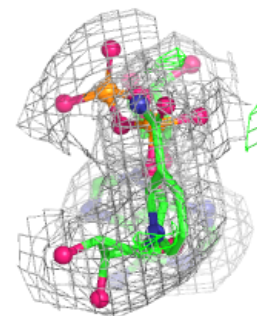
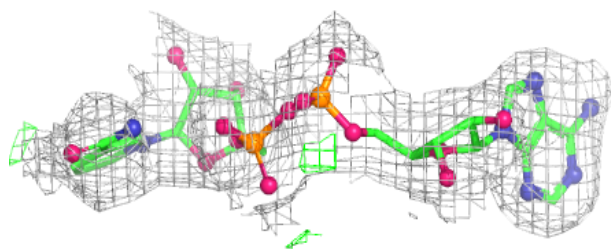
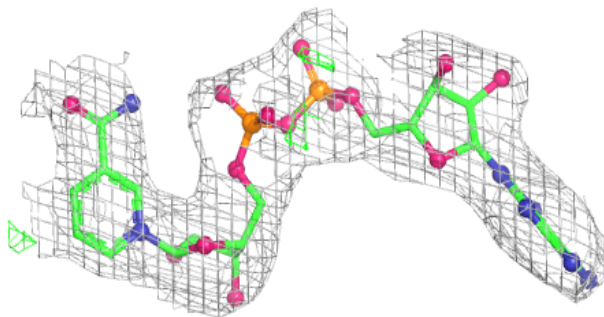
**Electron density around NAD F 401:**

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



Electron density around NAD C 401:

$2mF_o-DF_c$ (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.