



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:31 AM UTC

PDB ID : 4MVJ / pdb_00004mvj
Title : 2.85 Angstrom Resolution Crystal Structure of Glyceraldehyde 3-phosphate Dehydrogenase A (gapA) from Escherichia coli Modified by Acetyl Phosphate.
Authors : Minasov, G.; Kuhn, M.; Dubrovskaya, I.; Winsor, J.; Shuvalova, L.; Grimshaw, S.; Kwon, K.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2013-09-24
Resolution : 2.85 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:

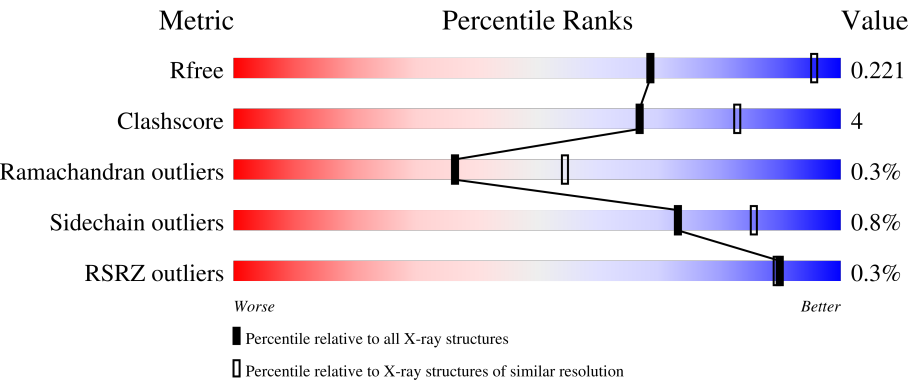
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 2.85 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.



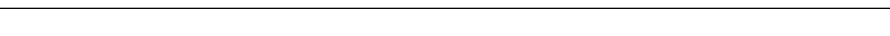
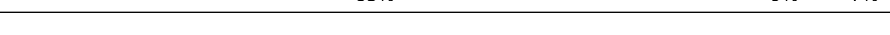
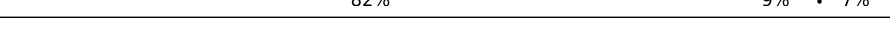
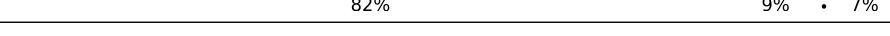
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	355	82% 9% • 7%
1	B	355	82% 9% • 6%
1	C	355	81% 10% • 7%
1	K	355	81% 11% • 7%
1	N	355	% 85% 8% • 7%

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Mol	Chain	Length	Quality of chain
2	D	355	
2	L	355	
3	E	355	
4	F	355	
4	J	355	
4	O	355	
5	G	355	
6	H	355	
7	I	355	
8	M	355	
9	P	355	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
12	PO4	A	407	-	-	X	-
12	PO4	B	409	-	X	-	-
9	SCY	P	150	-	-	X	-

2 Entry composition

There are 20 unique types of molecules in this entry. The entry contains 41155 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 A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	1	0
			2509	1576	435	487	11			
1	B	332	Total	C	N	O	S	0	1	0
			2514	1579	436	488	11			
1	C	331	Total	C	N	O	S	0	0	0
			2499	1570	432	486	11			
1	K	331	Total	C	N	O	S	0	1	0
			2509	1576	435	487	11			
1	N	331	Total	C	N	O	S	0	1	0
			2509	1576	435	487	11			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP C9QTS9
A	-22	HIS	-	expression tag	UNP C9QTS9
A	-21	HIS	-	expression tag	UNP C9QTS9
A	-20	HIS	-	expression tag	UNP C9QTS9
A	-19	HIS	-	expression tag	UNP C9QTS9
A	-18	HIS	-	expression tag	UNP C9QTS9
A	-17	HIS	-	expression tag	UNP C9QTS9
A	-16	SER	-	expression tag	UNP C9QTS9
A	-15	SER	-	expression tag	UNP C9QTS9
A	-14	GLY	-	expression tag	UNP C9QTS9
A	-13	VAL	-	expression tag	UNP C9QTS9
A	-12	ASP	-	expression tag	UNP C9QTS9
A	-11	LEU	-	expression tag	UNP C9QTS9
A	-10	GLY	-	expression tag	UNP C9QTS9
A	-9	THR	-	expression tag	UNP C9QTS9
A	-8	GLU	-	expression tag	UNP C9QTS9
A	-7	ASN	-	expression tag	UNP C9QTS9
A	-6	LEU	-	expression tag	UNP C9QTS9
A	-5	TYR	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	PHE	-	expression tag	UNP C9QTS9
A	-3	GLN	-	expression tag	UNP C9QTS9
A	-2	SER	-	expression tag	UNP C9QTS9
A	-1	ASN	-	expression tag	UNP C9QTS9
A	0	ALA	-	expression tag	UNP C9QTS9
B	-23	MET	-	expression tag	UNP C9QTS9
B	-22	HIS	-	expression tag	UNP C9QTS9
B	-21	HIS	-	expression tag	UNP C9QTS9
B	-20	HIS	-	expression tag	UNP C9QTS9
B	-19	HIS	-	expression tag	UNP C9QTS9
B	-18	HIS	-	expression tag	UNP C9QTS9
B	-17	HIS	-	expression tag	UNP C9QTS9
B	-16	SER	-	expression tag	UNP C9QTS9
B	-15	SER	-	expression tag	UNP C9QTS9
B	-14	GLY	-	expression tag	UNP C9QTS9
B	-13	VAL	-	expression tag	UNP C9QTS9
B	-12	ASP	-	expression tag	UNP C9QTS9
B	-11	LEU	-	expression tag	UNP C9QTS9
B	-10	GLY	-	expression tag	UNP C9QTS9
B	-9	THR	-	expression tag	UNP C9QTS9
B	-8	GLU	-	expression tag	UNP C9QTS9
B	-7	ASN	-	expression tag	UNP C9QTS9
B	-6	LEU	-	expression tag	UNP C9QTS9
B	-5	TYR	-	expression tag	UNP C9QTS9
B	-4	PHE	-	expression tag	UNP C9QTS9
B	-3	GLN	-	expression tag	UNP C9QTS9
B	-2	SER	-	expression tag	UNP C9QTS9
B	-1	ASN	-	expression tag	UNP C9QTS9
B	0	ALA	-	expression tag	UNP C9QTS9
C	-23	MET	-	expression tag	UNP C9QTS9
C	-22	HIS	-	expression tag	UNP C9QTS9
C	-21	HIS	-	expression tag	UNP C9QTS9
C	-20	HIS	-	expression tag	UNP C9QTS9
C	-19	HIS	-	expression tag	UNP C9QTS9
C	-18	HIS	-	expression tag	UNP C9QTS9
C	-17	HIS	-	expression tag	UNP C9QTS9
C	-16	SER	-	expression tag	UNP C9QTS9
C	-15	SER	-	expression tag	UNP C9QTS9
C	-14	GLY	-	expression tag	UNP C9QTS9
C	-13	VAL	-	expression tag	UNP C9QTS9
C	-12	ASP	-	expression tag	UNP C9QTS9
C	-11	LEU	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	GLY	-	expression tag	UNP C9QTS9
C	-9	THR	-	expression tag	UNP C9QTS9
C	-8	GLU	-	expression tag	UNP C9QTS9
C	-7	ASN	-	expression tag	UNP C9QTS9
C	-6	LEU	-	expression tag	UNP C9QTS9
C	-5	TYR	-	expression tag	UNP C9QTS9
C	-4	PHE	-	expression tag	UNP C9QTS9
C	-3	GLN	-	expression tag	UNP C9QTS9
C	-2	SER	-	expression tag	UNP C9QTS9
C	-1	ASN	-	expression tag	UNP C9QTS9
C	0	ALA	-	expression tag	UNP C9QTS9
K	-23	MET	-	expression tag	UNP C9QTS9
K	-22	HIS	-	expression tag	UNP C9QTS9
K	-21	HIS	-	expression tag	UNP C9QTS9
K	-20	HIS	-	expression tag	UNP C9QTS9
K	-19	HIS	-	expression tag	UNP C9QTS9
K	-18	HIS	-	expression tag	UNP C9QTS9
K	-17	HIS	-	expression tag	UNP C9QTS9
K	-16	SER	-	expression tag	UNP C9QTS9
K	-15	SER	-	expression tag	UNP C9QTS9
K	-14	GLY	-	expression tag	UNP C9QTS9
K	-13	VAL	-	expression tag	UNP C9QTS9
K	-12	ASP	-	expression tag	UNP C9QTS9
K	-11	LEU	-	expression tag	UNP C9QTS9
K	-10	GLY	-	expression tag	UNP C9QTS9
K	-9	THR	-	expression tag	UNP C9QTS9
K	-8	GLU	-	expression tag	UNP C9QTS9
K	-7	ASN	-	expression tag	UNP C9QTS9
K	-6	LEU	-	expression tag	UNP C9QTS9
K	-5	TYR	-	expression tag	UNP C9QTS9
K	-4	PHE	-	expression tag	UNP C9QTS9
K	-3	GLN	-	expression tag	UNP C9QTS9
K	-2	SER	-	expression tag	UNP C9QTS9
K	-1	ASN	-	expression tag	UNP C9QTS9
K	0	ALA	-	expression tag	UNP C9QTS9
N	-23	MET	-	expression tag	UNP C9QTS9
N	-22	HIS	-	expression tag	UNP C9QTS9
N	-21	HIS	-	expression tag	UNP C9QTS9
N	-20	HIS	-	expression tag	UNP C9QTS9
N	-19	HIS	-	expression tag	UNP C9QTS9
N	-18	HIS	-	expression tag	UNP C9QTS9
N	-17	HIS	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
N	-16	SER	-	expression tag	UNP C9QTS9
N	-15	SER	-	expression tag	UNP C9QTS9
N	-14	GLY	-	expression tag	UNP C9QTS9
N	-13	VAL	-	expression tag	UNP C9QTS9
N	-12	ASP	-	expression tag	UNP C9QTS9
N	-11	LEU	-	expression tag	UNP C9QTS9
N	-10	GLY	-	expression tag	UNP C9QTS9
N	-9	THR	-	expression tag	UNP C9QTS9
N	-8	GLU	-	expression tag	UNP C9QTS9
N	-7	ASN	-	expression tag	UNP C9QTS9
N	-6	LEU	-	expression tag	UNP C9QTS9
N	-5	TYR	-	expression tag	UNP C9QTS9
N	-4	PHE	-	expression tag	UNP C9QTS9
N	-3	GLN	-	expression tag	UNP C9QTS9
N	-2	SER	-	expression tag	UNP C9QTS9
N	-1	ASN	-	expression tag	UNP C9QTS9
N	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 2 is a protein called Glyceraldehyde-3-phosphate dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	331	Total	C	N	O	S	0	2	0
			2514	1578	437	488	11			
2	L	331	Total	C	N	O	S	0	1	0
			2506	1574	435	486	11			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-23	MET	-	expression tag	UNP C9QTS9
D	-22	HIS	-	expression tag	UNP C9QTS9
D	-21	HIS	-	expression tag	UNP C9QTS9
D	-20	HIS	-	expression tag	UNP C9QTS9
D	-19	HIS	-	expression tag	UNP C9QTS9
D	-18	HIS	-	expression tag	UNP C9QTS9
D	-17	HIS	-	expression tag	UNP C9QTS9
D	-16	SER	-	expression tag	UNP C9QTS9
D	-15	SER	-	expression tag	UNP C9QTS9
D	-14	GLY	-	expression tag	UNP C9QTS9
D	-13	VAL	-	expression tag	UNP C9QTS9
D	-12	ASP	-	expression tag	UNP C9QTS9
D	-11	LEU	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	GLY	-	expression tag	UNP C9QTS9
D	-9	THR	-	expression tag	UNP C9QTS9
D	-8	GLU	-	expression tag	UNP C9QTS9
D	-7	ASN	-	expression tag	UNP C9QTS9
D	-6	LEU	-	expression tag	UNP C9QTS9
D	-5	TYR	-	expression tag	UNP C9QTS9
D	-4	PHE	-	expression tag	UNP C9QTS9
D	-3	GLN	-	expression tag	UNP C9QTS9
D	-2	SER	-	expression tag	UNP C9QTS9
D	-1	ASN	-	expression tag	UNP C9QTS9
D	0	ALA	-	expression tag	UNP C9QTS9
L	-23	MET	-	expression tag	UNP C9QTS9
L	-22	HIS	-	expression tag	UNP C9QTS9
L	-21	HIS	-	expression tag	UNP C9QTS9
L	-20	HIS	-	expression tag	UNP C9QTS9
L	-19	HIS	-	expression tag	UNP C9QTS9
L	-18	HIS	-	expression tag	UNP C9QTS9
L	-17	HIS	-	expression tag	UNP C9QTS9
L	-16	SER	-	expression tag	UNP C9QTS9
L	-15	SER	-	expression tag	UNP C9QTS9
L	-14	GLY	-	expression tag	UNP C9QTS9
L	-13	VAL	-	expression tag	UNP C9QTS9
L	-12	ASP	-	expression tag	UNP C9QTS9
L	-11	LEU	-	expression tag	UNP C9QTS9
L	-10	GLY	-	expression tag	UNP C9QTS9
L	-9	THR	-	expression tag	UNP C9QTS9
L	-8	GLU	-	expression tag	UNP C9QTS9
L	-7	ASN	-	expression tag	UNP C9QTS9
L	-6	LEU	-	expression tag	UNP C9QTS9
L	-5	TYR	-	expression tag	UNP C9QTS9
L	-4	PHE	-	expression tag	UNP C9QTS9
L	-3	GLN	-	expression tag	UNP C9QTS9
L	-2	SER	-	expression tag	UNP C9QTS9
L	-1	ASN	-	expression tag	UNP C9QTS9
L	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 3 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	331	Total	C	N	O	S	0	2	0
			2520	1582	434	493	11			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-23	MET	-	expression tag	UNP C9QTS9
E	-22	HIS	-	expression tag	UNP C9QTS9
E	-21	HIS	-	expression tag	UNP C9QTS9
E	-20	HIS	-	expression tag	UNP C9QTS9
E	-19	HIS	-	expression tag	UNP C9QTS9
E	-18	HIS	-	expression tag	UNP C9QTS9
E	-17	HIS	-	expression tag	UNP C9QTS9
E	-16	SER	-	expression tag	UNP C9QTS9
E	-15	SER	-	expression tag	UNP C9QTS9
E	-14	GLY	-	expression tag	UNP C9QTS9
E	-13	VAL	-	expression tag	UNP C9QTS9
E	-12	ASP	-	expression tag	UNP C9QTS9
E	-11	LEU	-	expression tag	UNP C9QTS9
E	-10	GLY	-	expression tag	UNP C9QTS9
E	-9	THR	-	expression tag	UNP C9QTS9
E	-8	GLU	-	expression tag	UNP C9QTS9
E	-7	ASN	-	expression tag	UNP C9QTS9
E	-6	LEU	-	expression tag	UNP C9QTS9
E	-5	TYR	-	expression tag	UNP C9QTS9
E	-4	PHE	-	expression tag	UNP C9QTS9
E	-3	GLN	-	expression tag	UNP C9QTS9
E	-2	SER	-	expression tag	UNP C9QTS9
E	-1	ASN	-	expression tag	UNP C9QTS9
E	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 4 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	331	Total	C	N	O	S	0	3	0
			2531	1588	440	492	11			
4	J	330	Total	C	N	O	S	0	1	0
			2504	1573	434	487	10			
4	O	331	Total	C	N	O	S	0	0	0
			2502	1572	432	487	11			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-23	MET	-	expression tag	UNP C9QTS9
F	-22	HIS	-	expression tag	UNP C9QTS9
F	-21	HIS	-	expression tag	UNP C9QTS9
F	-20	HIS	-	expression tag	UNP C9QTS9
F	-19	HIS	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-18	HIS	-	expression tag	UNP C9QTS9
F	-17	HIS	-	expression tag	UNP C9QTS9
F	-16	SER	-	expression tag	UNP C9QTS9
F	-15	SER	-	expression tag	UNP C9QTS9
F	-14	GLY	-	expression tag	UNP C9QTS9
F	-13	VAL	-	expression tag	UNP C9QTS9
F	-12	ASP	-	expression tag	UNP C9QTS9
F	-11	LEU	-	expression tag	UNP C9QTS9
F	-10	GLY	-	expression tag	UNP C9QTS9
F	-9	THR	-	expression tag	UNP C9QTS9
F	-8	GLU	-	expression tag	UNP C9QTS9
F	-7	ASN	-	expression tag	UNP C9QTS9
F	-6	LEU	-	expression tag	UNP C9QTS9
F	-5	TYR	-	expression tag	UNP C9QTS9
F	-4	PHE	-	expression tag	UNP C9QTS9
F	-3	GLN	-	expression tag	UNP C9QTS9
F	-2	SER	-	expression tag	UNP C9QTS9
F	-1	ASN	-	expression tag	UNP C9QTS9
F	0	ALA	-	expression tag	UNP C9QTS9
J	-23	MET	-	expression tag	UNP C9QTS9
J	-22	HIS	-	expression tag	UNP C9QTS9
J	-21	HIS	-	expression tag	UNP C9QTS9
J	-20	HIS	-	expression tag	UNP C9QTS9
J	-19	HIS	-	expression tag	UNP C9QTS9
J	-18	HIS	-	expression tag	UNP C9QTS9
J	-17	HIS	-	expression tag	UNP C9QTS9
J	-16	SER	-	expression tag	UNP C9QTS9
J	-15	SER	-	expression tag	UNP C9QTS9
J	-14	GLY	-	expression tag	UNP C9QTS9
J	-13	VAL	-	expression tag	UNP C9QTS9
J	-12	ASP	-	expression tag	UNP C9QTS9
J	-11	LEU	-	expression tag	UNP C9QTS9
J	-10	GLY	-	expression tag	UNP C9QTS9
J	-9	THR	-	expression tag	UNP C9QTS9
J	-8	GLU	-	expression tag	UNP C9QTS9
J	-7	ASN	-	expression tag	UNP C9QTS9
J	-6	LEU	-	expression tag	UNP C9QTS9
J	-5	TYR	-	expression tag	UNP C9QTS9
J	-4	PHE	-	expression tag	UNP C9QTS9
J	-3	GLN	-	expression tag	UNP C9QTS9
J	-2	SER	-	expression tag	UNP C9QTS9
J	-1	ASN	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
J	0	ALA	-	expression tag	UNP C9QTS9
O	-23	MET	-	expression tag	UNP C9QTS9
O	-22	HIS	-	expression tag	UNP C9QTS9
O	-21	HIS	-	expression tag	UNP C9QTS9
O	-20	HIS	-	expression tag	UNP C9QTS9
O	-19	HIS	-	expression tag	UNP C9QTS9
O	-18	HIS	-	expression tag	UNP C9QTS9
O	-17	HIS	-	expression tag	UNP C9QTS9
O	-16	SER	-	expression tag	UNP C9QTS9
O	-15	SER	-	expression tag	UNP C9QTS9
O	-14	GLY	-	expression tag	UNP C9QTS9
O	-13	VAL	-	expression tag	UNP C9QTS9
O	-12	ASP	-	expression tag	UNP C9QTS9
O	-11	LEU	-	expression tag	UNP C9QTS9
O	-10	GLY	-	expression tag	UNP C9QTS9
O	-9	THR	-	expression tag	UNP C9QTS9
O	-8	GLU	-	expression tag	UNP C9QTS9
O	-7	ASN	-	expression tag	UNP C9QTS9
O	-6	LEU	-	expression tag	UNP C9QTS9
O	-5	TYR	-	expression tag	UNP C9QTS9
O	-4	PHE	-	expression tag	UNP C9QTS9
O	-3	GLN	-	expression tag	UNP C9QTS9
O	-2	SER	-	expression tag	UNP C9QTS9
O	-1	ASN	-	expression tag	UNP C9QTS9
O	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 5 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	G	331	Total	C	N	O	S	0	2	0
			2520	1582	437	490	11			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	-23	MET	-	expression tag	UNP C9QTS9
G	-22	HIS	-	expression tag	UNP C9QTS9
G	-21	HIS	-	expression tag	UNP C9QTS9
G	-20	HIS	-	expression tag	UNP C9QTS9
G	-19	HIS	-	expression tag	UNP C9QTS9
G	-18	HIS	-	expression tag	UNP C9QTS9
G	-17	HIS	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-16	SER	-	expression tag	UNP C9QTS9
G	-15	SER	-	expression tag	UNP C9QTS9
G	-14	GLY	-	expression tag	UNP C9QTS9
G	-13	VAL	-	expression tag	UNP C9QTS9
G	-12	ASP	-	expression tag	UNP C9QTS9
G	-11	LEU	-	expression tag	UNP C9QTS9
G	-10	GLY	-	expression tag	UNP C9QTS9
G	-9	THR	-	expression tag	UNP C9QTS9
G	-8	GLU	-	expression tag	UNP C9QTS9
G	-7	ASN	-	expression tag	UNP C9QTS9
G	-6	LEU	-	expression tag	UNP C9QTS9
G	-5	TYR	-	expression tag	UNP C9QTS9
G	-4	PHE	-	expression tag	UNP C9QTS9
G	-3	GLN	-	expression tag	UNP C9QTS9
G	-2	SER	-	expression tag	UNP C9QTS9
G	-1	ASN	-	expression tag	UNP C9QTS9
G	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 6 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	H	331	Total	C	N	O	S	0	0	0
			2505	1574	432	488	11			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	-23	MET	-	expression tag	UNP C9QTS9
H	-22	HIS	-	expression tag	UNP C9QTS9
H	-21	HIS	-	expression tag	UNP C9QTS9
H	-20	HIS	-	expression tag	UNP C9QTS9
H	-19	HIS	-	expression tag	UNP C9QTS9
H	-18	HIS	-	expression tag	UNP C9QTS9
H	-17	HIS	-	expression tag	UNP C9QTS9
H	-16	SER	-	expression tag	UNP C9QTS9
H	-15	SER	-	expression tag	UNP C9QTS9
H	-14	GLY	-	expression tag	UNP C9QTS9
H	-13	VAL	-	expression tag	UNP C9QTS9
H	-12	ASP	-	expression tag	UNP C9QTS9
H	-11	LEU	-	expression tag	UNP C9QTS9
H	-10	GLY	-	expression tag	UNP C9QTS9
H	-9	THR	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-8	GLU	-	expression tag	UNP C9QTS9
H	-7	ASN	-	expression tag	UNP C9QTS9
H	-6	LEU	-	expression tag	UNP C9QTS9
H	-5	TYR	-	expression tag	UNP C9QTS9
H	-4	PHE	-	expression tag	UNP C9QTS9
H	-3	GLN	-	expression tag	UNP C9QTS9
H	-2	SER	-	expression tag	UNP C9QTS9
H	-1	ASN	-	expression tag	UNP C9QTS9
H	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 7 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	I	330	Total	C	N	O	S	0	1	0
			2507	1575	434	488	10			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	-23	MET	-	expression tag	UNP C9QTS9
I	-22	HIS	-	expression tag	UNP C9QTS9
I	-21	HIS	-	expression tag	UNP C9QTS9
I	-20	HIS	-	expression tag	UNP C9QTS9
I	-19	HIS	-	expression tag	UNP C9QTS9
I	-18	HIS	-	expression tag	UNP C9QTS9
I	-17	HIS	-	expression tag	UNP C9QTS9
I	-16	SER	-	expression tag	UNP C9QTS9
I	-15	SER	-	expression tag	UNP C9QTS9
I	-14	GLY	-	expression tag	UNP C9QTS9
I	-13	VAL	-	expression tag	UNP C9QTS9
I	-12	ASP	-	expression tag	UNP C9QTS9
I	-11	LEU	-	expression tag	UNP C9QTS9
I	-10	GLY	-	expression tag	UNP C9QTS9
I	-9	THR	-	expression tag	UNP C9QTS9
I	-8	GLU	-	expression tag	UNP C9QTS9
I	-7	ASN	-	expression tag	UNP C9QTS9
I	-6	LEU	-	expression tag	UNP C9QTS9
I	-5	TYR	-	expression tag	UNP C9QTS9
I	-4	PHE	-	expression tag	UNP C9QTS9
I	-3	GLN	-	expression tag	UNP C9QTS9
I	-2	SER	-	expression tag	UNP C9QTS9
I	-1	ASN	-	expression tag	UNP C9QTS9

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Chain	Residue	Modelled	Actual	Comment	Reference
I	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 8 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	M	331	Total	C	N	O	S	0	1	0
			2508	1575	433	489	11			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	-23	MET	-	expression tag	UNP C9QTS9
M	-22	HIS	-	expression tag	UNP C9QTS9
M	-21	HIS	-	expression tag	UNP C9QTS9
M	-20	HIS	-	expression tag	UNP C9QTS9
M	-19	HIS	-	expression tag	UNP C9QTS9
M	-18	HIS	-	expression tag	UNP C9QTS9
M	-17	HIS	-	expression tag	UNP C9QTS9
M	-16	SER	-	expression tag	UNP C9QTS9
M	-15	SER	-	expression tag	UNP C9QTS9
M	-14	GLY	-	expression tag	UNP C9QTS9
M	-13	VAL	-	expression tag	UNP C9QTS9
M	-12	ASP	-	expression tag	UNP C9QTS9
M	-11	LEU	-	expression tag	UNP C9QTS9
M	-10	GLY	-	expression tag	UNP C9QTS9
M	-9	THR	-	expression tag	UNP C9QTS9
M	-8	GLU	-	expression tag	UNP C9QTS9
M	-7	ASN	-	expression tag	UNP C9QTS9
M	-6	LEU	-	expression tag	UNP C9QTS9
M	-5	TYR	-	expression tag	UNP C9QTS9
M	-4	PHE	-	expression tag	UNP C9QTS9
M	-3	GLN	-	expression tag	UNP C9QTS9
M	-2	SER	-	expression tag	UNP C9QTS9
M	-1	ASN	-	expression tag	UNP C9QTS9
M	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 9 is a protein called Glyceraldehyde-3-phosphate Dehydrogenase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	P	330	Total	C	N	O	S	0	1	0
			2507	1575	434	488	10			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	-23	MET	-	expression tag	UNP C9QTS9
P	-22	HIS	-	expression tag	UNP C9QTS9
P	-21	HIS	-	expression tag	UNP C9QTS9
P	-20	HIS	-	expression tag	UNP C9QTS9
P	-19	HIS	-	expression tag	UNP C9QTS9
P	-18	HIS	-	expression tag	UNP C9QTS9
P	-17	HIS	-	expression tag	UNP C9QTS9
P	-16	SER	-	expression tag	UNP C9QTS9
P	-15	SER	-	expression tag	UNP C9QTS9
P	-14	GLY	-	expression tag	UNP C9QTS9
P	-13	VAL	-	expression tag	UNP C9QTS9
P	-12	ASP	-	expression tag	UNP C9QTS9
P	-11	LEU	-	expression tag	UNP C9QTS9
P	-10	GLY	-	expression tag	UNP C9QTS9
P	-9	THR	-	expression tag	UNP C9QTS9
P	-8	GLU	-	expression tag	UNP C9QTS9
P	-7	ASN	-	expression tag	UNP C9QTS9
P	-6	LEU	-	expression tag	UNP C9QTS9
P	-5	TYR	-	expression tag	UNP C9QTS9
P	-4	PHE	-	expression tag	UNP C9QTS9
P	-3	GLN	-	expression tag	UNP C9QTS9
P	-2	SER	-	expression tag	UNP C9QTS9
P	-1	ASN	-	expression tag	UNP C9QTS9
P	0	ALA	-	expression tag	UNP C9QTS9

- Molecule 10 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	1	Total Na 1 1	0	0
10	B	2	Total Na 2 2	0	0
10	C	1	Total Na 1 1	0	0
10	D	1	Total Na 1 1	0	0
10	E	1	Total Na 1 1	0	0
10	G	1	Total Na 1 1	0	0
10	H	1	Total Na 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	I	1	Total 1	Na 1	0	0
10	K	1	Total 1	Na 1	0	0
10	M	1	Total 1	Na 1	0	0
10	N	1	Total 1	Na 1	0	0

- Molecule 11 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

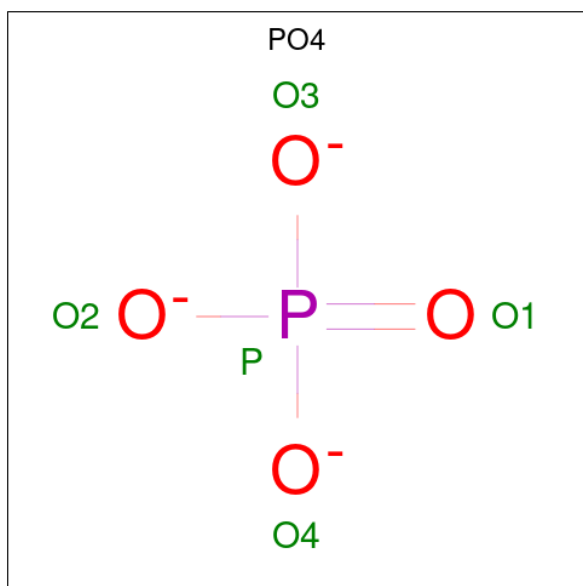
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	4	Total 4	Cl 4	0	0
11	B	6	Total 6	Cl 6	0	0
11	C	5	Total 5	Cl 5	0	0
11	D	5	Total 5	Cl 5	0	0
11	E	5	Total 5	Cl 5	0	0
11	F	1	Total 1	Cl 1	0	0
11	G	4	Total 4	Cl 4	0	0
11	H	1	Total 1	Cl 1	0	0
11	I	2	Total 2	Cl 2	0	0
11	J	2	Total 2	Cl 2	0	0
11	K	6	Total 6	Cl 6	0	0
11	L	1	Total 1	Cl 1	0	0
11	M	2	Total 2	Cl 2	0	0
11	N	1	Total 1	Cl 1	0	0
11	O	2	Total 2	Cl 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	P	2	Total	Cl	0	0
			2	2		

- Molecule 12 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



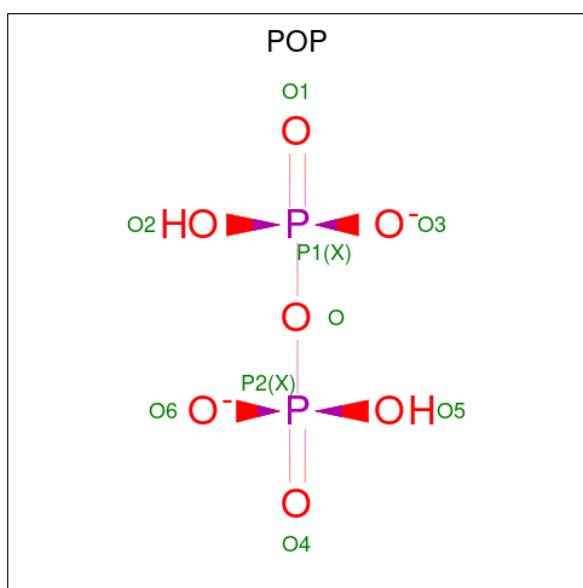
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	A	1	Total	O	P	0	0
			5	4	1		
12	A	1	Total	O	P	0	0
			5	4	1		
12	A	1	Total	O	P	0	0
			5	4	1		
12	B	1	Total	O	P	0	0
			5	4	1		
12	D	1	Total	O	P	0	0
			5	4	1		
12	D	1	Total	O	P	0	0
			5	4	1		
12	E	1	Total	O	P	0	0
			5	4	1		
12	F	1	Total	O	P	0	0
			5	4	1		
12	H	1	Total	O	P	0	0
			5	4	1		
12	I	1	Total	O	P	0	0
			5	4	1		

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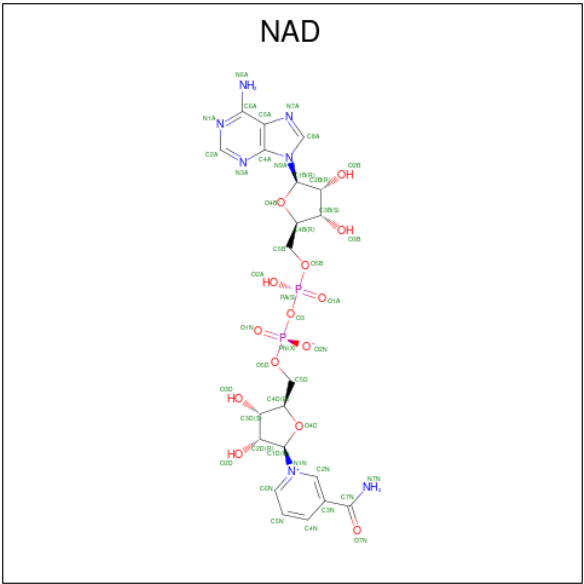
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	J	1	Total	O	P	0	0
			5	4	1		
12	J	1	Total	O	P	0	0
			5	4	1		
12	L	1	Total	O	P	0	0
			5	4	1		
12	P	1	Total	O	P	0	0
			5	4	1		
12	P	1	Total	O	P	0	0
			5	4	1		

- Molecule 13 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



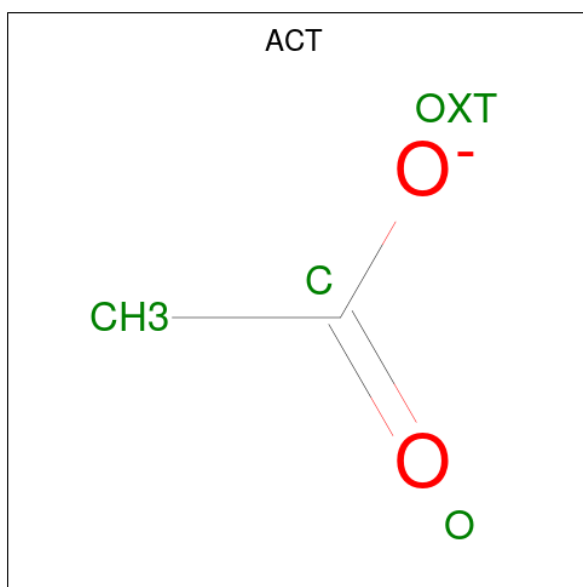
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	A	1	Total	O	P	0	0
			9	7	2		
13	B	1	Total	O	P	0	0
			9	7	2		
13	F	1	Total	O	P	0	0
			9	7	2		
13	G	1	Total	O	P	0	0
			9	7	2		
13	J	1	Total	O	P	0	0
			9	7	2		
13	M	1	Total	O	P	0	0
			9	7	2		

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



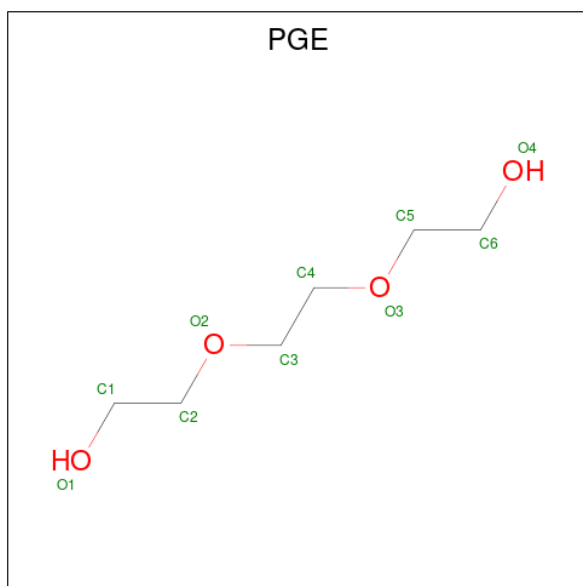
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
14	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
14	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
14	I	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
14	K	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
14	O	1	Total	C	N	O	P	0	0
			44	21	7	14	2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

- Molecule 15 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



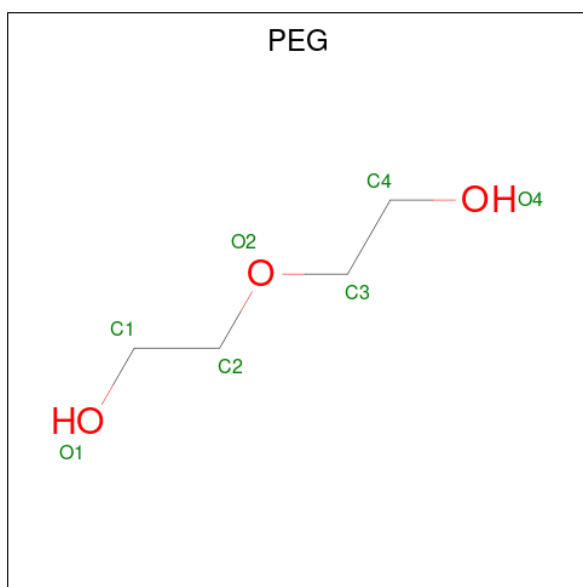
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
15	H	1	Total	C	O	0	0
			4	2	2		

- Molecule 16 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



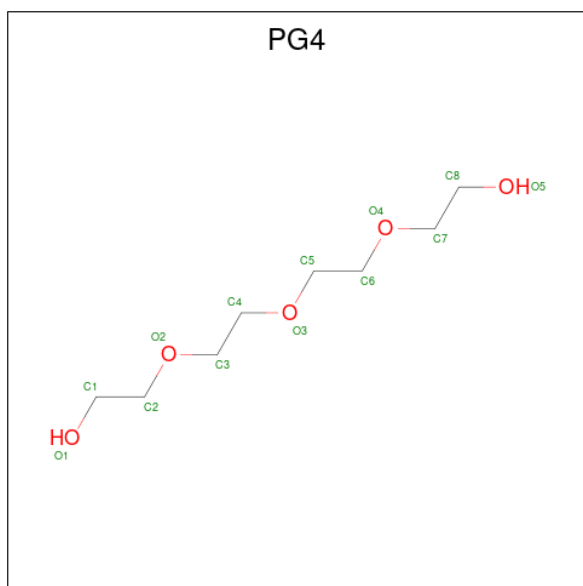
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
16	H	1	Total	C	O	0	0
			10	6	4		

- Molecule 17 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



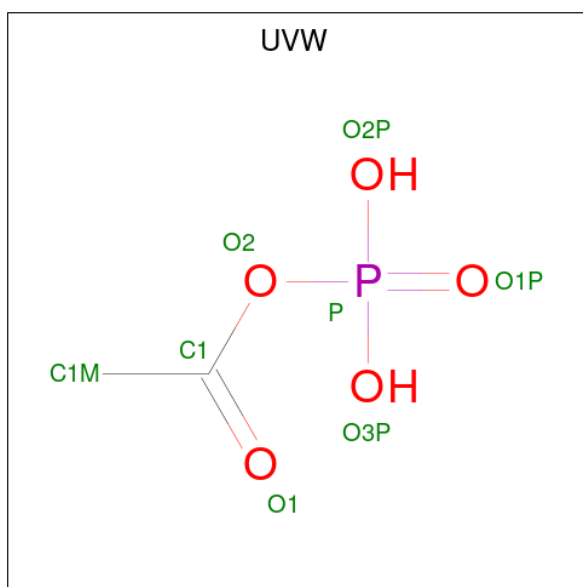
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
17	H	1	Total	C	O	0	0
			7	4	3		

- Molecule 18 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
18	J	1	Total	C	O	0	0
			13	8	5		

- Molecule 19 is ACETYLPHOSPHATE (CCD ID: UVW) (formula: $C_2H_5O_5P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
19	M	1	Total	C	O	P	0	0
			8	2	5	1		

- Molecule 20 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
20	A	39	Total	O	0	0
			39	39		
20	B	35	Total	O	0	3
			38	38		
20	C	61	Total	O	0	0
			61	61		
20	D	37	Total	O	0	1
			38	38		
20	E	23	Total	O	0	0
			23	23		
20	F	27	Total	O	0	0
			27	27		
20	G	33	Total	O	0	2
			35	35		
20	H	21	Total	O	0	1
			22	22		
20	I	22	Total	O	0	1
			23	23		
20	J	28	Total	O	0	0
			28	28		
20	K	26	Total	O	0	0
			26	26		

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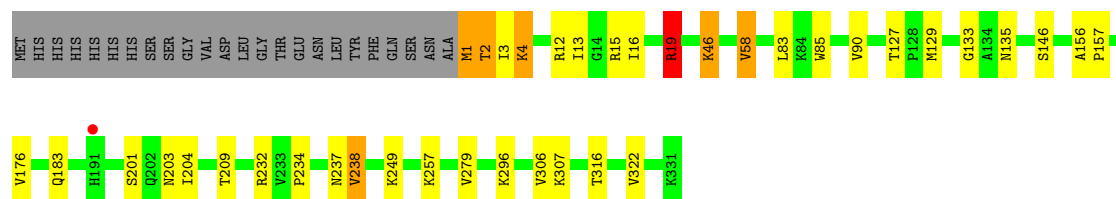
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
20	L	20	Total 20	O 20	0	0
20	M	24	Total 24	O 24	0	0
20	N	11	Total 11	O 11	0	0
20	O	20	Total 20	O 20	0	0
20	P	16	Total 16	O 16	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

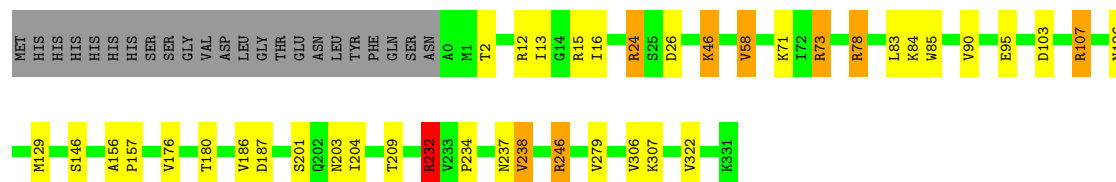
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A

Chain A: 



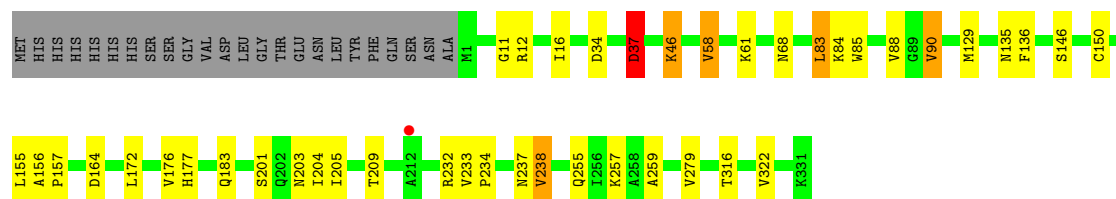
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A

Chain B: 



- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A

Chain C: 



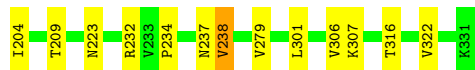
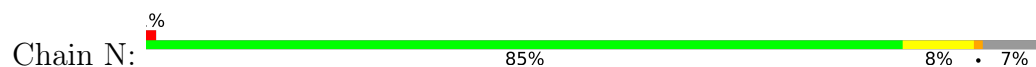
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A

Chain K: 

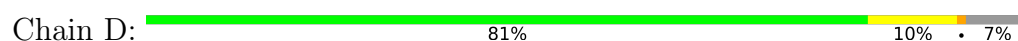




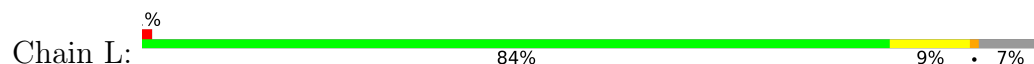
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A



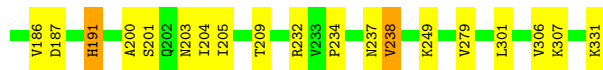
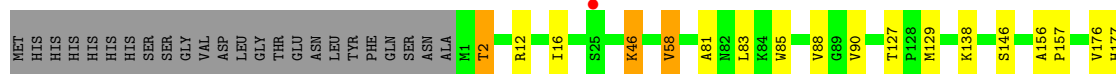
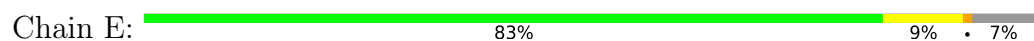
- Molecule 2: Glyceraldehyde-3-phosphate dehydrogenase A



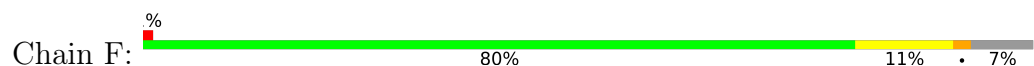
- Molecule 2: Glyceraldehyde-3-phosphate dehydrogenase A

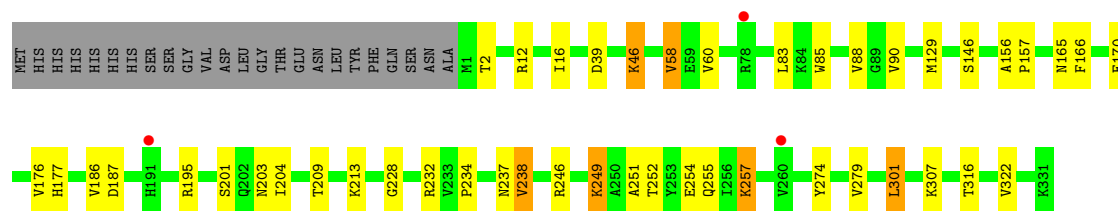


- Molecule 3: Glyceraldehyde-3-phosphate Dehydrogenase A



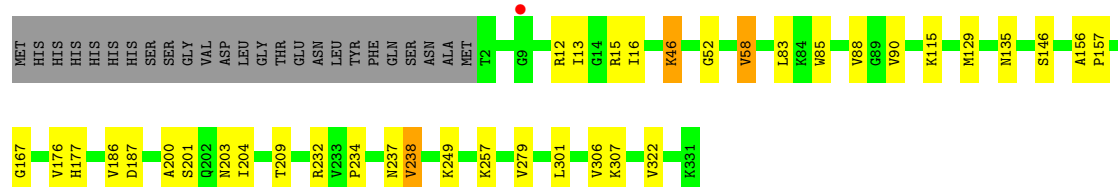
- Molecule 4: Glyceraldehyde-3-phosphate Dehydrogenase A





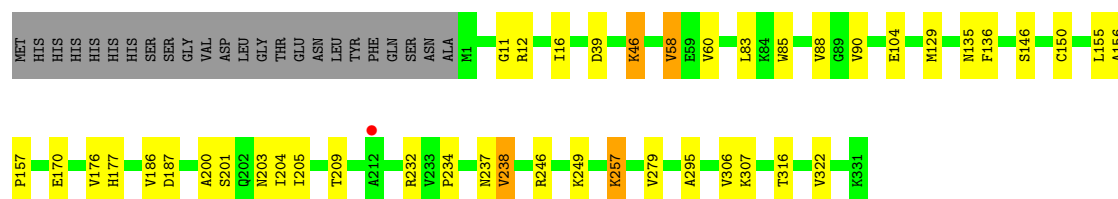
• Molecule 4: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain J: 82% 10% • 7%



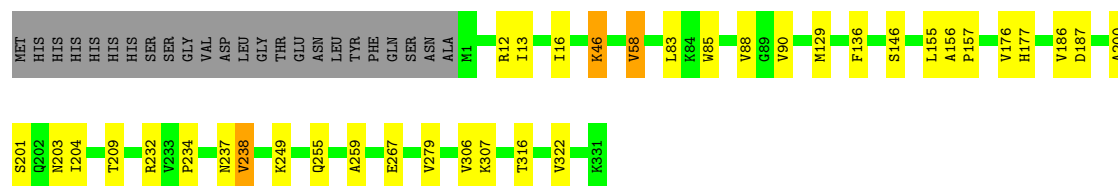
• Molecule 4: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain O: 81% 11% • 7%



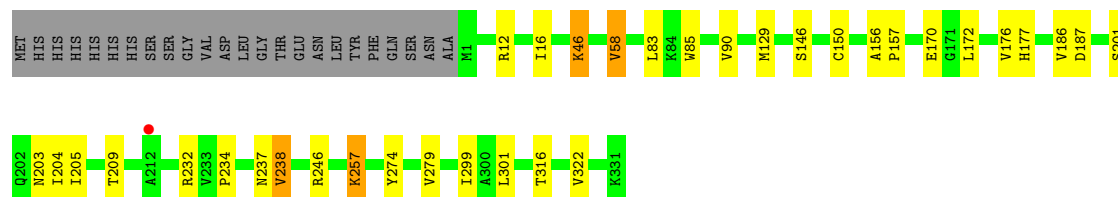
• Molecule 5: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain G: 83% 10% • 7%



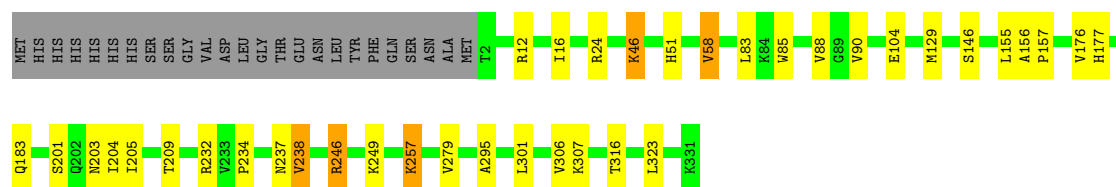
• Molecule 6: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain H: 83% 9% • 7%



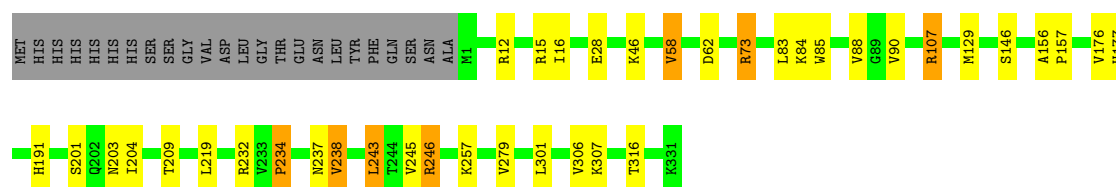
• Molecule 7: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain I:  82% 9% 7%



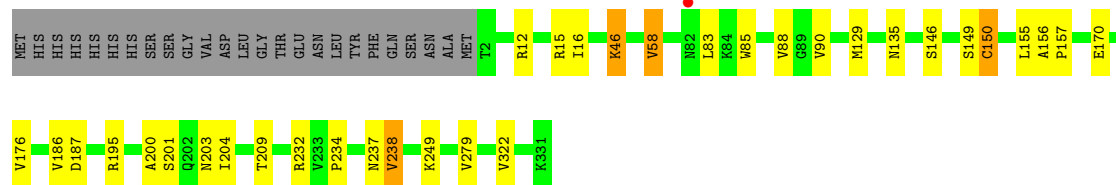
• Molecule 8: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain M:  82% 9% 7%



• Molecule 9: Glyceraldehyde-3-phosphate Dehydrogenase A

Chain P:  83% 8% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	145.88Å 69.69Å 271.92Å 90.00° 98.80° 90.00°	Depositor
Resolution (Å)	29.91 – 2.85 29.91 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.91-2.85) 99.8 (29.91-2.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.85Å)	Xtriage
Refinement program	REFMAC 5.8.0046	Depositor
R, R_{free}	0.186 , 0.224 0.184 , 0.221	Depositor DCC
R_{free} test set	6389 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	64.0	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	41155	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.16% 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: NA, ACT, PEG, ALY, PGE, CL, SCY, PO4, NAD, UVW, POP, PG4

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	0.93	2/2536 (0.1%)	1.49	21/3432 (0.6%)
1	B	0.98	3/2541 (0.1%)	1.42	23/3439 (0.7%)
1	C	0.99	1/2525 (0.0%)	1.11	13/3417 (0.4%)
1	K	0.91	3/2536 (0.1%)	1.30	10/3432 (0.3%)
1	N	0.74	0/2536	1.05	7/3432 (0.2%)
2	D	0.92	2/2554 (0.1%)	1.24	18/3457 (0.5%)
2	L	0.79	0/2546	1.10	12/3446 (0.3%)
3	E	0.88	3/2533 (0.1%)	1.10	11/3427 (0.3%)
4	F	0.81	0/2545	1.07	7/3443 (0.2%)
4	J	0.84	1/2518 (0.0%)	1.06	10/3408 (0.3%)
4	O	0.80	0/2515	1.08	10/3403 (0.3%)
5	G	0.85	0/2537	1.07	8/3432 (0.2%)
6	H	0.83	0/2508	1.07	7/3392 (0.2%)
7	I	0.79	0/2508	1.14	12/3394 (0.4%)
8	M	0.89	1/2534 (0.0%)	1.30	18/3429 (0.5%)
9	P	0.77	0/2511	1.10	15/3397 (0.4%)
All	All	0.86	16/40483 (0.0%)	1.18	202/54780 (0.4%)

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	A	0	1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	24	ARG	CD-NE	-9.15	1.33	1.46
1	K	24	ARG	CZ-NH1	-7.72	1.22	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	19	ARG	CD-NE	-7.70	1.35	1.46
1	B	78	ARG	CA-CB	6.22	1.63	1.53
1	B	107	ARG	CZ-NH1	-5.89	1.24	1.32
1	K	24	ARG	CZ-NH2	5.88	1.41	1.33
1	C	233	VAL	CA-CB	-5.87	1.50	1.55
2	D	62	ASP	CA-CB	5.86	1.62	1.53
3	E	2	THR	CA-CB	5.83	1.61	1.53
3	E	2	THR	N-CA	5.63	1.53	1.46
2	D	22	GLN	CA-C	5.47	1.60	1.52
4	J	249	LYS	C-O	5.40	1.31	1.23
1	B	15	ARG	CZ-NH1	-5.18	1.25	1.32
3	E	2	THR	CB-OG1	5.18	1.52	1.43
1	A	19	ARG	NE-CZ	-5.12	1.27	1.33
8	M	62	ASP	CA-CB	5.03	1.60	1.53

All (202) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	19	ARG	NE-CZ-NH1	-29.64	91.86	121.50
1	K	24	ARG	NE-CZ-NH1	-29.52	91.98	121.50
1	K	24	ARG	NE-CZ-NH2	28.95	145.26	119.20
1	A	19	ARG	NE-CZ-NH2	28.33	144.69	119.20
1	A	15	ARG	NE-CZ-NH2	22.11	139.10	119.20
1	B	15	ARG	NE-CZ-NH2	21.87	138.89	119.20
2	D	15	ARG	NE-CZ-NH2	21.65	138.68	119.20
8	M	15	ARG	NE-CZ-NH2	21.32	138.39	119.20
1	B	15	ARG	NE-CZ-NH1	-20.93	100.57	121.50
1	A	15	ARG	NE-CZ-NH1	-20.80	100.70	121.50
2	D	15	ARG	NE-CZ-NH1	-20.69	100.81	121.50
8	M	15	ARG	NE-CZ-NH1	-20.68	100.82	121.50
1	A	19	ARG	CD-NE-CZ	19.92	152.29	124.40
1	B	107	ARG	NE-CZ-NH1	-18.48	103.02	121.50
1	B	107	ARG	NE-CZ-NH2	17.69	135.12	119.20
1	B	73	ARG	NE-CZ-NH1	-15.35	106.15	121.50
1	A	19	ARG	CG-CD-NE	15.33	145.72	112.00
8	M	73	ARG	NE-CZ-NH1	-14.84	106.66	121.50
8	M	73	ARG	NE-CZ-NH2	14.48	132.24	119.20
8	M	73	ARG	CG-CD-NE	14.47	143.84	112.00
1	B	73	ARG	CG-CD-NE	14.28	143.42	112.00
1	B	73	ARG	NE-CZ-NH2	14.27	132.04	119.20
7	I	246	ARG	NE-CZ-NH2	13.20	131.08	119.20
2	L	329	ILE	CA-CB-CG1	11.63	130.17	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	24	ARG	CD-NE-CZ	11.44	140.42	124.40
1	A	15	ARG	CD-NE-CZ	10.53	139.14	124.40
1	B	73	ARG	CD-NE-CZ	10.51	139.12	124.40
1	B	15	ARG	CD-NE-CZ	10.46	139.05	124.40
7	I	246	ARG	NE-CZ-NH1	-10.46	111.04	121.50
2	D	15	ARG	CD-NE-CZ	10.24	138.74	124.40
8	M	15	ARG	CD-NE-CZ	10.14	138.59	124.40
8	M	73	ARG	CD-NE-CZ	9.96	138.34	124.40
1	A	232	ARG	NE-CZ-NH1	-9.53	111.97	121.50
7	I	246	ARG	CG-CD-NE	9.49	132.89	112.00
9	P	232	ARG	NE-CZ-NH1	-9.14	112.36	121.50
2	D	232	ARG	NE-CZ-NH1	-8.98	112.52	121.50
7	I	246	ARG	CD-NE-CZ	8.98	136.97	124.40
2	L	329	ILE	CB-CG1-CD1	8.92	132.53	113.80
6	H	203	ASN	N-CA-C	8.75	123.09	109.52
1	C	203	ASN	N-CA-C	8.75	123.08	109.52
1	N	203	ASN	N-CA-C	8.73	123.05	109.52
4	F	203	ASN	N-CA-C	8.71	123.02	109.52
4	O	203	ASN	N-CA-C	8.64	122.92	109.52
1	B	232	ARG	NE-CZ-NH1	-8.63	112.87	121.50
1	K	203	ASN	N-CA-C	8.55	122.78	109.52
1	A	203	ASN	N-CA-C	8.52	122.72	109.52
4	O	205	ILE	N-CA-C	8.47	116.67	107.60
2	D	203	ASN	N-CA-C	8.46	122.87	109.50
1	B	203	ASN	N-CA-C	8.44	122.60	109.52
7	I	204	ILE	N-CA-C	-8.44	95.67	107.99
4	J	203	ASN	N-CA-C	8.41	122.55	109.52
9	P	204	ILE	N-CA-C	-8.38	95.76	107.99
3	E	203	ASN	N-CA-C	8.35	122.46	109.52
9	P	203	ASN	N-CA-C	8.27	122.33	109.52
1	B	204	ILE	N-CA-C	-8.20	96.02	107.99
5	G	203	ASN	N-CA-C	8.13	122.12	109.52
8	M	204	ILE	N-CA-C	-7.99	96.33	107.99
5	G	204	ILE	N-CA-C	-7.97	96.36	107.99
8	M	203	ASN	N-CA-C	7.89	121.75	109.52
3	E	205	ILE	N-CA-C	7.86	116.01	107.60
2	L	203	ASN	N-CA-C	7.81	121.62	109.52
1	N	204	ILE	N-CA-C	-7.76	96.66	107.99
3	E	331	LYS	CG-CD-CE	7.65	128.90	111.30
1	K	204	ILE	N-CA-C	-7.63	96.85	107.99
8	M	279	VAL	N-CA-C	7.60	118.34	110.05
1	A	204	ILE	N-CA-C	-7.59	96.90	107.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	O	204	ILE	N-CA-C	-7.58	96.92	107.99
1	A	232	ARG	NE-CZ-NH2	7.55	126.00	119.20
8	M	15	ARG	CG-CD-NE	-7.52	95.45	112.00
2	L	204	ILE	N-CA-C	-7.51	97.02	107.99
1	B	15	ARG	CG-CD-NE	-7.47	95.56	112.00
1	C	205	ILE	N-CA-C	7.46	115.59	107.60
2	D	232	ARG	NE-CZ-NH2	7.42	125.88	119.20
7	I	205	ILE	N-CA-C	7.39	115.51	107.60
4	F	204	ILE	N-CA-C	-7.37	97.23	107.99
7	I	203	ASN	N-CA-C	7.34	120.91	109.52
1	B	232	ARG	NE-CZ-NH2	7.33	125.80	119.20
9	P	232	ARG	NE-CZ-NH2	7.31	125.78	119.20
4	O	279	VAL	N-CA-C	7.31	118.02	110.05
3	E	279	VAL	N-CA-C	7.17	117.86	110.05
1	A	279	VAL	N-CA-C	7.17	117.86	110.05
1	C	176	VAL	N-CA-C	-7.16	96.57	107.73
1	A	232	ARG	CD-NE-CZ	7.13	134.38	124.40
9	P	232	ARG	CD-NE-CZ	7.07	134.30	124.40
4	J	204	ILE	N-CA-C	-7.04	97.72	107.99
1	N	279	VAL	N-CA-C	7.02	117.70	110.05
1	C	204	ILE	N-CA-C	-7.01	97.06	107.51
6	H	204	ILE	N-CA-C	-7.01	97.76	107.99
2	D	204	ILE	N-CA-C	-7.00	97.77	107.99
2	D	28	GLU	CA-CB-CG	6.96	128.03	114.10
6	H	205	ILE	N-CA-C	6.96	115.66	107.73
2	D	279	VAL	N-CA-C	6.92	117.59	110.05
8	M	176	VAL	N-CA-C	-6.91	96.95	107.73
1	B	24	ARG	NE-CZ-NH2	6.74	125.27	119.20
7	I	279	VAL	N-CA-C	6.73	117.39	110.05
4	J	279	VAL	N-CA-C	6.73	117.38	110.05
3	E	204	ILE	N-CA-C	-6.70	97.53	107.51
2	L	322	VAL	N-CA-C	-6.57	103.92	110.62
1	K	176	VAL	N-CA-C	-6.57	97.49	107.73
2	L	329	ILE	CG1-CB-CG2	-6.56	91.02	110.70
4	F	316	THR	N-CA-C	6.55	118.22	111.14
1	B	279	VAL	N-CA-C	6.52	117.16	110.05
4	J	176	VAL	N-CA-C	-6.48	97.62	107.73
2	D	316	THR	N-CA-C	6.46	118.11	111.14
1	B	322	VAL	N-CA-C	-6.40	104.27	110.42
9	P	195	ARG	NE-CZ-NH1	-6.40	115.10	121.50
9	P	176	VAL	N-CA-C	-6.39	97.76	107.73
2	L	279	VAL	N-CA-C	6.33	116.95	110.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	M	28	GLU	CB-CG-CD	-6.31	101.87	112.60
2	D	71	LYS	CA-CB-CG	-6.29	101.52	114.10
5	G	176	VAL	N-CA-C	-6.25	97.98	107.73
4	O	176	VAL	N-CA-C	-6.23	98.02	107.73
2	L	176	VAL	N-CA-C	-6.22	98.03	107.73
4	F	279	VAL	N-CA-C	6.19	116.80	110.05
1	B	24	ARG	NE-CZ-NH1	-6.18	115.32	121.50
4	F	176	VAL	N-CA-C	-6.16	98.12	107.73
1	B	71	LYS	CD-CE-NZ	6.14	131.54	111.90
3	E	176	VAL	N-CA-C	-6.14	98.15	107.73
5	G	316	THR	N-CA-C	6.00	117.61	111.14
5	G	322	VAL	N-CA-C	-5.97	104.53	110.62
4	O	155	LEU	N-CA-C	5.97	117.86	111.36
5	G	279	VAL	N-CA-C	5.96	116.55	110.05
4	O	200	ALA	N-CA-C	5.96	117.78	111.28
3	E	138	LYS	CB-CG-CD	5.95	124.99	111.30
1	K	322	VAL	N-CA-C	-5.95	104.70	110.42
4	F	322	VAL	N-CA-C	-5.94	104.71	110.42
1	B	126	ASN	CB-CG-ND2	-5.89	107.56	116.40
9	P	195	ARG	CB-CG-CD	-5.89	97.74	111.30
2	D	176	VAL	N-CA-C	-5.89	98.54	107.73
1	B	2	THR	N-CA-C	-5.89	100.96	110.32
7	I	176	VAL	N-CA-C	-5.89	98.54	107.73
3	E	127	THR	N-CA-C	-5.88	101.01	109.48
6	H	176	VAL	N-CA-C	-5.87	98.58	107.73
4	J	115	LYS	CG-CD-CE	5.83	124.71	111.30
1	C	155	LEU	N-CA-C	5.81	117.69	111.36
1	N	155	LEU	N-CA-C	5.81	117.42	111.14
9	P	279	VAL	N-CA-C	5.79	116.36	110.05
1	C	322	VAL	N-CA-C	-5.78	104.87	110.42
1	N	176	VAL	N-CA-C	-5.76	98.75	107.73
1	A	15	ARG	CG-CD-NE	-5.74	99.37	112.00
1	A	127	THR	N-CA-C	-5.74	101.22	109.48
2	D	15	ARG	CG-CD-NE	-5.73	99.40	112.00
1	N	316	THR	N-CA-C	5.68	117.28	111.14
2	D	234	PRO	N-CA-C	5.66	120.12	112.48
2	L	234	PRO	N-CA-C	5.62	120.07	112.48
4	J	322	VAL	N-CA-C	-5.61	105.03	110.42
7	I	155	LEU	N-CA-C	5.57	117.43	111.36
7	I	51	HIS	N-CA-C	5.57	119.07	112.72
5	G	200	ALA	N-CA-C	5.56	117.34	111.28
6	H	322	VAL	N-CA-C	-5.56	104.95	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	135	ASN	N-CA-C	5.56	119.31	112.03
2	L	155	LEU	N-CA-C	5.56	117.42	111.36
4	J	249	LYS	N-CA-C	-5.55	103.58	110.41
7	I	316	THR	N-CA-C	5.54	117.13	111.14
1	A	322	VAL	N-CA-C	-5.53	105.11	110.42
6	H	279	VAL	N-CA-C	5.53	116.08	110.05
1	C	279	VAL	N-CA-C	5.50	116.05	110.05
8	M	234	PRO	N-CA-C	5.50	119.90	112.48
4	F	301	LEU	N-CA-C	-5.48	104.95	111.69
8	M	316	THR	N-CA-C	5.47	117.05	111.14
4	O	322	VAL	N-CA-C	-5.47	105.04	110.62
3	E	301	LEU	N-CA-C	-5.44	105.00	111.69
5	G	155	LEU	N-CA-C	5.42	117.27	111.36
1	A	249	LYS	CB-CA-C	-5.42	100.92	109.80
1	C	257	LYS	CG-CD-CE	5.41	123.75	111.30
2	D	22	GLN	CB-CG-CD	5.38	121.75	112.60
9	P	195	ARG	NE-CZ-NH2	5.38	124.04	119.20
1	C	316	THR	N-CA-C	5.34	116.91	111.14
4	O	316	THR	N-CA-C	5.33	116.89	111.14
3	E	191	HIS	N-CA-C	5.31	117.75	111.33
3	E	200	ALA	N-CA-C	5.30	117.06	111.28
1	A	3	ILE	N-CA-C	5.30	116.34	108.23
1	K	135	ASN	N-CA-C	5.29	118.97	112.03
4	J	135	ASN	N-CA-C	5.29	118.96	112.03
2	D	246	ARG	CG-CD-NE	5.29	123.63	112.00
4	J	15	ARG	NE-CZ-NH1	5.28	126.78	121.50
9	P	200	ALA	N-CA-C	5.27	117.02	111.28
2	D	155	LEU	N-CA-C	5.25	117.09	111.36
1	K	200	ALA	N-CA-C	5.25	117.01	111.28
4	O	135	ASN	N-CA-C	5.25	118.91	112.03
1	K	155	LEU	N-CA-C	5.24	117.07	111.36
1	N	322	VAL	N-CA-C	-5.23	105.39	110.42
1	C	83	LEU	N-CA-C	5.23	118.76	112.38
2	L	316	THR	N-CA-C	5.22	116.78	111.14
1	B	246	ARG	CD-NE-CZ	-5.21	117.10	124.40
8	M	246	ARG	CG-CD-NE	5.17	123.38	112.00
1	A	176	VAL	N-CA-C	-5.17	99.66	107.73
9	P	155	LEU	N-CA-C	5.17	116.72	111.14
2	D	135	ASN	N-CA-C	5.14	118.77	112.03
1	C	37	ASP	N-CA-CB	5.13	117.70	110.36
1	A	316	THR	N-CA-C	5.12	116.67	111.14
1	B	176	VAL	N-CA-C	-5.12	99.74	107.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	VAL	N-CA-C	5.09	115.66	109.30
9	P	135	ASN	N-CA-C	5.08	118.68	112.03
6	H	316	THR	N-CA-C	5.07	116.61	111.14
9	P	322	VAL	N-CA-C	-5.06	105.56	110.42
1	A	135	ASN	N-CA-C	5.02	118.61	112.03
4	J	200	ALA	N-CA-C	5.02	116.75	111.28
8	M	219	LEU	CA-C-N	5.01	125.04	119.32
8	M	219	LEU	C-N-CA	5.01	125.04	119.32
2	L	281	SER	N-CA-C	5.01	117.40	111.33
9	P	15	ARG	NE-CZ-NH1	5.01	126.51	121.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	19	ARG	Sidechain

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	2509	0	2522	21	0
1	B	2514	0	2527	22	0
1	C	2499	0	2516	25	0
1	K	2509	0	2522	26	0
1	N	2509	0	2522	14	0
2	D	2514	0	2526	23	0
2	L	2506	0	2521	15	0
3	E	2520	0	2527	14	0
4	F	2531	0	2538	30	0
4	J	2504	0	2511	17	0
4	O	2502	0	2517	25	0
5	G	2520	0	2529	20	0
6	H	2505	0	2519	20	0
7	I	2507	0	2511	22	0
8	M	2508	0	2521	19	0
9	P	2507	0	2513	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	1	0	0	0	0
10	B	2	0	0	0	0
10	C	1	0	0	0	0
10	D	1	0	0	0	0
10	E	1	0	0	0	0
10	G	1	0	0	0	0
10	H	1	0	0	0	0
10	I	1	0	0	0	0
10	K	1	0	0	0	0
10	M	1	0	0	0	0
10	N	1	0	0	0	0
11	A	4	0	0	2	0
11	B	6	0	0	0	0
11	C	5	0	0	1	0
11	D	5	0	0	1	0
11	E	5	0	0	0	0
11	F	1	0	0	0	0
11	G	4	0	0	1	0
11	H	1	0	0	0	0
11	I	2	0	0	0	0
11	J	2	0	0	0	0
11	K	6	0	0	0	0
11	L	1	0	0	0	0
11	M	2	0	0	0	0
11	N	1	0	0	0	0
11	O	2	0	0	1	0
11	P	2	0	0	0	0
12	A	15	0	0	2	0
12	B	5	0	0	0	0
12	D	10	0	0	2	0
12	E	5	0	0	0	0
12	F	5	0	0	0	0
12	H	5	0	0	0	0
12	I	5	0	0	0	0
12	J	10	0	0	0	0
12	L	5	0	0	0	0
12	P	10	0	0	0	0
13	A	9	0	0	1	0
13	B	9	0	0	2	0
13	F	9	0	0	0	0
13	G	9	0	0	1	0
13	J	9	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	9	0	0	0	0
14	C	44	0	26	4	0
14	D	44	0	26	3	0
14	E	44	0	26	0	0
14	I	44	0	26	0	0
14	K	44	0	26	4	0
14	O	44	0	26	4	0
14	P	44	0	26	2	0
15	H	4	0	3	0	0
16	H	10	0	14	0	0
17	H	7	0	10	0	0
18	J	13	0	18	0	0
19	M	8	0	3	0	0
20	A	39	0	0	1	0
20	B	38	0	0	1	0
20	C	61	0	0	3	0
20	D	38	0	0	1	0
20	E	23	0	0	1	0
20	F	27	0	0	2	0
20	G	35	0	0	2	0
20	H	22	0	0	0	0
20	I	23	0	0	1	0
20	J	28	0	0	0	0
20	K	26	0	0	3	0
20	L	20	0	0	1	0
20	M	24	0	0	0	0
20	N	11	0	0	1	0
20	O	20	0	0	2	0
20	P	16	0	0	0	0
All	All	41155	0	40572	304	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (304) 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)
7:I:104:GLU:OE2	4:O:104:GLU:OE2	1.85	0.94
1:K:71:LYS:HE3	20:K:522:HOH:O	1.76	0.83
8:M:243:LEU:HD22	8:M:245:VAL:HG13	1.61	0.82
1:K:254:GLU:HB2	20:K:511:HOH:O	1.80	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:GLN:HB3	20:A:512:HOH:O	1.83	0.79
4:F:166:PHE:HA	4:F:249:LYS:HG2	1.65	0.77
7:I:24:ARG:HH11	7:I:323:LEU:HB3	1.50	0.74
1:B:246:ARG:HD3	2:D:246:ARG:NH1	2.05	0.71
7:I:246:ARG:NE	1:K:246:ARG:HD2	2.06	0.71
1:B:13:ILE:N	13:B:410:POP:O4	2.24	0.69
8:M:243:LEU:CD2	8:M:245:VAL:HG13	2.24	0.67
7:I:183:GLN:HB3	20:I:503:HOH:O	1.93	0.67
9:P:150:SCY:HE2	14:P:401:NAD:C5N	2.25	0.66
1:A:1:MET:HA	1:A:1:MET:CE	2.29	0.63
1:B:83:LEU:HD13	1:B:85:TRP:CZ2	2.34	0.62
2:D:12:ARG:N	14:D:401:NAD:O2A	2.23	0.62
8:M:83:LEU:HD13	8:M:85:TRP:CZ2	2.35	0.62
9:P:83:LEU:HD13	9:P:85:TRP:CZ2	2.35	0.62
2:D:83:LEU:HD13	2:D:85:TRP:CZ2	2.35	0.61
4:J:13:ILE:HG12	13:J:406:POP:O6	2.01	0.61
4:J:83:LEU:HD13	4:J:85:TRP:CZ2	2.36	0.61
4:F:83:LEU:HD13	4:F:85:TRP:CZ2	2.35	0.61
5:G:83:LEU:HD13	5:G:85:TRP:CZ2	2.36	0.61
1:B:246:ARG:HD3	2:D:246:ARG:HH11	1.66	0.61
6:H:257:ALY:HG3	6:H:274:TYR:OH	2.01	0.60
2:L:83:LEU:HD13	2:L:85:TRP:CZ2	2.36	0.60
3:E:83:LEU:HD13	3:E:85:TRP:CZ2	2.36	0.60
6:H:83:LEU:HD13	6:H:85:TRP:CZ2	2.36	0.60
7:I:83:LEU:HD13	7:I:85:TRP:CZ2	2.36	0.60
4:O:83:LEU:HD13	4:O:85:TRP:CZ2	2.37	0.60
1:A:83:LEU:HD13	1:A:85:TRP:CZ2	2.36	0.60
1:C:183:GLN:HB3	20:C:514:HOH:O	2.02	0.60
1:K:83:LEU:HD13	1:K:85:TRP:CZ2	2.37	0.60
2:D:43:TYR:HD1	12:D:409:PO4:O3	1.85	0.59
1:C:83:LEU:HD13	1:C:85:TRP:CZ2	2.38	0.59
1:N:83:LEU:HD13	1:N:85:TRP:CZ2	2.37	0.59
4:F:251:ALA:HA	4:F:255:GLN:OE1	2.03	0.58
1:A:257:LYS:NZ	12:A:407:PO4:O1	2.36	0.58
4:F:257:ALY:HG3	4:F:274:TYR:OH	2.04	0.58
1:A:13:ILE:HG12	13:A:409:POP:O3	2.05	0.57
7:I:24:ARG:NH1	7:I:323:LEU:HB3	2.18	0.57
5:G:13:ILE:HG12	13:G:406:POP:O5	2.04	0.57
8:M:73:ARG:NH1	8:M:84:LYS:O	2.37	0.56
5:G:267:GLU:HB2	20:G:518:HOH:O	2.05	0.56
1:A:2:THR:O	1:A:4:LYS:NZ	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:ARG:NH1	1:B:84:LYS:O	2.39	0.55
1:K:120:THR:O	14:K:401:NAD:H1D	2.06	0.55
1:K:177:HIS:HB3	1:K:232:ARG:HD3	1.88	0.55
4:F:166:PHE:HA	4:F:249:LYS:CG	2.37	0.55
1:B:24:ARG:NH2	1:B:26:ASP:OD2	2.41	0.54
4:J:177:HIS:HB3	4:J:232:ARG:HD3	1.89	0.54
2:D:71:LYS:HG2	2:D:72:ILE:N	2.21	0.54
1:A:19:ARG:HD3	11:A:405:CL:CL	2.45	0.53
1:A:296:LYS:HA	12:A:407:PO4:O3	2.08	0.53
8:M:191:HIS:H	8:M:191:HIS:CD2	2.24	0.53
2:L:177:HIS:HB3	2:L:232:ARG:HD3	1.91	0.53
7:I:246:ARG:CD	1:K:246:ARG:HD2	2.38	0.53
9:P:150:SCY:H	9:P:150:SCY:CD	2.19	0.53
2:D:46:LYS:HG2	2:D:58:VAL:CG1	2.39	0.53
1:K:237:ASN:OD1	1:K:314:ASN:OD1	2.27	0.52
1:N:177:HIS:HB3	1:N:232:ARG:HD3	1.91	0.52
1:K:150:CYS:SG	14:K:401:NAD:H4N	2.49	0.52
1:A:1:MET:HA	1:A:1:MET:HE3	1.91	0.52
8:M:107:ARG:O	8:M:107:ARG:HG3	2.09	0.52
4:F:301:LEU:CD2	6:H:170:GLU:HB2	2.40	0.52
7:I:177:HIS:HB3	7:I:232:ARG:HD3	1.90	0.52
4:F:177:HIS:HB3	4:F:232:ARG:HD3	1.91	0.52
8:M:177:HIS:HB3	8:M:232:ARG:HD3	1.92	0.52
3:E:177:HIS:HB3	3:E:232:ARG:HD3	1.91	0.51
1:C:177:HIS:HB3	1:C:232:ARG:HD3	1.91	0.51
1:C:164:ASP:O	5:G:249:LYS:NZ	2.41	0.51
8:M:246:ARG:NE	4:O:246:ARG:HD2	2.25	0.51
4:O:177:HIS:HB3	4:O:232:ARG:HD3	1.91	0.51
6:H:177:HIS:HB3	6:H:232:ARG:HD3	1.91	0.51
9:P:149:SER:HB2	9:P:150:SCY:OCD	2.10	0.51
3:E:81:ALA:HA	20:E:506:HOH:O	2.12	0.50
4:O:306:VAL:HG12	4:O:307:LYS:N	2.27	0.50
4:J:12:ARG:O	4:J:16:ILE:HG12	2.12	0.50
1:K:12:ARG:O	1:K:16:ILE:HG12	2.12	0.50
1:C:12:ARG:O	1:C:16:ILE:HG12	2.12	0.50
1:C:34:ASP:OD1	14:C:401:NAD:H1B	2.11	0.50
1:A:12:ARG:O	1:A:16:ILE:HG12	2.12	0.49
5:G:12:ARG:O	5:G:16:ILE:HG12	2.12	0.49
5:G:177:HIS:HB3	5:G:232:ARG:HD3	1.94	0.49
5:G:46:ALY:HG3	5:G:58:VAL:CG1	2.43	0.49
1:K:46:ALY:HG3	1:K:58:VAL:CG1	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:TRP:HB3	1:B:90:VAL:HB	1.95	0.49
3:E:12:ARG:O	3:E:16:ILE:HG12	2.12	0.49
4:F:12:ARG:O	4:F:16:ILE:HG12	2.12	0.49
6:H:12:ARG:O	6:H:16:ILE:HG12	2.12	0.49
2:D:129:MET:HG2	2:D:146:SER:HB3	1.95	0.49
4:F:165:ASN:O	4:F:249:LYS:HD2	2.12	0.49
2:L:85:TRP:HB3	2:L:90:VAL:HB	1.95	0.49
1:A:85:TRP:HB3	1:A:90:VAL:HB	1.94	0.49
5:G:85:TRP:HB3	5:G:90:VAL:HB	1.95	0.49
8:M:12:ARG:O	8:M:16:ILE:HG12	2.13	0.49
9:P:85:TRP:HB3	9:P:90:VAL:HB	1.95	0.49
6:H:46:ALY:HG3	6:H:58:VAL:CG1	2.42	0.49
2:L:12:ARG:O	2:L:16:ILE:HG12	2.13	0.49
1:N:46:ALY:HG3	1:N:58:VAL:CG1	2.43	0.49
4:O:136:PHE:HB2	11:O:403:CL:CL	2.50	0.49
7:I:156:ALA:HB3	7:I:157:PRO:HD3	1.95	0.49
9:P:150:SCY:HE2	14:P:401:NAD:C6N	2.42	0.49
1:C:156:ALA:HB3	1:C:157:PRO:HD3	1.95	0.48
1:C:11:GLY:HA3	14:C:401:NAD:O5B	2.12	0.48
1:K:156:ALA:HB3	1:K:157:PRO:HD3	1.95	0.48
5:G:267:GLU:CB	20:G:518:HOH:O	2.61	0.48
1:N:12:ARG:O	1:N:16:ILE:HG12	2.12	0.48
1:B:12:ARG:O	1:B:16:ILE:HG12	2.14	0.48
5:G:136:PHE:HB2	11:G:403:CL:CL	2.51	0.48
7:I:12:ARG:O	7:I:16:ILE:HG12	2.13	0.48
2:D:12:ARG:O	2:D:16:ILE:HG12	2.12	0.48
2:L:129:MET:HG2	2:L:146:SER:HB3	1.96	0.48
3:E:85:TRP:HB3	3:E:90:VAL:HB	1.95	0.48
4:F:246:ARG:HD2	6:H:246:ARG:CD	2.44	0.48
4:J:46:ALY:HG3	4:J:58:VAL:CG1	2.44	0.48
1:N:85:TRP:HB3	1:N:90:VAL:HB	1.96	0.48
6:H:156:ALA:HB3	6:H:157:PRO:HD3	1.96	0.48
7:I:85:TRP:HB3	7:I:90:VAL:HB	1.95	0.48
4:O:12:ARG:O	4:O:16:ILE:HG12	2.13	0.48
9:P:12:ARG:O	9:P:16:ILE:HG12	2.12	0.48
2:D:170:GLU:OE2	2:D:246:ARG:NH1	2.47	0.48
1:B:12:ARG:N	13:B:410:POP:O4	2.48	0.47
2:D:156:ALA:HB3	2:D:157:PRO:HD3	1.96	0.47
4:J:301:LEU:CD2	2:L:170:GLU:HB2	2.43	0.47
8:M:46:LYS:HG2	8:M:58:VAL:CG1	2.43	0.47
6:H:129:MET:HG2	6:H:146:SER:HB3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:O:11:GLY:HA3	14:O:401:NAD:O5B	2.14	0.47
6:H:85:TRP:HB3	6:H:90:VAL:HB	1.96	0.47
3:E:156:ALA:HB3	3:E:157:PRO:HD3	1.96	0.47
4:F:85:TRP:HB3	4:F:90:VAL:HB	1.97	0.47
5:G:156:ALA:HB3	5:G:157:PRO:HD3	1.96	0.47
9:P:129:MET:HG2	9:P:146:SER:HB3	1.96	0.47
2:D:150:CYS:SG	14:D:401:NAD:H4N	2.54	0.47
4:J:52:GLY:HA2	20:K:526:HOH:O	2.14	0.47
1:A:156:ALA:HB3	1:A:157:PRO:HD3	1.97	0.47
1:A:306:VAL:HG12	1:A:307:LYS:N	2.30	0.47
1:C:129:MET:HG2	1:C:146:SER:HB3	1.97	0.47
2:D:136:PHE:HB2	11:D:406:CL:CL	2.51	0.47
4:J:85:TRP:HB3	4:J:90:VAL:HB	1.95	0.47
8:M:85:TRP:HB3	8:M:90:VAL:HB	1.96	0.47
8:M:306:VAL:HG12	8:M:307:LYS:N	2.29	0.47
4:O:46:ALY:HG3	4:O:58:VAL:CG1	2.45	0.47
4:F:195:ARG:HG3	20:F:502:HOH:O	2.15	0.47
4:F:252:THR:O	4:F:255:GLN:HB2	2.15	0.47
4:F:301:LEU:HD23	6:H:170:GLU:HB2	1.97	0.47
5:G:129:MET:HG2	5:G:146:SER:HB3	1.96	0.47
4:J:129:MET:HG2	4:J:146:SER:HB3	1.97	0.47
4:O:156:ALA:HB3	4:O:157:PRO:HD3	1.97	0.47
1:A:129:MET:HG2	1:A:146:SER:HB3	1.97	0.46
1:B:129:MET:HG2	1:B:146:SER:HB3	1.97	0.46
4:F:246:ARG:CD	6:H:246:ARG:HD2	2.45	0.46
1:N:129:MET:HG2	1:N:146:SER:HB3	1.98	0.46
1:B:95:GLU:HG3	20:B:501:HOH:O	2.14	0.46
6:H:201:SER:HA	6:H:234:PRO:HB3	1.97	0.46
2:D:85:TRP:HB3	2:D:90:VAL:HB	1.97	0.46
4:F:307:LYS:HE2	6:H:172:LEU:HB3	1.97	0.46
7:I:129:MET:HG2	7:I:146:SER:HB3	1.98	0.46
8:M:129:MET:HG2	8:M:146:SER:HB3	1.96	0.46
4:F:201:SER:HA	4:F:234:PRO:HB3	1.97	0.46
1:C:259:ALA:HA	5:G:255:GLN:NE2	2.31	0.46
3:E:46:ALY:HG3	3:E:58:VAL:CG1	2.46	0.46
3:E:129:MET:HG2	3:E:146:SER:HB3	1.97	0.46
4:F:129:MET:HG2	4:F:146:SER:HB3	1.97	0.46
7:I:46:ALY:HG3	7:I:58:VAL:CG1	2.46	0.46
2:D:150:CYS:SG	14:D:401:NAD:C4N	3.04	0.46
1:N:156:ALA:HB3	1:N:157:PRO:HD3	1.97	0.46
7:I:246:ARG:HD3	1:K:246:ARG:HD2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:201:SER:HA	1:K:234:PRO:HB3	1.97	0.46
4:O:85:TRP:HB3	4:O:90:VAL:HB	1.96	0.46
9:P:46:ALY:HG3	9:P:58:VAL:CG1	2.46	0.46
1:C:85:TRP:HB3	1:C:90:VAL:HB	1.98	0.46
4:J:201:SER:HA	4:J:234:PRO:HB3	1.98	0.46
1:K:85:TRP:HB3	1:K:90:VAL:HB	1.96	0.46
2:L:201:SER:HA	2:L:234:PRO:HB3	1.98	0.46
4:F:46:ALY:HG3	4:F:58:VAL:CG1	2.46	0.46
1:A:201:SER:HA	1:A:234:PRO:HB3	1.98	0.45
4:F:237:ASN:O	4:F:238:VAL:HB	2.16	0.45
4:F:246:ARG:HD2	6:H:246:ARG:HD2	1.98	0.45
1:B:46:ALY:HG3	1:B:58:VAL:CG1	2.46	0.45
4:J:156:ALA:HB3	4:J:157:PRO:HD3	1.97	0.45
1:A:46:ALY:HG3	1:A:58:VAL:CG1	2.46	0.45
1:C:46:ALY:HG3	1:C:58:VAL:CG1	2.47	0.45
1:A:1:MET:HA	1:A:1:MET:HE2	1.98	0.45
1:C:136:PHE:HB2	11:C:405:CL:CL	2.53	0.45
7:I:301:LEU:CD2	1:K:170:GLU:HB2	2.47	0.45
1:K:129:MET:HG2	1:K:146:SER:HB3	1.99	0.45
4:O:201:SER:HA	4:O:234:PRO:HB3	1.98	0.45
1:B:24:ARG:NH2	1:B:26:ASP:OD1	2.48	0.45
1:K:237:ASN:O	1:K:238:VAL:HB	2.17	0.45
8:M:301:LEU:CD2	4:O:170:GLU:HB2	2.46	0.45
1:C:201:SER:HA	1:C:234:PRO:HB3	1.99	0.45
4:F:156:ALA:HB3	4:F:157:PRO:HD3	1.97	0.45
7:I:201:SER:HA	7:I:234:PRO:HB3	1.99	0.45
5:G:186:VAL:O	5:G:187:ASP:C	2.60	0.45
2:L:237:ASN:O	2:L:238:VAL:HB	2.18	0.45
8:M:156:ALA:HB3	8:M:157:PRO:HD3	1.98	0.45
4:O:12:ARG:HG2	14:O:401:NAD:O2A	2.17	0.45
1:B:156:ALA:HB3	1:B:157:PRO:HD3	1.98	0.44
1:N:223:ASN:HB2	20:N:510:HOH:O	2.17	0.44
4:O:249:LYS:HE3	20:O:518:HOH:O	2.17	0.44
2:D:1:MET:HG2	2:D:2:THR:H	1.82	0.44
4:O:129:MET:HG2	4:O:146:SER:HB3	1.98	0.44
1:K:40:TYR:CD1	2:L:194:TRP:CE2	3.05	0.44
8:M:201:SER:HA	8:M:234:PRO:HB3	1.99	0.44
4:F:170:GLU:HB2	6:H:301:LEU:CD2	2.48	0.44
7:I:306:VAL:HG12	7:I:307:LYS:N	2.32	0.44
2:L:156:ALA:HB3	2:L:157:PRO:HD3	1.99	0.44
1:C:150:CYS:SG	14:C:401:NAD:H4N	2.57	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:P:201:SER:HA	9:P:234:PRO:HB3	1.98	0.44
1:A:307:LYS:HE2	1:C:172:LEU:HB3	2.00	0.43
1:B:237:ASN:O	1:B:238:VAL:HB	2.18	0.43
5:G:306:VAL:HG12	5:G:307:LYS:N	2.31	0.43
1:N:209:THR:HG23	1:N:209:THR:O	2.18	0.43
4:F:228:GLY:HA2	6:H:299:ILE:CD1	2.47	0.43
4:J:306:VAL:HG12	4:J:307:LYS:N	2.34	0.43
12:D:408:PO4:O3	20:D:537:HOH:O	2.21	0.43
4:O:237:ASN:O	4:O:238:VAL:HB	2.18	0.43
1:C:255:GLN:NE2	5:G:259:ALA:HA	2.32	0.43
8:M:209:THR:HG23	8:M:209:THR:O	2.18	0.43
6:H:237:ASN:O	6:H:238:VAL:HB	2.19	0.43
1:N:201:SER:HA	1:N:234:PRO:HB3	1.99	0.43
1:N:306:VAL:HG12	1:N:307:LYS:N	2.34	0.43
1:A:237:ASN:O	1:A:238:VAL:HB	2.17	0.43
4:F:186:VAL:O	4:F:187:ASP:C	2.62	0.43
7:I:257:ALY:HE2	7:I:295:ALA:HB1	2.01	0.43
4:J:209:THR:O	4:J:209:THR:HG23	2.19	0.43
3:E:237:ASN:O	3:E:238:VAL:HB	2.19	0.43
4:F:166:PHE:CA	4:F:249:LYS:HG2	2.42	0.43
8:M:237:ASN:O	8:M:238:VAL:HB	2.18	0.43
4:F:254:GLU:HG3	20:F:517:HOH:O	2.18	0.43
1:B:306:VAL:HG12	1:B:307:LYS:N	2.34	0.43
1:C:68:ASN:HA	20:C:547:HOH:O	2.19	0.43
3:E:186:VAL:O	3:E:187:ASP:C	2.62	0.43
9:P:156:ALA:HB3	9:P:157:PRO:HD3	2.00	0.43
2:D:306:VAL:HG12	2:D:307:LYS:N	2.34	0.43
9:P:150:SCY:CD	9:P:150:SCY:N	2.82	0.43
4:J:237:ASN:O	4:J:238:VAL:HB	2.19	0.42
4:J:307:LYS:HE2	2:L:172:LEU:HB3	2.01	0.42
2:L:209:THR:HG23	2:L:209:THR:O	2.19	0.42
4:O:257:ALY:HE2	4:O:295:ALA:HB1	2.01	0.42
9:P:186:VAL:O	9:P:187:ASP:C	2.62	0.42
9:P:237:ASN:O	9:P:238:VAL:HB	2.19	0.42
2:D:237:ASN:O	2:D:238:VAL:HB	2.19	0.42
7:I:237:ASN:O	7:I:238:VAL:HB	2.18	0.42
2:L:19:ARG:HD3	20:L:514:HOH:O	2.18	0.42
4:O:209:THR:O	4:O:209:THR:HG23	2.19	0.42
1:A:133:GLY:N	11:A:403:CL:CL	2.83	0.42
1:C:209:THR:O	1:C:209:THR:HG23	2.19	0.42
4:F:209:THR:HG23	4:F:209:THR:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:G:201:SER:HA	5:G:234:PRO:HB3	2.01	0.42
2:L:186:VAL:O	2:L:187:ASP:C	2.62	0.42
1:B:180:THR:OG1	1:B:232:ARG:NH1	2.52	0.42
1:B:201:SER:HA	1:B:234:PRO:HB3	2.02	0.42
1:K:186:VAL:O	1:K:187:ASP:C	2.62	0.42
2:L:88:VAL:HG13	2:L:90:VAL:HG23	2.02	0.42
2:D:209:THR:HG23	2:D:209:THR:O	2.19	0.42
3:E:209:THR:O	3:E:209:THR:HG23	2.20	0.42
5:G:237:ASN:O	5:G:238:VAL:HB	2.19	0.42
1:A:209:THR:HG23	1:A:209:THR:O	2.20	0.42
1:N:301:LEU:CD2	9:P:170:GLU:HB2	2.50	0.42
1:B:103:ASP:O	1:B:107:ARG:HB2	2.20	0.42
1:B:209:THR:HG23	1:B:209:THR:O	2.18	0.42
1:C:88:VAL:HG13	1:C:90:VAL:HG23	2.01	0.42
1:K:13:ILE:N	14:K:401:NAD:O2N	2.30	0.42
1:C:61:LYS:HD2	20:C:522:HOH:O	2.18	0.42
5:G:209:THR:HG23	5:G:209:THR:O	2.19	0.42
7:I:307:LYS:HE2	1:K:172:LEU:HB3	2.01	0.42
3:E:201:SER:HA	3:E:234:PRO:HB3	2.01	0.42
4:O:88:VAL:HG13	4:O:90:VAL:HG23	2.02	0.42
4:O:150:CYS:SG	14:O:401:NAD:C4N	3.08	0.42
1:C:237:ASN:O	1:C:238:VAL:HB	2.19	0.41
4:J:186:VAL:O	4:J:187:ASP:C	2.62	0.41
1:K:209:THR:HG23	1:K:209:THR:O	2.20	0.41
6:H:186:VAL:O	6:H:187:ASP:C	2.63	0.41
4:O:150:CYS:HB3	14:O:401:NAD:H4N	2.01	0.41
4:O:249:LYS:CE	20:O:518:HOH:O	2.68	0.41
1:C:37:ASP:OD1	1:C:37:ASP:N	2.47	0.41
4:F:88:VAL:HG13	4:F:90:VAL:HG23	2.02	0.41
1:K:13:ILE:HG12	14:K:401:NAD:O2N	2.19	0.41
9:P:209:THR:HG23	9:P:209:THR:O	2.20	0.41
2:D:180:THR:OG1	2:D:232:ARG:NH1	2.53	0.41
1:B:186:VAL:O	1:B:187:ASP:C	2.61	0.41
3:E:306:VAL:HG12	3:E:307:LYS:N	2.35	0.41
7:I:88:VAL:HG13	7:I:90:VAL:HG23	2.02	0.41
4:J:88:VAL:HG13	4:J:90:VAL:HG23	2.02	0.41
1:N:237:ASN:O	1:N:238:VAL:HB	2.21	0.41
6:H:209:THR:HG23	6:H:209:THR:O	2.20	0.41
7:I:209:THR:HG23	7:I:209:THR:O	2.20	0.41
1:B:246:ARG:HD3	2:D:246:ARG:CZ	2.51	0.41
3:E:88:VAL:HG13	3:E:90:VAL:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:306:VAL:HG12	1:K:307:LYS:N	2.36	0.41
1:N:88:VAL:HG13	1:N:90:VAL:HG23	2.02	0.41
8:M:88:VAL:HG13	8:M:90:VAL:HG23	2.02	0.41
4:O:186:VAL:O	4:O:187:ASP:C	2.64	0.41
9:P:88:VAL:HG13	9:P:90:VAL:HG23	2.02	0.41
4:F:39:ASP:HA	4:F:60:VAL:HG21	2.01	0.41
1:K:79:ASP:HB3	1:K:82:ASN:HD21	1.86	0.41
1:C:150:CYS:SG	14:C:401:NAD:C4N	3.09	0.40
2:D:186:VAL:O	2:D:187:ASP:C	2.63	0.40
5:G:88:VAL:HG13	5:G:90:VAL:HG23	2.03	0.40
1:C:83:LEU:O	1:C:84:LYS:C	2.65	0.40
2:D:88:VAL:HG13	2:D:90:VAL:HG23	2.03	0.40
4:O:39:ASP:HA	4:O:60:VAL:HG21	2.04	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	329/355 (93%)	314 (95%)	14 (4%)	1 (0%)	36	54
1	B	330/355 (93%)	316 (96%)	13 (4%)	1 (0%)	36	54
1	C	328/355 (92%)	315 (96%)	12 (4%)	1 (0%)	36	54
1	K	329/355 (93%)	314 (95%)	14 (4%)	1 (0%)	36	54
1	N	329/355 (93%)	315 (96%)	13 (4%)	1 (0%)	36	54
2	D	331/355 (93%)	317 (96%)	13 (4%)	1 (0%)	36	54
2	L	330/355 (93%)	315 (96%)	14 (4%)	1 (0%)	36	54
3	E	329/355 (93%)	314 (95%)	14 (4%)	1 (0%)	36	54
4	F	330/355 (93%)	316 (96%)	13 (4%)	1 (0%)	36	54
4	J	327/355 (92%)	313 (96%)	12 (4%)	2 (1%)	21	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	O	327/355 (92%)	312 (95%)	14 (4%)	1 (0%)	36	54
5	G	329/355 (93%)	315 (96%)	13 (4%)	1 (0%)	36	54
6	H	326/355 (92%)	311 (95%)	14 (4%)	1 (0%)	36	54
7	I	326/355 (92%)	311 (95%)	14 (4%)	1 (0%)	36	54
8	M	329/355 (93%)	313 (95%)	15 (5%)	1 (0%)	36	54
9	P	326/355 (92%)	312 (96%)	13 (4%)	1 (0%)	36	54
All	All	5255/5680 (92%)	5023 (96%)	215 (4%)	17 (0%)	36	54

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	238	VAL
1	A	238	VAL
1	B	238	VAL
2	D	238	VAL
3	E	238	VAL
4	F	238	VAL
5	G	238	VAL
6	H	238	VAL
7	I	238	VAL
4	J	238	VAL
1	K	238	VAL
2	L	238	VAL
8	M	238	VAL
1	N	238	VAL
4	O	238	VAL
9	P	238	VAL
4	J	167	GLY

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	266/286 (93%)	262 (98%)	4 (2%)	57	77
1	B	266/286 (93%)	263 (99%)	3 (1%)	65	81
1	C	265/286 (93%)	263 (99%)	2 (1%)	73	86
1	K	266/286 (93%)	265 (100%)	1 (0%)	84	92
1	N	266/286 (93%)	265 (100%)	1 (0%)	84	92
2	D	268/287 (93%)	265 (99%)	3 (1%)	65	81
2	L	267/287 (93%)	264 (99%)	3 (1%)	65	81
3	E	266/285 (93%)	263 (99%)	3 (1%)	65	81
4	F	267/285 (94%)	263 (98%)	4 (2%)	57	77
4	J	264/285 (93%)	263 (100%)	1 (0%)	84	92
4	O	264/285 (93%)	263 (100%)	1 (0%)	84	92
5	G	266/285 (93%)	265 (100%)	1 (0%)	84	92
6	H	263/284 (93%)	262 (100%)	1 (0%)	84	92
7	I	263/284 (93%)	262 (100%)	1 (0%)	84	92
8	M	266/286 (93%)	263 (99%)	3 (1%)	65	81
9	P	263/284 (93%)	262 (100%)	1 (0%)	84	92
All	All	4246/4567 (93%)	4213 (99%)	33 (1%)	73	86

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	THR
1	A	4	LYS
1	A	58	VAL
1	B	58	VAL
1	B	78	ARG
1	B	232	ARG
1	C	37	ASP
1	C	58	VAL
2	D	22	GLN
2	D	58	VAL
2	D	232	ARG
3	E	2	THR
3	E	58	VAL

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Mol	Chain	Res	Type
3	E	191	HIS
4	F	2	THR
4	F	58	VAL
4	F	213	LYS
4	F	249	LYS
5	G	58	VAL
6	H	58	VAL
7	I	58	VAL
4	J	58	VAL
1	K	58	VAL
2	L	58	VAL
2	L	329	ILE
2	L	331	LYS
8	M	58	VAL
8	M	107	ARG
8	M	243	LEU
1	N	58	VAL
4	O	58	VAL
9	P	58	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	A	202	GLN
1	B	126	ASN
1	B	202	GLN
1	B	328	HIS
1	C	202	GLN
1	C	328	HIS
2	D	202	GLN
3	E	202	GLN
5	G	147	ASN
5	G	202	GLN
5	G	328	HIS
6	H	202	GLN
7	I	147	ASN
7	I	202	GLN
4	J	202	GLN
4	J	328	HIS
2	L	202	GLN
8	M	191	HIS

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Mol	Chain	Res	Type
8	M	202	GLN
1	N	202	GLN
4	O	202	GLN
9	P	202	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

25 non-standard protein/DNA/RNA residues 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
1	ALY	C	46	1	10,11,12	1.38	1 (10%)	7,12,14	3.97	3 (42%)
5	SCY	G	150	5	7,8,9	0.96	0	4,9,11	1.09	0
7	ALY	I	257	7	10,11,12	1.38	2 (20%)	7,12,14	6.68	5 (71%)
4	ALY	J	46	4	10,11,12	1.26	1 (10%)	7,12,14	3.64	3 (42%)
4	ALY	J	257	4	10,11,12	1.42	2 (20%)	7,12,14	7.23	5 (71%)
5	ALY	G	46	5	10,11,12	1.06	0	7,12,14	3.56	3 (42%)
6	ALY	H	257	6	10,11,12	1.32	2 (20%)	7,12,14	6.04	3 (42%)
3	ALY	E	46	3	10,11,12	1.10	0	7,12,14	3.86	3 (42%)
9	ALY	P	249	9	10,11,12	0.43	0	7,12,14	3.24	3 (42%)
4	ALY	F	46	4	10,11,12	0.96	0	7,12,14	3.72	3 (42%)
7	ALY	I	249	7	10,11,12	0.60	0	7,12,14	2.80	4 (57%)
4	ALY	F	257	4	10,11,12	1.56	2 (20%)	7,12,14	6.09	3 (42%)
9	ALY	P	46	9	10,11,12	1.08	1 (10%)	7,12,14	3.55	3 (42%)
9	SCY	P	150	9	7,8,9	1.14	1 (14%)	4,9,11	1.18	0
3	ALY	E	249	3	10,11,12	0.66	0	7,12,14	3.23	3 (42%)
8	ALY	M	257	8	10,11,12	1.62	2 (20%)	7,12,14	6.54	5 (71%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	ALY	N	46	1	10,11,12	1.21	1 (10%)	7,12,14	3.36	3 (42%)
4	ALY	O	46	4	10,11,12	0.89	0	7,12,14	3.60	3 (42%)
6	ALY	H	46	6	10,11,12	1.53	2 (20%)	7,12,14	3.09	3 (42%)
4	ALY	O	257	4	10,11,12	1.30	2 (20%)	7,12,14	6.47	5 (71%)
1	ALY	A	46	1	10,11,12	1.20	1 (10%)	7,12,14	3.66	3 (42%)
1	ALY	B	46	1	10,11,12	1.38	2 (20%)	7,12,14	3.92	3 (42%)
6	SCY	H	150	6	7,8,9	1.01	0	4,9,11	1.36	1 (25%)
1	ALY	K	46	1	10,11,12	1.06	0	7,12,14	3.54	3 (42%)
7	ALY	I	46	7	10,11,12	1.41	2 (20%)	7,12,14	3.84	3 (42%)

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
1	ALY	C	46	1	-	4/9/10/12	-
5	SCY	G	150	5	-	2/5/7/9	-
7	ALY	I	257	7	-	2/9/10/12	-
4	ALY	J	46	4	-	4/9/10/12	-
4	ALY	J	257	4	-	2/9/10/12	-
5	ALY	G	46	5	-	4/9/10/12	-
6	ALY	H	257	6	-	5/9/10/12	-
3	ALY	E	46	3	-	4/9/10/12	-
9	ALY	P	249	9	-	4/9/10/12	-
4	ALY	F	46	4	-	4/9/10/12	-
7	ALY	I	249	7	-	3/9/10/12	-
4	ALY	F	257	4	-	5/9/10/12	-
9	ALY	P	46	9	-	4/9/10/12	-
9	SCY	P	150	9	-	3/5/7/9	-
3	ALY	E	249	3	-	4/9/10/12	-
8	ALY	M	257	8	-	2/9/10/12	-
1	ALY	N	46	1	-	4/9/10/12	-
4	ALY	O	46	4	-	4/9/10/12	-
6	ALY	H	46	6	-	4/9/10/12	-
4	ALY	O	257	4	-	2/9/10/12	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	ALY	A	46	1	-	4/9/10/12	-
1	ALY	B	46	1	-	4/9/10/12	-
6	SCY	H	150	6	-	2/5/7/9	-
1	ALY	K	46	1	-	4/9/10/12	-
7	ALY	I	46	7	-	4/9/10/12	-

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	M	257	ALY	CE-NZ	3.24	1.53	1.46
4	J	257	ALY	CE-NZ	3.21	1.53	1.46
6	H	46	ALY	CE-NZ	3.14	1.53	1.46
4	F	257	ALY	CE-NZ	3.14	1.53	1.46
4	F	257	ALY	CH-NZ	2.91	1.41	1.34
7	I	46	ALY	CE-NZ	2.88	1.52	1.46
6	H	46	ALY	OH-CH	-2.77	1.17	1.23
1	B	46	ALY	CE-NZ	2.61	1.52	1.46
8	M	257	ALY	CH-NZ	2.58	1.41	1.34
7	I	46	ALY	OH-CH	-2.54	1.17	1.23
7	I	257	ALY	CE-NZ	2.48	1.51	1.46
1	B	46	ALY	OH-CH	-2.47	1.17	1.23
9	P	46	ALY	CE-NZ	2.47	1.51	1.46
6	H	257	ALY	CE-NZ	2.39	1.51	1.46
1	N	46	ALY	CE-NZ	2.38	1.51	1.46
6	H	257	ALY	CH-NZ	2.32	1.40	1.34
4	O	257	ALY	CE-NZ	2.31	1.51	1.46
4	J	46	ALY	CE-NZ	2.30	1.51	1.46
1	C	46	ALY	CB-CA	2.25	1.57	1.53
9	P	150	SCY	CB-SG	-2.22	1.76	1.81
1	A	46	ALY	CE-NZ	2.19	1.51	1.46
7	I	257	ALY	CB-CA	-2.17	1.50	1.53
4	J	257	ALY	CH-NZ	2.10	1.39	1.34
4	O	257	ALY	CB-CA	-2.10	1.50	1.53

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	I	257	ALY	CE-NZ-CH	15.96	146.13	122.56
4	J	257	ALY	CE-NZ-CH	15.91	146.05	122.56
4	O	257	ALY	CE-NZ-CH	15.47	145.41	122.56
4	F	257	ALY	CE-NZ-CH	15.01	144.72	122.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	H	257	ALY	CE-NZ-CH	14.80	144.41	122.56
8	M	257	ALY	CE-NZ-CH	13.03	141.79	122.56
1	B	46	ALY	CE-NZ-CH	9.75	136.95	122.56
1	C	46	ALY	CE-NZ-CH	9.73	136.93	122.56
7	I	46	ALY	CE-NZ-CH	9.49	136.57	122.56
3	E	46	ALY	CE-NZ-CH	9.47	136.54	122.56
4	F	46	ALY	CE-NZ-CH	8.84	135.61	122.56
4	J	46	ALY	CE-NZ-CH	8.68	135.38	122.56
4	O	46	ALY	CE-NZ-CH	8.44	135.03	122.56
1	A	46	ALY	CE-NZ-CH	8.44	135.02	122.56
5	G	46	ALY	CE-NZ-CH	8.42	134.99	122.56
1	K	46	ALY	CE-NZ-CH	8.30	134.82	122.56
9	P	46	ALY	CE-NZ-CH	7.93	134.27	122.56
3	E	249	ALY	CE-NZ-CH	7.64	133.84	122.56
9	P	249	ALY	CE-NZ-CH	7.62	133.81	122.56
8	M	257	ALY	CH3-CH-NZ	6.71	127.57	116.12
1	N	46	ALY	CE-NZ-CH	6.34	131.92	122.56
4	J	257	ALY	CH3-CH-NZ	6.21	126.71	116.12
8	M	257	ALY	CD-CE-NZ	-6.10	95.08	112.20
6	H	46	ALY	CE-NZ-CH	5.97	131.38	122.56
4	J	257	ALY	CD-CE-NZ	-5.90	95.65	112.20
7	I	249	ALY	CE-NZ-CH	5.14	130.15	122.56
8	M	257	ALY	OH-CH-CH3	-5.07	113.03	122.05
1	N	46	ALY	CH3-CH-NZ	4.85	124.40	116.12
4	O	257	ALY	CD-CE-NZ	-4.83	98.65	112.20
7	I	257	ALY	CD-CE-NZ	-4.62	99.23	112.20
6	H	257	ALY	OH-CH-CH3	-4.53	113.99	122.05
8	M	257	ALY	CG-CD-CE	4.45	134.08	113.56
4	F	257	ALY	OH-CH-CH3	-4.40	114.22	122.05
4	J	257	ALY	OH-CH-CH3	-4.39	114.24	122.05
6	H	46	ALY	CH3-CH-NZ	4.39	123.61	116.12
7	I	249	ALY	CD-CE-NZ	-4.32	100.09	112.20
4	J	257	ALY	CG-CD-CE	4.25	133.18	113.56
9	P	46	ALY	CH3-CH-NZ	3.89	122.76	116.12
7	I	257	ALY	OH-CH-CH3	-3.84	115.21	122.05
1	N	46	ALY	OH-CH-CH3	-3.73	115.41	122.05
1	A	46	ALY	CH3-CH-NZ	3.49	122.08	116.12
4	O	46	ALY	CH3-CH-NZ	3.40	121.92	116.12
7	I	257	ALY	OH-CH-NZ	3.37	130.62	121.78
4	F	257	ALY	OH-CH-NZ	3.34	130.54	121.78
1	K	46	ALY	CH3-CH-NZ	3.31	121.77	116.12
4	J	46	ALY	CH3-CH-NZ	3.31	121.76	116.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	H	257	ALY	OH-CH-NZ	3.27	130.36	121.78
4	F	46	ALY	CH3-CH-NZ	3.25	121.67	116.12
6	H	46	ALY	OH-CH-CH3	-3.25	116.27	122.05
4	O	257	ALY	OH-CH-CH3	-3.22	116.33	122.05
1	A	46	ALY	OH-CH-CH3	-3.11	116.51	122.05
4	O	257	ALY	OH-CH-NZ	3.04	129.76	121.78
9	P	46	ALY	OH-CH-CH3	-3.04	116.64	122.05
5	G	46	ALY	CH3-CH-NZ	3.03	121.30	116.12
3	E	46	ALY	CH3-CH-NZ	3.00	121.24	116.12
1	C	46	ALY	CH3-CH-NZ	2.93	121.12	116.12
9	P	249	ALY	CH3-CH-NZ	2.91	121.10	116.12
4	O	257	ALY	CG-CD-CE	2.90	126.96	113.56
3	E	249	ALY	CH3-CH-NZ	2.84	120.96	116.12
7	I	257	ALY	CG-CD-CE	2.83	126.65	113.56
5	G	46	ALY	OH-CH-CH3	-2.72	117.20	122.05
7	I	46	ALY	CH3-CH-NZ	2.72	120.76	116.12
4	F	46	ALY	OH-CH-CH3	-2.67	117.29	122.05
4	O	46	ALY	OH-CH-CH3	-2.66	117.32	122.05
1	B	46	ALY	CH3-CH-NZ	2.63	120.60	116.12
1	K	46	ALY	OH-CH-CH3	-2.57	117.47	122.05
1	C	46	ALY	OH-CH-CH3	-2.42	117.74	122.05
4	J	46	ALY	OH-CH-CH3	-2.42	117.75	122.05
6	H	150	SCY	OCD-CD-CE	-2.30	114.24	123.12
7	I	46	ALY	OH-CH-CH3	-2.16	118.20	122.05
7	I	249	ALY	OH-CH-CH3	-2.12	118.27	122.05
3	E	46	ALY	OH-CH-CH3	-2.12	118.29	122.05
1	B	46	ALY	OH-CH-CH3	-2.10	118.31	122.05
7	I	249	ALY	CH3-CH-NZ	2.06	119.63	116.12
9	P	249	ALY	OH-CH-CH3	-2.05	118.41	122.05
3	E	249	ALY	OH-CH-CH3	-2.02	118.47	122.05

There are no chirality outliers.

All (88) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	257	ALY	C-CA-CB-CG
5	G	150	SCY	OCD-CD-SG-CB
5	G	150	SCY	CE-CD-SG-CB
6	H	150	SCY	OCD-CD-SG-CB
6	H	150	SCY	CE-CD-SG-CB
6	H	257	ALY	C-CA-CB-CG
9	P	150	SCY	CA-CB-SG-CD

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Mol	Chain	Res	Type	Atoms
9	P	150	SCY	OCD-CD-SG-CB
9	P	150	SCY	CE-CD-SG-CB
4	O	257	ALY	CG-CD-CE-NZ
3	E	249	ALY	CG-CD-CE-NZ
4	F	46	ALY	CG-CD-CE-NZ
5	G	46	ALY	CG-CD-CE-NZ
1	K	46	ALY	CG-CD-CE-NZ
4	O	46	ALY	CG-CD-CE-NZ
1	A	46	ALY	OH-CH-NZ-CE
1	A	46	ALY	CH3-CH-NZ-CE
1	B	46	ALY	OH-CH-NZ-CE
1	B	46	ALY	CH3-CH-NZ-CE
1	C	46	ALY	OH-CH-NZ-CE
1	C	46	ALY	CH3-CH-NZ-CE
3	E	46	ALY	OH-CH-NZ-CE
3	E	46	ALY	CH3-CH-NZ-CE
3	E	249	ALY	OH-CH-NZ-CE
3	E	249	ALY	CH3-CH-NZ-CE
4	F	46	ALY	OH-CH-NZ-CE
4	F	46	ALY	CH3-CH-NZ-CE
5	G	46	ALY	OH-CH-NZ-CE
5	G	46	ALY	CH3-CH-NZ-CE
6	H	46	ALY	OH-CH-NZ-CE
6	H	46	ALY	CH3-CH-NZ-CE
7	I	46	ALY	OH-CH-NZ-CE
7	I	46	ALY	CH3-CH-NZ-CE
7	I	249	ALY	OH-CH-NZ-CE
7	I	249	ALY	CH3-CH-NZ-CE
4	J	46	ALY	OH-CH-NZ-CE
4	J	46	ALY	CH3-CH-NZ-CE
4	J	257	ALY	OH-CH-NZ-CE
4	J	257	ALY	CH3-CH-NZ-CE
1	K	46	ALY	OH-CH-NZ-CE
1	K	46	ALY	CH3-CH-NZ-CE
8	M	257	ALY	OH-CH-NZ-CE
8	M	257	ALY	CH3-CH-NZ-CE
1	N	46	ALY	OH-CH-NZ-CE
1	N	46	ALY	CH3-CH-NZ-CE
4	O	46	ALY	OH-CH-NZ-CE
4	O	46	ALY	CH3-CH-NZ-CE
9	P	46	ALY	OH-CH-NZ-CE
9	P	46	ALY	CH3-CH-NZ-CE

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Mol	Chain	Res	Type	Atoms
9	P	249	ALY	OH-CH-NZ-CE
9	P	249	ALY	CH3-CH-NZ-CE
1	C	46	ALY	CG-CD-CE-NZ
3	E	46	ALY	CG-CD-CE-NZ
6	H	46	ALY	CG-CD-CE-NZ
7	I	257	ALY	CG-CD-CE-NZ
1	N	46	ALY	CG-CD-CE-NZ
9	P	46	ALY	CG-CD-CE-NZ
9	P	249	ALY	CG-CD-CE-NZ
1	B	46	ALY	CG-CD-CE-NZ
4	J	46	ALY	CG-CD-CE-NZ
1	A	46	ALY	CG-CD-CE-NZ
7	I	46	ALY	CG-CD-CE-NZ
4	F	257	ALY	CG-CD-CE-NZ
6	H	257	ALY	CG-CD-CE-NZ
4	F	46	ALY	CE-CD-CG-CB
5	G	46	ALY	CE-CD-CG-CB
4	F	257	ALY	CA-CB-CG-CD
6	H	257	ALY	CA-CB-CG-CD
1	C	46	ALY	CE-CD-CG-CB
4	O	46	ALY	CE-CD-CG-CB
1	B	46	ALY	CE-CD-CG-CB
3	E	46	ALY	CE-CD-CG-CB
1	K	46	ALY	CE-CD-CG-CB
4	J	46	ALY	CE-CD-CG-CB
1	N	46	ALY	CE-CD-CG-CB
6	H	46	ALY	CE-CD-CG-CB
7	I	46	ALY	CE-CD-CG-CB
9	P	46	ALY	CE-CD-CG-CB
1	A	46	ALY	CE-CD-CG-CB
7	I	249	ALY	CA-CB-CG-CD
4	F	257	ALY	CE-CD-CG-CB
6	H	257	ALY	CE-CD-CG-CB
4	F	257	ALY	N-CA-CB-CG
6	H	257	ALY	N-CA-CB-CG
7	I	257	ALY	CE-CD-CG-CB
9	P	249	ALY	CE-CD-CG-CB
3	E	249	ALY	CE-CD-CG-CB
4	O	257	ALY	CE-CD-CG-CB

There are no ring outliers.

18 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	46	ALY	1	0
7	I	257	ALY	1	0
4	J	46	ALY	1	0
5	G	46	ALY	1	0
6	H	257	ALY	1	0
3	E	46	ALY	1	0
4	F	46	ALY	1	0
4	F	257	ALY	1	0
9	P	46	ALY	1	0
9	P	150	SCY	5	0
1	N	46	ALY	1	0
4	O	46	ALY	1	0
6	H	46	ALY	1	0
4	O	257	ALY	1	0
1	A	46	ALY	1	0
1	B	46	ALY	1	0
1	K	46	ALY	1	0
7	I	46	ALY	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 94 ligands modelled in this entry, 61 are monoatomic - leaving 33 for Mogul analysis.

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	$\# Z > 2$
12	PO4	D	408	-	4,4,4	1.56	1 (25%)	6,6,6	1.26	1 (16%)
14	NAD	O	401	-	46,48,48	1.31	6 (13%)	64,73,73	1.91	11 (17%)
13	POP	B	410	-	6,8,8	0.80	0	12,13,13	1.44	2 (16%)
12	PO4	E	408	-	4,4,4	0.44	0	6,6,6	1.12	1 (16%)
12	PO4	L	402	-	4,4,4	0.82	0	6,6,6	0.69	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
12	PO4	H	403	-	4,4,4	0.97	0	6,6,6	0.84	0
19	UVW	M	404	-	6,7,7	1.22	1 (16%)	7,10,10	1.33	1 (14%)
12	PO4	D	409	-	4,4,4	1.25	0	6,6,6	0.61	0
12	PO4	J	404	-	4,4,4	1.09	0	6,6,6	0.58	0
16	PGE	H	405	-	9,9,9	0.72	0	8,8,8	0.36	0
13	POP	G	406	-	6,8,8	1.13	0	12,13,13	0.92	0
12	PO4	F	402	-	4,4,4	0.92	0	6,6,6	0.56	0
13	POP	F	403	-	6,8,8	0.74	0	12,13,13	1.35	0
12	PO4	P	405	-	4,4,4	1.15	0	6,6,6	1.01	0
14	NAD	K	401	-	46,48,48	1.39	7 (15%)	64,73,73	1.85	15 (23%)
12	PO4	A	406	-	4,4,4	0.52	0	6,6,6	1.95	1 (16%)
13	POP	A	409	-	6,8,8	0.90	0	12,13,13	1.63	3 (25%)
17	PEG	H	406	-	6,6,6	0.52	0	5,5,5	0.65	0
12	PO4	J	403	-	4,4,4	0.89	0	6,6,6	0.72	0
12	PO4	P	404	-	4,4,4	0.72	0	6,6,6	0.91	0
13	POP	J	406	-	6,8,8	0.98	0	12,13,13	1.31	1 (8%)
12	PO4	A	408	-	4,4,4	0.64	0	6,6,6	0.89	0
14	NAD	C	401	-	46,48,48	1.77	9 (19%)	64,73,73	1.90	14 (21%)
18	PG4	J	405	-	12,12,12	0.62	0	11,11,11	0.24	0
14	NAD	I	401	-	46,48,48	1.39	7 (15%)	64,73,73	1.60	10 (15%)
13	POP	M	405	-	6,8,8	0.48	0	12,13,13	1.68	4 (33%)
15	ACT	H	404	-	3,3,3	0.93	0	3,3,3	0.71	0
12	PO4	A	407	-	4,4,4	1.08	0	6,6,6	1.02	0
14	NAD	P	401	-	46,48,48	1.44	8 (17%)	64,73,73	1.69	16 (25%)
14	NAD	D	401	-	46,48,48	1.30	6 (13%)	64,73,73	1.78	16 (25%)
12	PO4	B	409	-	4,4,4	0.99	0	6,6,6	2.35	4 (66%)
12	PO4	I	405	-	4,4,4	0.99	0	6,6,6	0.84	0
14	NAD	E	401	-	46,48,48	1.25	4 (8%)	64,73,73	1.75	13 (20%)

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
19	UVW	M	404	-	-	0/3/5/5	-
14	NAD	I	401	-	-	19/30/62/62	0/5/5/5
14	NAD	K	401	-	-	8/30/62/62	0/5/5/5
14	NAD	O	401	-	-	18/30/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
13	POP	M	405	-	-	3/6/6/6	-
13	POP	A	409	-	-	3/6/6/6	-
13	POP	B	410	-	-	3/6/6/6	-
17	PEG	H	406	-	-	2/4/4/4	-
14	NAD	P	401	-	-	19/30/62/62	0/5/5/5
13	POP	J	406	-	-	1/6/6/6	-
14	NAD	D	401	-	-	18/30/62/62	0/5/5/5
14	NAD	E	401	-	-	12/30/62/62	0/5/5/5
13	POP	G	406	-	-	0/6/6/6	-
16	PGE	H	405	-	-	0/7/7/7	-
13	POP	F	403	-	-	2/6/6/6	-
14	NAD	C	401	-	-	14/30/62/62	0/5/5/5
18	PG4	J	405	-	-	4/10/10/10	-

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	C	401	NAD	C5A-C4A	5.67	1.49	1.39
14	P	401	NAD	C5A-C4A	4.87	1.47	1.39
14	D	401	NAD	C5A-C4A	4.80	1.47	1.39
14	I	401	NAD	C5A-C4A	4.73	1.47	1.39
14	K	401	NAD	C5A-C4A	4.67	1.47	1.39
14	E	401	NAD	C5A-C4A	4.42	1.47	1.39
14	C	401	NAD	O4D-C1D	4.42	1.46	1.40
14	C	401	NAD	C5A-C6A	4.19	1.52	1.41
14	O	401	NAD	PA-O3	4.11	1.63	1.59
14	P	401	NAD	O4D-C1D	4.09	1.46	1.40
14	K	401	NAD	O4D-C1D	4.04	1.46	1.40
14	I	401	NAD	O4D-C1D	3.76	1.45	1.40
14	C	401	NAD	PA-O3	3.60	1.63	1.59
14	O	401	NAD	C5A-C4A	3.47	1.45	1.39
14	K	401	NAD	C5A-C6A	3.24	1.50	1.41
14	C	401	NAD	PN-O3	3.23	1.63	1.59
14	P	401	NAD	C8A-N7A	3.21	1.37	1.31
14	I	401	NAD	C8A-N7A	3.16	1.37	1.31
14	C	401	NAD	C8A-N7A	3.06	1.37	1.31
14	O	401	NAD	C5A-C6A	3.01	1.49	1.41
14	O	401	NAD	O4D-C1D	2.93	1.44	1.40
12	D	408	PO4	P-O1	2.91	1.57	1.50
14	I	401	NAD	PA-O3	2.90	1.62	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	E	401	NAD	C5A-C6A	2.90	1.49	1.41
14	D	401	NAD	PN-O3	2.90	1.62	1.59
14	C	401	NAD	C1B-N9A	2.75	1.53	1.46
14	D	401	NAD	C8A-N7A	2.71	1.36	1.31
14	E	401	NAD	C8A-N7A	2.66	1.36	1.31
14	D	401	NAD	O4D-C1D	2.63	1.44	1.40
19	M	404	UVW	P-O2	2.61	1.64	1.59
14	P	401	NAD	PN-O3	2.56	1.62	1.59
14	O	401	NAD	PN-O3	2.56	1.62	1.59
14	I	401	NAD	C5A-C6A	2.52	1.48	1.41
14	O	401	NAD	C8A-N7A	2.50	1.36	1.31
14	K	401	NAD	C8A-N7A	2.48	1.36	1.31
14	I	401	NAD	PN-O3	2.47	1.62	1.59
14	D	401	NAD	C5A-C6A	2.46	1.47	1.41
14	P	401	NAD	C4A-N9A	-2.43	1.32	1.37
14	K	401	NAD	C4A-N9A	-2.33	1.32	1.37
14	E	401	NAD	PN-O3	2.26	1.61	1.59
14	P	401	NAD	C5A-C6A	2.26	1.47	1.41
14	P	401	NAD	PA-O3	2.21	1.61	1.59
14	D	401	NAD	C5A-N7A	-2.21	1.35	1.39
14	P	401	NAD	C5A-N7A	-2.19	1.35	1.39
14	I	401	NAD	C5A-N7A	-2.16	1.35	1.39
14	C	401	NAD	C3N-C7N	2.15	1.53	1.50
14	K	401	NAD	C5A-N7A	-2.11	1.35	1.39
14	K	401	NAD	PN-O3	2.07	1.61	1.59
14	C	401	NAD	C2A-N3A	2.07	1.37	1.33

All (113) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	C	401	NAD	C5A-C4A-N3A	-6.65	117.56	126.72
14	O	401	NAD	C5A-C4A-N3A	-6.30	118.04	126.72
14	K	401	NAD	C5A-C4A-N3A	-6.17	118.22	126.72
14	I	401	NAD	C5A-C4A-N3A	-6.08	118.35	126.72
14	D	401	NAD	C5A-C4A-N3A	-5.96	118.51	126.72
14	E	401	NAD	C5A-C4A-N3A	-5.48	119.17	126.72
14	C	401	NAD	C4A-C5A-N7A	-4.83	105.06	110.58
14	D	401	NAD	N3A-C4A-N9A	4.71	135.18	127.17
14	I	401	NAD	N3A-C4A-N9A	4.65	135.07	127.17
14	O	401	NAD	C4A-C5A-N7A	-4.61	105.31	110.58
14	O	401	NAD	C2A-N3A-C4A	4.55	122.93	111.83
14	P	401	NAD	C5A-C4A-N3A	-4.54	120.47	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	O	401	NAD	N3A-C4A-N9A	4.48	134.78	127.17
14	K	401	NAD	N3A-C4A-N9A	4.35	134.57	127.17
14	O	401	NAD	N3A-C2A-N1A	-4.34	122.01	128.58
14	C	401	NAD	C2A-N3A-C4A	4.34	122.43	111.83
14	K	401	NAD	C4A-C5A-N7A	-4.25	105.72	110.58
14	E	401	NAD	N3A-C2A-N1A	-4.21	122.21	128.58
14	C	401	NAD	N3A-C4A-N9A	4.11	134.16	127.17
14	D	401	NAD	N3A-C2A-N1A	-4.11	122.36	128.58
14	E	401	NAD	C2A-N3A-C4A	4.09	121.81	111.83
14	D	401	NAD	C2A-N3A-C4A	4.08	121.80	111.83
14	E	401	NAD	N3A-C4A-N9A	4.05	134.06	127.17
14	K	401	NAD	C2A-N3A-C4A	3.98	121.54	111.83
14	P	401	NAD	N3A-C2A-N1A	-3.96	122.58	128.58
14	O	401	NAD	C5A-N7A-C8A	3.82	109.45	103.45
14	I	401	NAD	C2A-N3A-C4A	3.79	121.08	111.83
12	A	406	PO4	O4-P-O2	3.71	119.45	107.91
14	E	401	NAD	C4A-C5A-N7A	-3.70	106.35	110.58
14	K	401	NAD	O2N-PN-O3	-3.69	97.30	107.27
14	C	401	NAD	O3-PN-O1N	-3.66	99.70	110.70
14	E	401	NAD	C3N-C7N-N7N	-3.60	113.30	117.74
14	I	401	NAD	C4A-C5A-N7A	-3.60	106.46	110.58
13	M	405	POP	O6-P2-O5	3.55	121.11	107.80
14	P	401	NAD	N3A-C4A-N9A	3.50	133.12	127.17
14	O	401	NAD	N9A-C8A-N7A	-3.46	109.03	113.94
14	P	401	NAD	C2A-N3A-C4A	3.46	120.27	111.83
12	B	409	PO4	O4-P-O1	3.45	123.15	110.95
14	I	401	NAD	N3A-C2A-N1A	-3.35	123.52	128.58
14	O	401	NAD	C4A-N9A-C8A	3.14	109.04	105.74
14	E	401	NAD	C5A-N7A-C8A	3.09	108.31	103.45
14	O	401	NAD	C6A-C5A-N7A	3.09	138.05	132.09
14	C	401	NAD	O3-PA-O1A	-3.09	101.41	110.70
14	C	401	NAD	N3A-C2A-N1A	-3.08	123.92	128.58
14	K	401	NAD	C6N-N1N-C2N	-3.07	119.27	121.88
14	C	401	NAD	C5A-N7A-C8A	3.07	108.27	103.45
14	K	401	NAD	N3A-C2A-N1A	-3.03	123.99	128.58
14	P	401	NAD	O3-PN-O1N	-3.01	101.66	110.70
14	P	401	NAD	C4A-N9A-C8A	2.99	108.88	105.74
14	C	401	NAD	C6A-C5A-N7A	2.98	137.84	132.09
14	D	401	NAD	C4A-C5A-N7A	-2.96	107.20	110.58
12	B	409	PO4	O3-P-O2	2.92	116.99	107.91
14	C	401	NAD	O2N-PN-O1N	2.90	125.93	112.44
14	I	401	NAD	C5A-N7A-C8A	2.87	107.96	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	A	409	POP	O-P2-O4	-2.87	95.95	111.04
14	K	401	NAD	C5A-N7A-C8A	2.84	107.91	103.45
14	P	401	NAD	C4A-C5A-N7A	-2.81	107.37	110.58
14	I	401	NAD	C4A-N9A-C8A	2.79	108.67	105.74
14	E	401	NAD	C6A-C5A-N7A	2.77	137.43	132.09
14	E	401	NAD	N9A-C8A-N7A	-2.75	110.04	113.94
14	D	401	NAD	O2N-PN-O1N	2.72	125.12	112.44
13	B	410	POP	O6-P2-O5	2.66	117.77	107.80
14	E	401	NAD	C4A-N9A-C8A	2.65	108.52	105.74
14	K	401	NAD	C4A-N9A-C8A	2.61	108.48	105.74
19	M	404	UVW	O2-P-O1P	-2.59	101.17	109.47
14	E	401	NAD	O2N-PN-O1N	2.58	124.45	112.44
14	K	401	NAD	C5N-C4N-C3N	-2.56	117.85	120.36
14	P	401	NAD	C2A-N1A-C6A	2.55	122.92	118.73
14	I	401	NAD	N9A-C8A-N7A	-2.55	110.32	113.94
12	B	409	PO4	O4-P-O2	-2.54	100.01	107.91
14	D	401	NAD	C2B-C3B-C4B	2.50	107.44	102.61
14	C	401	NAD	C6N-N1N-C2N	-2.50	119.75	121.88
14	P	401	NAD	C2N-C3N-C4N	2.50	121.16	118.26
13	A	409	POP	O6-P2-O4	2.46	120.43	110.83
14	D	401	NAD	O2A-PA-O3	-2.45	100.64	107.27
13	A	409	POP	O3-P1-O2	2.44	116.96	107.80
14	D	401	NAD	C4A-N9A-C8A	2.40	108.26	105.74
14	D	401	NAD	C5A-N7A-C8A	2.40	107.23	103.45
12	B	409	PO4	O3-P-O1	-2.34	102.69	110.95
14	K	401	NAD	O3-PA-O1A	-2.33	103.70	110.70
14	P	401	NAD	C6N-N1N-C2N	-2.33	119.90	121.88
14	K	401	NAD	O2N-PN-O1N	2.32	123.24	112.44
13	M	405	POP	O-P1-O1	-2.31	98.91	111.04
13	J	406	POP	O6-P2-O5	2.30	116.41	107.80
14	K	401	NAD	N9A-C8A-N7A	-2.29	110.69	113.94
14	P	401	NAD	N9A-C8A-N7A	-2.27	110.72	113.94
14	C	401	NAD	C5A-C4A-N9A	2.26	108.28	105.81
14	K	401	NAD	C2B-C1B-N9A	-2.25	107.71	113.30
12	E	408	PO4	O4-P-O1	-2.24	103.04	110.95
14	O	401	NAD	C2A-N1A-C6A	2.22	122.38	118.73
13	M	405	POP	O2-P1-O1	2.22	119.47	110.83
14	P	401	NAD	C6A-C5A-N7A	2.21	136.35	132.09
14	D	401	NAD	C3N-C7N-N7N	2.20	120.45	117.74
14	C	401	NAD	C2N-N1N-C1D	2.20	123.98	119.13
14	E	401	NAD	C2A-N1A-C6A	2.19	122.33	118.73
14	O	401	NAD	C6N-N1N-C2N	-2.19	120.02	121.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	B	410	POP	O3-P1-O1	2.18	119.33	110.83
14	E	401	NAD	C2N-C3N-C4N	2.18	120.79	118.26
14	D	401	NAD	C2A-N1A-C6A	2.17	122.29	118.73
12	D	408	PO4	O4-P-O1	-2.16	103.32	110.95
14	P	401	NAD	O2N-PN-O1N	2.15	122.47	112.44
13	M	405	POP	O-P2-O4	-2.15	99.73	111.04
14	P	401	NAD	C5N-C4N-C3N	-2.13	118.27	120.36
14	I	401	NAD	C6A-C5A-N7A	2.09	136.13	132.09
14	D	401	NAD	O3D-C3D-C2D	-2.09	105.12	111.82
14	D	401	NAD	O7N-C7N-N7N	-2.09	119.60	122.62
14	I	401	NAD	O3D-C3D-C2D	-2.08	105.14	111.82
14	K	401	NAD	C6A-C5A-N7A	2.08	136.10	132.09
14	C	401	NAD	C2B-C1B-N9A	2.08	118.46	113.30
14	D	401	NAD	N9A-C8A-N7A	-2.06	111.01	113.94
14	D	401	NAD	O2B-C2B-C1B	2.04	117.13	110.10
14	P	401	NAD	C5A-N7A-C8A	2.02	106.63	103.45
14	P	401	NAD	O4B-C1B-N9A	2.01	111.96	108.09

There are no chirality outliers.

All (126) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
13	A	409	POP	P2-O-P1-O2
13	B	410	POP	P1-O-P2-O5
13	F	403	POP	P2-O-P1-O3
13	M	405	POP	P2-O-P1-O2
13	M	405	POP	P1-O-P2-O6
14	C	401	NAD	C5B-O5B-PA-O1A
14	C	401	NAD	C5B-O5B-PA-O3
14	C	401	NAD	O4D-C1D-N1N-C2N
14	C	401	NAD	O4D-C1D-N1N-C6N
14	D	401	NAD	C5B-O5B-PA-O1A
14	D	401	NAD	C5B-O5B-PA-O3
14	D	401	NAD	O4B-C4B-C5B-O5B
14	D	401	NAD	C5D-O5D-PN-O3
14	D	401	NAD	C5D-O5D-PN-O1N
14	D	401	NAD	C5D-O5D-PN-O2N
14	D	401	NAD	O4D-C1D-N1N-C2N
14	E	401	NAD	C5B-O5B-PA-O1A
14	E	401	NAD	C5B-O5B-PA-O2A
14	E	401	NAD	C5B-O5B-PA-O3
14	E	401	NAD	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
14	E	401	NAD	C5D-O5D-PN-O1N
14	E	401	NAD	O4D-C4D-C5D-O5D
14	E	401	NAD	O4D-C1D-N1N-C2N
14	I	401	NAD	C5B-O5B-PA-O1A
14	I	401	NAD	C5B-O5B-PA-O3
14	I	401	NAD	O4D-C4D-C5D-O5D
14	I	401	NAD	C3D-C4D-C5D-O5D
14	I	401	NAD	O4D-C1D-N1N-C2N
14	K	401	NAD	C5B-O5B-PA-O2A
14	K	401	NAD	O4D-C1D-N1N-C2N
14	O	401	NAD	C5B-O5B-PA-O1A
14	O	401	NAD	C5B-O5B-PA-O3
14	O	401	NAD	O4D-C1D-N1N-C2N
14	O	401	NAD	O4D-C1D-N1N-C6N
14	O	401	NAD	C2D-C1D-N1N-C2N
14	O	401	NAD	C2D-C1D-N1N-C6N
14	O	401	NAD	C2N-C3N-C7N-O7N
14	O	401	NAD	C2N-C3N-C7N-N7N
14	O	401	NAD	C4N-C3N-C7N-N7N
14	P	401	NAD	C5B-O5B-PA-O2A
14	P	401	NAD	C5B-O5B-PA-O3
14	P	401	NAD	C5D-O5D-PN-O3
14	P	401	NAD	C5D-O5D-PN-O1N
14	P	401	NAD	C5D-O5D-PN-O2N
14	P	401	NAD	O4D-C1D-N1N-C2N
14	P	401	NAD	O4D-C1D-N1N-C6N
14	P	401	NAD	C2D-C1D-N1N-C2N
14	P	401	NAD	C2D-C1D-N1N-C6N
14	P	401	NAD	C2N-C3N-C7N-O7N
14	P	401	NAD	C2N-C3N-C7N-N7N
14	P	401	NAD	C4N-C3N-C7N-O7N
14	P	401	NAD	C4N-C3N-C7N-N7N
14	O	401	NAD	C4N-C3N-C7N-O7N
17	H	406	PEG	C4-C3-O2-C2
14	D	401	NAD	C3B-C4B-C5B-O5B
14	E	401	NAD	C3B-C4B-C5B-O5B
14	P	401	NAD	O4B-C4B-C5B-O5B
14	C	401	NAD	C2N-C3N-C7N-O7N
14	C	401	NAD	C2N-C3N-C7N-N7N
14	D	401	NAD	C2N-C3N-C7N-O7N
18	J	405	PG4	O3-C5-C6-O4
14	C	401	NAD	C4N-C3N-C7N-N7N

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Mol	Chain	Res	Type	Atoms
14	E	401	NAD	C3D-C4D-C5D-O5D
14	I	401	NAD	O4B-C4B-C5B-O5B
14	O	401	NAD	O4B-C4B-C5B-O5B
14	O	401	NAD	O4D-C4D-C5D-O5D
14	O	401	NAD	C3D-C4D-C5D-O5D
14	C	401	NAD	C4N-C3N-C7N-O7N
18	J	405	PG4	O1-C1-C2-O2
14	I	401	NAD	C3B-C4B-C5B-O5B
14	O	401	NAD	C3B-C4B-C5B-O5B
14	P	401	NAD	C3B-C4B-C5B-O5B
14	D	401	NAD	C4N-C3N-C7N-O7N
14	C	401	NAD	O4B-C4B-C5B-O5B
14	D	401	NAD	C4N-C3N-C7N-N7N
14	D	401	NAD	C2N-C3N-C7N-N7N
14	C	401	NAD	C3B-C4B-C5B-O5B
13	J	406	POP	P2-O-P1-O1
14	P	401	NAD	PN-O3-PA-O1A
14	D	401	NAD	PN-O3-PA-O5B
14	O	401	NAD	PN-O3-PA-O5B
18	J	405	PG4	O4-C7-C8-O5
14	P	401	NAD	C4B-C5B-O5B-PA
14	D	401	NAD	PA-O3-PN-O1N
14	D	401	NAD	PA-O3-PN-O2N
14	I	401	NAD	PA-O3-PN-O2N
14	K	401	NAD	PA-O3-PN-O1N
14	C	401	NAD	C5B-O5B-PA-O2A
14	D	401	NAD	C5B-O5B-PA-O2A
14	I	401	NAD	C5B-O5B-PA-O2A
14	I	401	NAD	C5D-O5D-PN-O3
14	I	401	NAD	C5D-O5D-PN-O1N
14	I	401	NAD	C5D-O5D-PN-O2N
14	K	401	NAD	C5B-O5B-PA-O1A
14	K	401	NAD	C5B-O5B-PA-O3
14	O	401	NAD	C5B-O5B-PA-O2A
14	I	401	NAD	C4N-C3N-C7N-O7N
17	H	406	PEG	O2-C3-C4-O4
14	E	401	NAD	PA-O3-PN-O2N
14	I	401	NAD	C2N-C3N-C7N-O7N
14	K	401	NAD	O4D-C4D-C5D-O5D
13	M	405	POP	P2-O-P1-O1
14	D	401	NAD	C4B-C5B-O5B-PA
14	D	401	NAD	O4D-C1D-N1N-C6N

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Mol	Chain	Res	Type	Atoms
14	E	401	NAD	O4D-C1D-N1N-C6N
14	I	401	NAD	O4D-C1D-N1N-C6N
14	K	401	NAD	O4D-C1D-N1N-C6N
14	E	401	NAD	PA-O3-PN-O1N
14	O	401	NAD	PA-O3-PN-O2N
14	P	401	NAD	PN-O3-PA-O5B
13	A	409	POP	P2-O-P1-O3
13	A	409	POP	P1-O-P2-O6
13	B	410	POP	P1-O-P2-O6
13	F	403	POP	P1-O-P2-O5
14	I	401	NAD	C4N-C3N-C7N-N7N
18	J	405	PG4	C6-C5-O3-C4
14	I	401	NAD	C2N-C3N-C7N-N7N
14	O	401	NAD	C4D-C5D-O5D-PN
14	P	401	NAD	C2B-C1B-N9A-C8A
14	I	401	NAD	PA-O3-PN-O1N
14	K	401	NAD	PA-O3-PN-O2N
13	B	410	POP	P1-O-P2-O4
14	C	401	NAD	O4B-C1B-N9A-C8A
14	C	401	NAD	C2B-C1B-N9A-C8A
14	I	401	NAD	C2B-C1B-N9A-C8A
14	C	401	NAD	PA-O3-PN-O1N

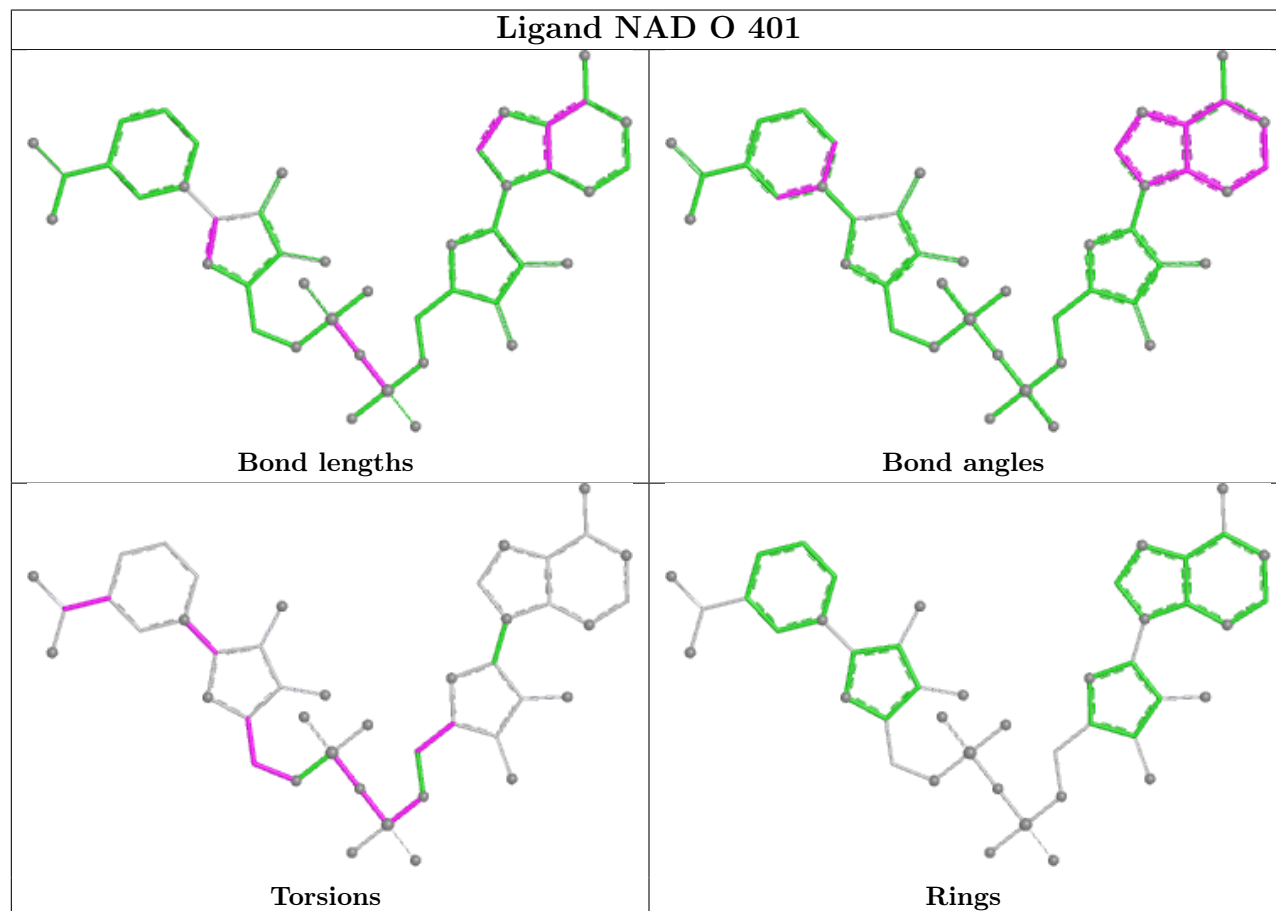
There are no ring outliers.

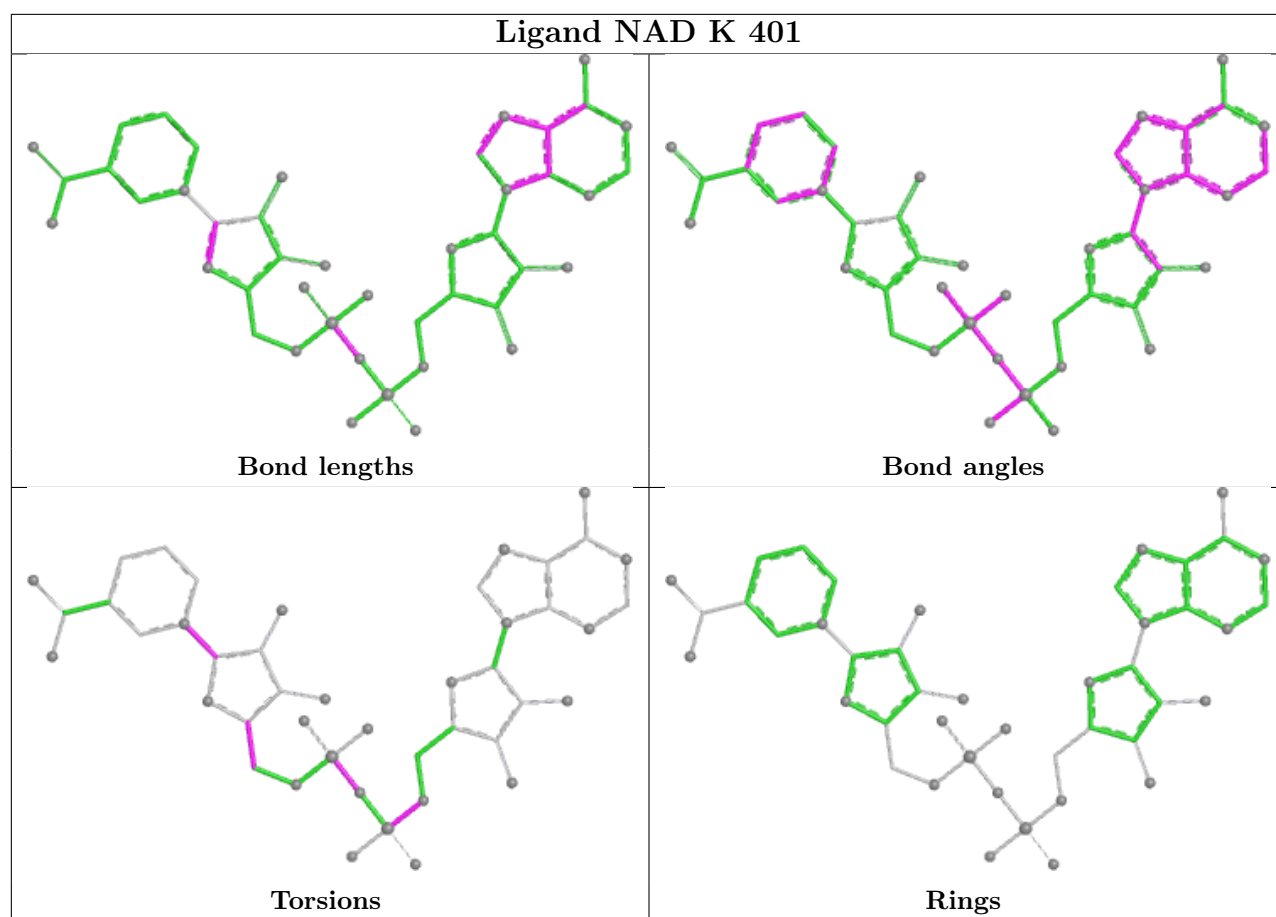
12 monomers are involved in 26 short contacts:

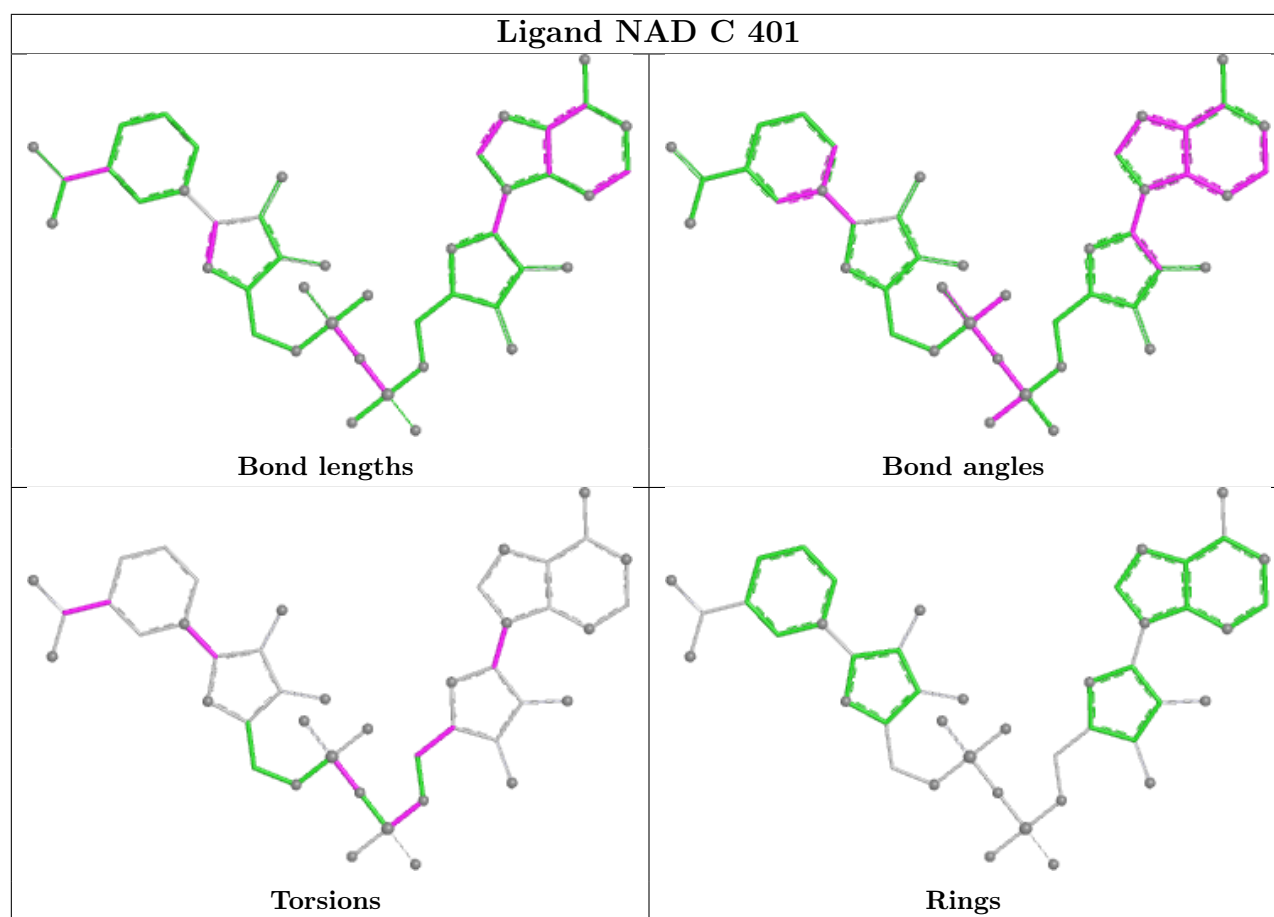
Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	D	408	PO4	1	0
14	O	401	NAD	4	0
13	B	410	POP	2	0
12	D	409	PO4	1	0
13	G	406	POP	1	0
14	K	401	NAD	4	0
13	A	409	POP	1	0
13	J	406	POP	1	0
14	C	401	NAD	4	0
12	A	407	PO4	2	0
14	P	401	NAD	2	0
14	D	401	NAD	3	0

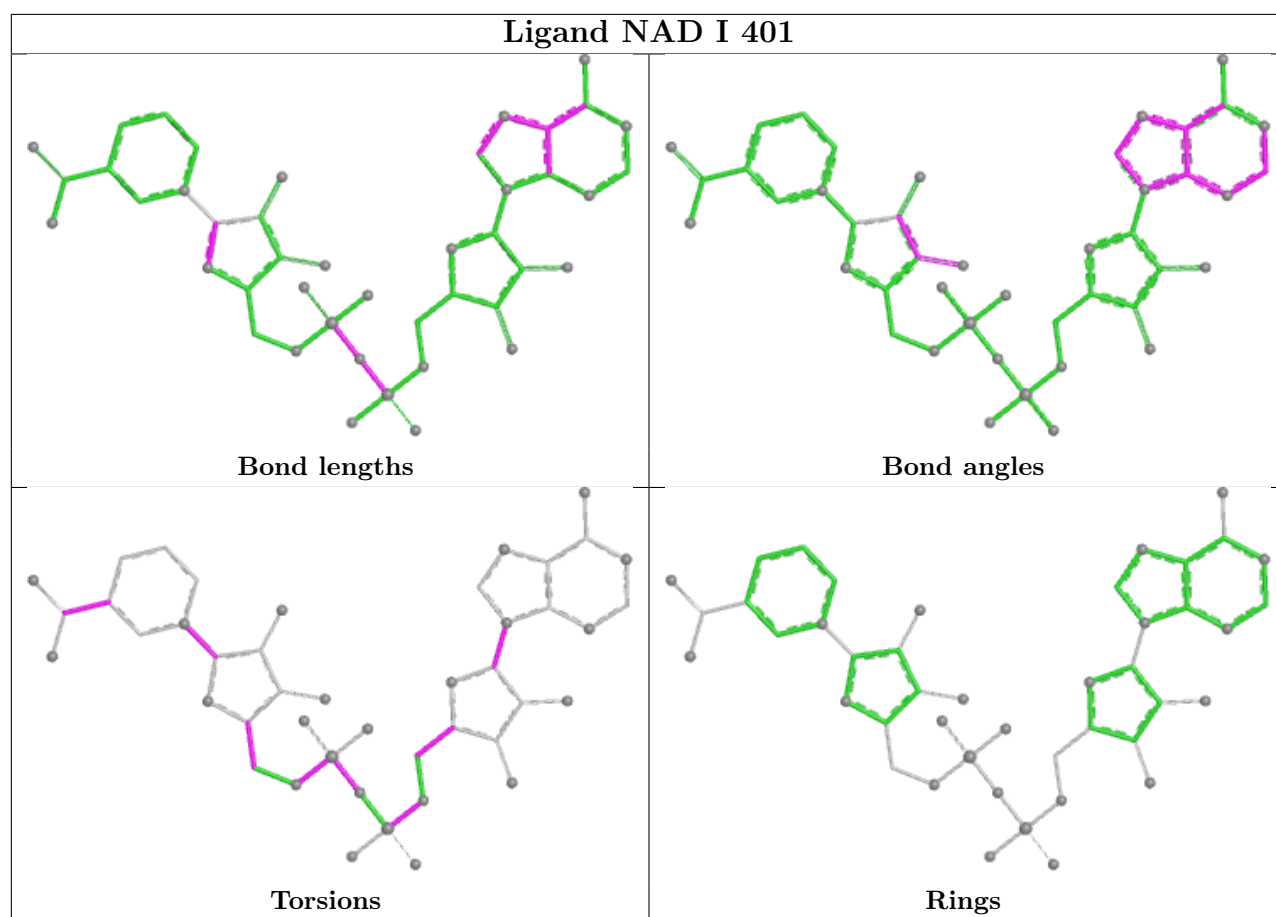
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

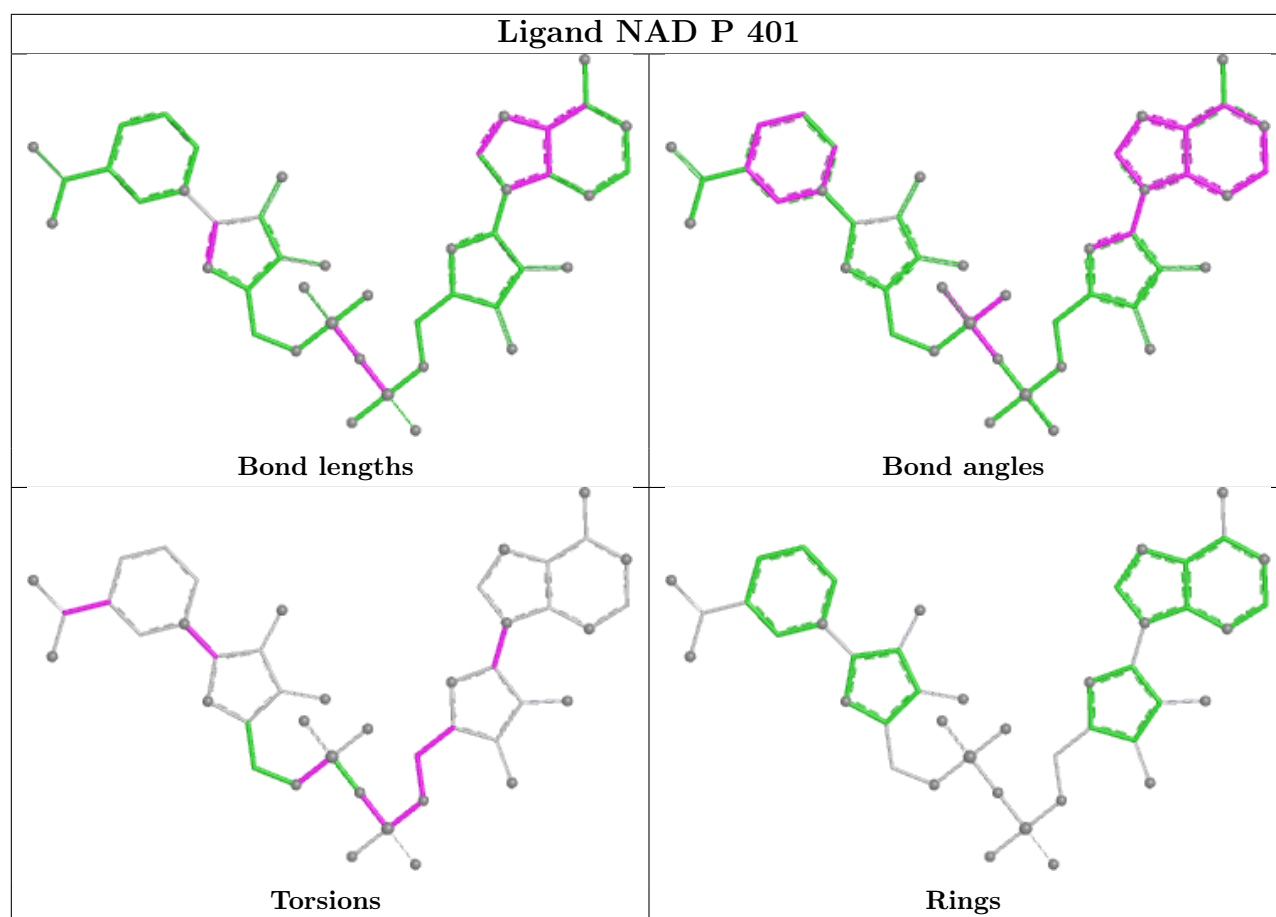
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. 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.

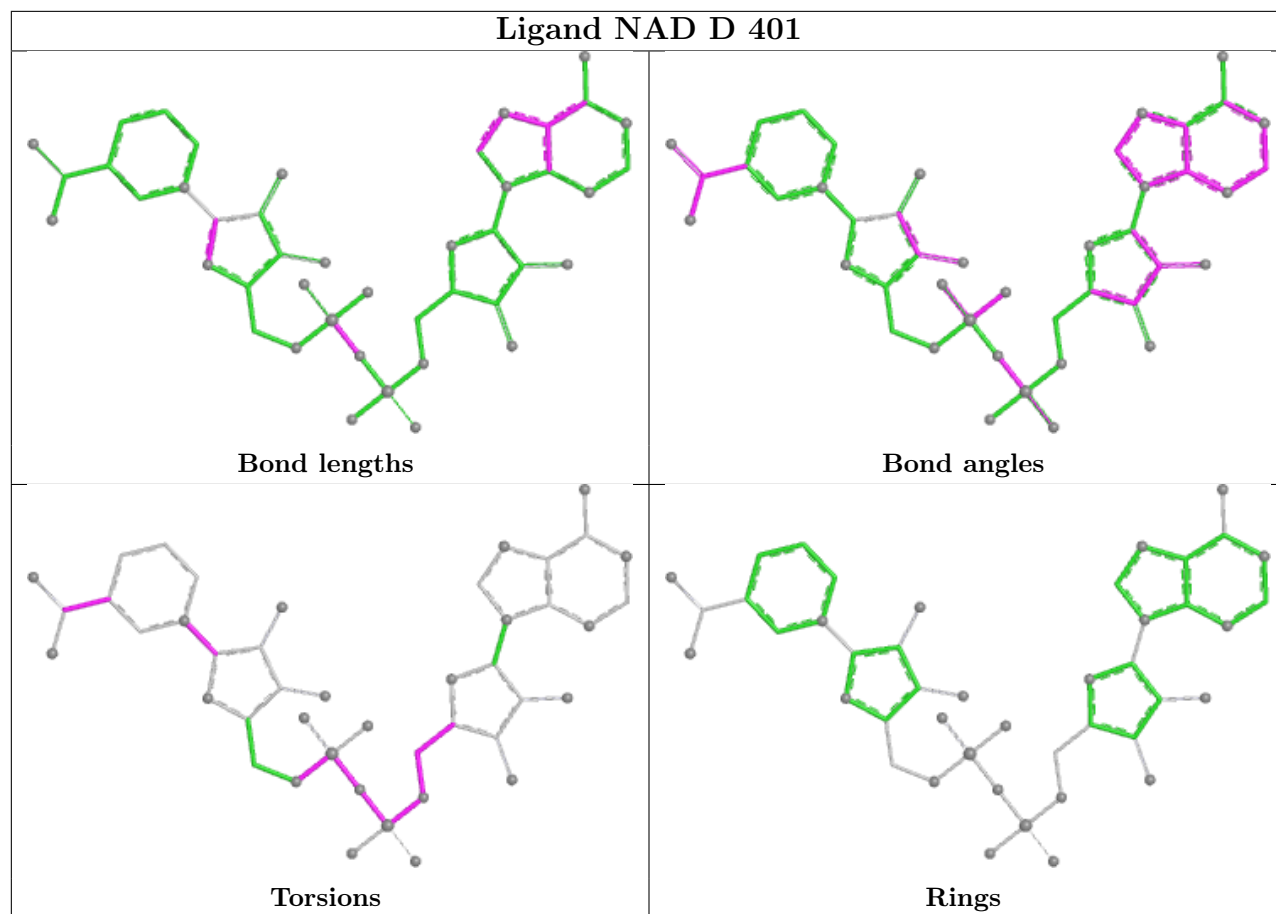


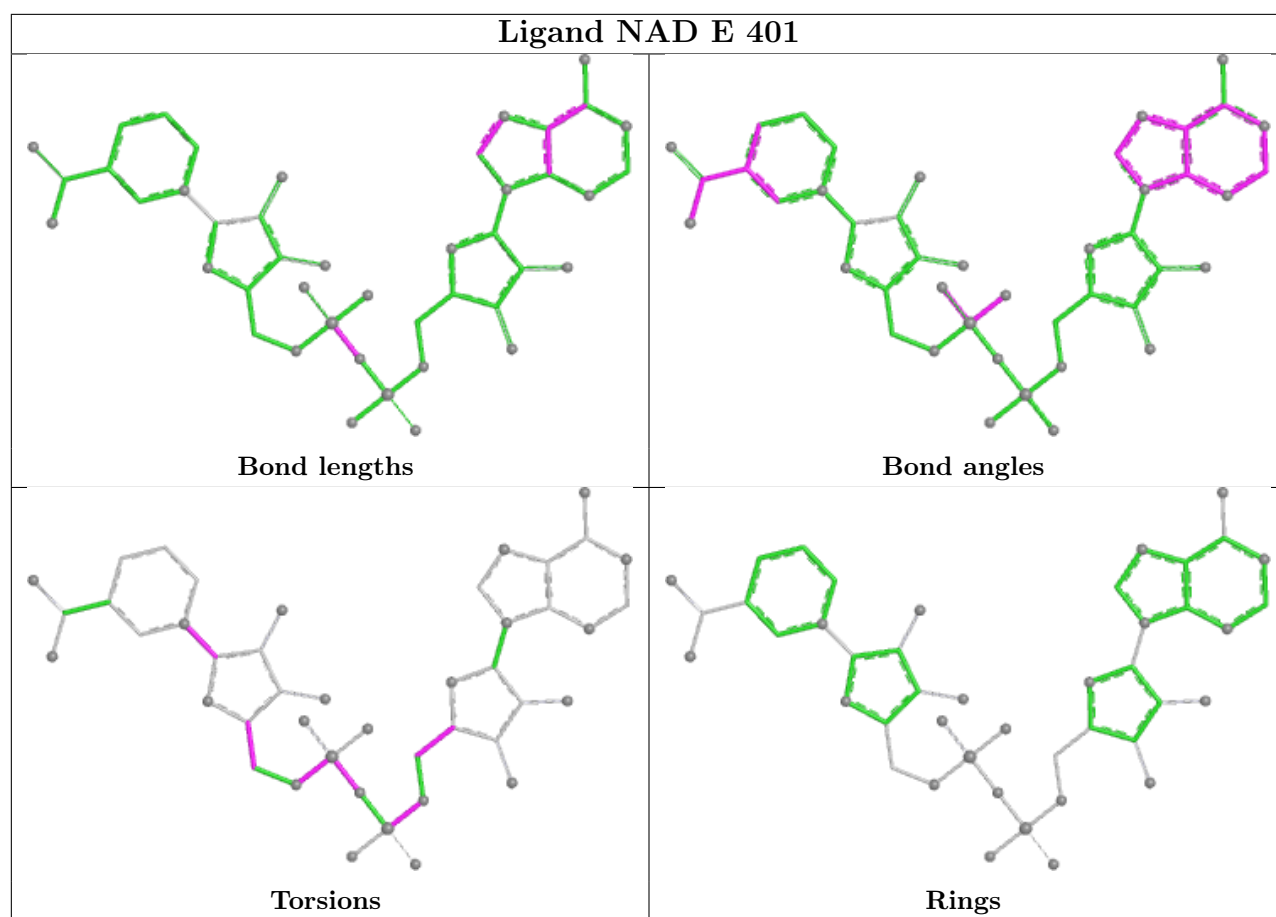












5.7 Other polymers [i](#)

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	330/355 (92%)	-0.54	1 (0%) 90 89	35, 59, 92, 123	1 (0%)
1	B	331/355 (93%)	-0.60	0 100 100	36, 58, 93, 134	1 (0%)
1	C	330/355 (92%)	-0.67	1 (0%) 90 89	30, 47, 80, 126	0
1	K	330/355 (92%)	-0.56	0 100 100	34, 59, 92, 116	1 (0%)
1	N	330/355 (92%)	-0.08	2 (0%) 85 82	46, 97, 150, 182	1 (0%)
2	D	331/355 (93%)	-0.53	1 (0%) 90 89	20, 65, 97, 153	3 (0%)
2	L	331/355 (93%)	-0.11	3 (0%) 81 76	43, 97, 163, 193	1 (0%)
3	E	329/355 (92%)	-0.48	1 (0%) 90 89	25, 62, 115, 160	2 (0%)
4	F	329/355 (92%)	-0.11	3 (0%) 81 76	42, 86, 123, 159	3 (0%)
4	J	328/355 (92%)	-0.37	1 (0%) 90 89	32, 80, 126, 158	1 (0%)
4	O	329/355 (92%)	-0.50	1 (0%) 90 89	42, 66, 106, 126	0
5	G	329/355 (92%)	-0.48	0 100 100	34, 63, 102, 133	2 (0%)
6	H	328/355 (92%)	-0.35	1 (0%) 90 89	43, 70, 102, 134	0
7	I	327/355 (92%)	-0.33	0 100 100	33, 80, 123, 158	1 (0%)
8	M	330/355 (92%)	-0.55	0 100 100	40, 59, 98, 129	1 (0%)
9	P	327/355 (92%)	-0.23	1 (0%) 90 89	45, 83, 123, 159	1 (0%)
All	All	5269/5680 (92%)	-0.41	16 (0%) 90 89	20, 69, 124, 193	19 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	78[A]	ARG	3.3
2	L	191[A]	HIS	3.1
1	N	191[A]	HIS	2.8
2	L	93	VAL	2.7
2	L	67	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
4	O	212	ALA	2.6
4	J	9	GLY	2.4
1	N	161	VAL	2.3
4	F	260	VAL	2.3
3	E	25	SER	2.3
2	D	130	PHE	2.3
4	F	191[A]	HIS	2.2
9	P	82	ASN	2.1
1	A	191[A]	HIS	2.0
6	H	212	ALA	2.0
1	C	212	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
6	SCY	H	150	9/10	0.76	0.19	58,73,109,115	0
9	ALY	P	249	12/13	0.76	0.12	98,125,137,140	0
9	SCY	P	150	9/10	0.79	0.13	68,81,91,105	0
7	ALY	I	249	12/13	0.84	0.12	107,139,150,153	0
4	ALY	F	46	12/13	0.86	0.15	58,66,79,79	0
4	ALY	J	257	12/13	0.86	0.11	88,97,100,100	0
5	ALY	G	46	12/13	0.87	0.14	56,63,75,75	0
4	ALY	F	257	12/13	0.87	0.11	56,79,84,84	0
1	ALY	N	46	12/13	0.88	0.12	59,65,74,75	0
4	ALY	O	257	12/13	0.89	0.12	61,67,88,91	0
1	ALY	K	46	12/13	0.89	0.12	54,62,68,71	0
5	SCY	G	150	9/10	0.89	0.12	52,58,94,96	0
7	ALY	I	257	12/13	0.90	0.10	80,94,97,97	0
6	ALY	H	257	12/13	0.90	0.10	70,75,80,85	0
4	ALY	O	46	12/13	0.90	0.10	63,70,90,91	0
8	ALY	M	257	12/13	0.91	0.09	55,61,82,86	0
9	ALY	P	46	12/13	0.91	0.09	46,51,60,66	0
3	ALY	E	46	12/13	0.91	0.11	53,65,80,84	0
6	ALY	H	46	12/13	0.91	0.10	41,48,56,58	0
7	ALY	I	46	12/13	0.92	0.09	49,59,62,63	0
3	ALY	E	249	12/13	0.92	0.10	60,78,121,128	0
1	ALY	A	46	12/13	0.92	0.09	55,60,66,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	ALY	B	46	12/13	0.93	0.09	34,39,55,60	0
1	ALY	C	46	12/13	0.93	0.09	35,42,49,52	0
4	ALY	J	46	12/13	0.93	0.07	35,49,57,60	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
12	PO4	A	408	5/5	0.53	0.24	45,48,51,75	5
12	PO4	H	403	5/5	0.53	0.22	53,56,60,80	5
15	ACT	H	404	4/4	0.53	0.24	65,67,79,80	4
12	PO4	F	402	5/5	0.62	0.16	49,52,67,76	5
12	PO4	P	405	5/5	0.64	0.16	45,52,58,76	5
12	PO4	P	404	5/5	0.66	0.19	43,45,46,69	5
10	NA	M	401	1/1	0.67	0.13	85,85,85,85	0
19	UVW	M	404	8/8	0.69	0.09	83,101,127,139	0
12	PO4	L	402	5/5	0.72	0.17	44,53,61,73	5
12	PO4	J	404	5/5	0.74	0.15	40,42,49,65	5
12	PO4	J	403	5/5	0.74	0.09	52,58,65,92	5
12	PO4	I	405	5/5	0.75	0.17	32,39,46,66	5
12	PO4	D	408	5/5	0.76	0.25	24,24,34,51	5
13	POP	B	410	9/9	0.76	0.09	54,91,117,118	0
13	POP	G	406	9/9	0.78	0.14	37,49,59,68	9
18	PG4	J	405	13/13	0.79	0.09	79,101,107,118	0
17	PEG	H	406	7/7	0.79	0.10	80,97,114,116	0
13	POP	M	405	9/9	0.80	0.12	30,41,54,56	9
11	CL	I	403	1/1	0.82	0.14	88,88,88,88	0
12	PO4	A	406	5/5	0.82	0.08	62,75,83,121	0
10	NA	I	402	1/1	0.82	0.09	91,91,91,91	0
11	CL	A	404	1/1	0.82	0.08	91,91,91,91	0
12	PO4	E	408	5/5	0.82	0.13	49,68,77,104	5
16	PGE	H	405	10/10	0.84	0.09	77,98,114,121	0
11	CL	B	405	1/1	0.84	0.08	82,82,82,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
14	NAD	C	401	44/44	0.85	0.12	52,86,101,113	0
11	CL	A	405	1/1	0.85	0.11	70,70,70,70	0
10	NA	E	402	1/1	0.87	0.08	79,79,79,79	0
11	CL	D	406	1/1	0.87	0.08	71,71,71,71	0
10	NA	G	401	1/1	0.87	0.09	78,78,78,78	0
13	POP	A	409	9/9	0.87	0.10	31,41,46,47	9
14	NAD	K	401	44/44	0.87	0.17	38,53,63,66	44
13	POP	F	403	9/9	0.88	0.11	41,57,65,74	9
11	CL	O	403	1/1	0.88	0.07	61,61,61,61	0
11	CL	G	402	1/1	0.89	0.07	87,87,87,87	0
11	CL	G	403	1/1	0.89	0.08	74,74,74,74	0
10	NA	N	401	1/1	0.89	0.15	89,89,89,89	0
14	NAD	D	401	44/44	0.89	0.17	40,56,69,76	44
11	CL	K	407	1/1	0.89	0.13	38,38,38,38	1
11	CL	B	407	1/1	0.90	0.09	81,81,81,81	0
12	PO4	D	409	5/5	0.90	0.25	62,75,97,102	0
11	CL	C	405	1/1	0.90	0.08	70,70,70,70	0
13	POP	J	406	9/9	0.90	0.10	28,40,44,48	9
14	NAD	P	401	44/44	0.90	0.14	34,51,58,62	44
11	CL	B	408	1/1	0.91	0.07	67,67,67,67	0
14	NAD	E	401	44/44	0.91	0.15	38,55,72,83	44
11	CL	D	407	1/1	0.91	0.07	66,66,66,66	0
14	NAD	O	401	44/44	0.91	0.15	34,58,85,88	44
10	NA	B	402	1/1	0.91	0.15	71,71,71,71	0
11	CL	J	402	1/1	0.92	0.07	69,69,69,69	0
11	CL	D	404	1/1	0.92	0.07	70,70,70,70	0
11	CL	N	402	1/1	0.92	0.06	65,65,65,65	0
10	NA	H	401	1/1	0.92	0.05	69,69,69,69	0
11	CL	P	403	1/1	0.92	0.07	59,59,59,59	0
11	CL	H	402	1/1	0.93	0.08	76,76,76,76	0
11	CL	E	407	1/1	0.93	0.06	73,73,73,73	0
12	PO4	B	409	5/5	0.93	0.10	27,34,38,38	5
14	NAD	I	401	44/44	0.93	0.13	32,49,62,69	44
11	CL	F	401	1/1	0.93	0.05	79,79,79,79	0
11	CL	K	405	1/1	0.93	0.08	59,59,59,59	0
11	CL	A	403	1/1	0.94	0.06	71,71,71,71	0
11	CL	A	402	1/1	0.95	0.09	66,66,66,66	0
11	CL	O	402	1/1	0.95	0.13	93,93,93,93	0
11	CL	B	406	1/1	0.95	0.05	80,80,80,80	0
11	CL	E	404	1/1	0.95	0.05	73,73,73,73	0
11	CL	C	407	1/1	0.95	0.16	54,54,54,54	1
11	CL	K	403	1/1	0.95	0.05	64,64,64,64	0

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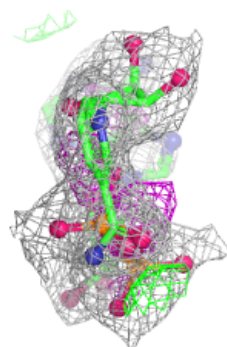
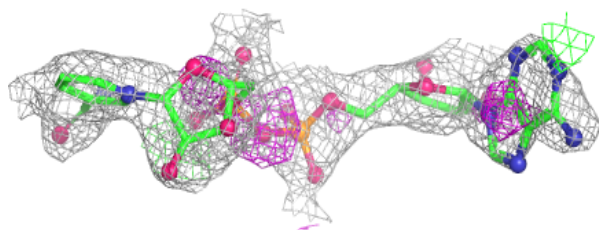
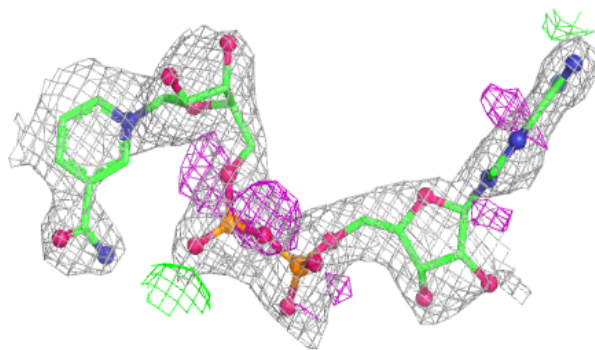
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
11	CL	D	403	1/1	0.95	0.05	71,71,71,71	0
11	CL	K	406	1/1	0.95	0.09	70,70,70,70	0
10	NA	B	401	1/1	0.95	0.06	53,53,53,53	0
11	CL	K	408	1/1	0.95	0.12	86,86,86,86	0
11	CL	L	401	1/1	0.95	0.06	53,53,53,53	0
11	CL	M	402	1/1	0.95	0.06	57,57,57,57	0
10	NA	C	402	1/1	0.96	0.04	67,67,67,67	0
11	CL	E	405	1/1	0.96	0.08	79,79,79,79	0
11	CL	G	404	1/1	0.96	0.08	75,75,75,75	0
12	PO4	A	407	5/5	0.96	0.10	36,39,67,71	5
11	CL	M	403	1/1	0.96	0.10	68,68,68,68	0
10	NA	D	402	1/1	0.96	0.04	65,65,65,65	0
11	CL	E	403	1/1	0.96	0.06	60,60,60,60	0
11	CL	B	404	1/1	0.97	0.06	71,71,71,71	0
11	CL	I	404	1/1	0.97	0.11	67,67,67,67	0
11	CL	J	401	1/1	0.97	0.06	67,67,67,67	0
10	NA	A	401	1/1	0.97	0.04	54,54,54,54	0
11	CL	C	404	1/1	0.97	0.06	57,57,57,57	0
11	CL	D	405	1/1	0.97	0.04	73,73,73,73	0
11	CL	B	403	1/1	0.97	0.06	51,51,51,51	0
11	CL	K	404	1/1	0.98	0.04	51,51,51,51	0
11	CL	P	402	1/1	0.98	0.04	57,57,57,57	0
11	CL	C	403	1/1	0.98	0.09	52,52,52,52	0
11	CL	G	405	1/1	0.98	0.06	46,46,46,46	0
11	CL	C	406	1/1	0.98	0.04	46,46,46,46	0
10	NA	K	402	1/1	0.98	0.09	59,59,59,59	0
11	CL	E	406	1/1	0.99	0.04	40,40,40,40	0

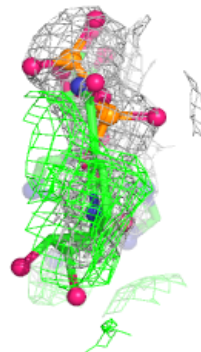
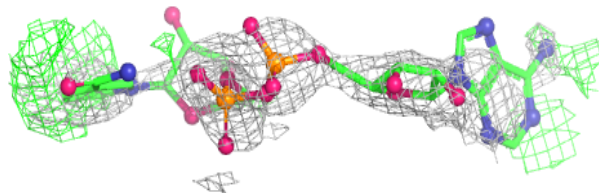
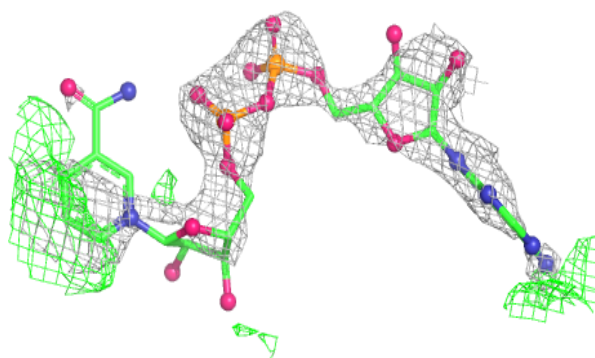
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD 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)

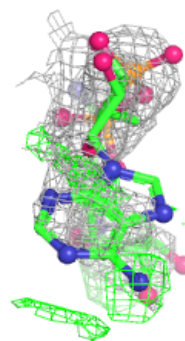
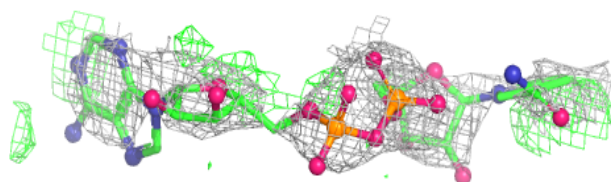
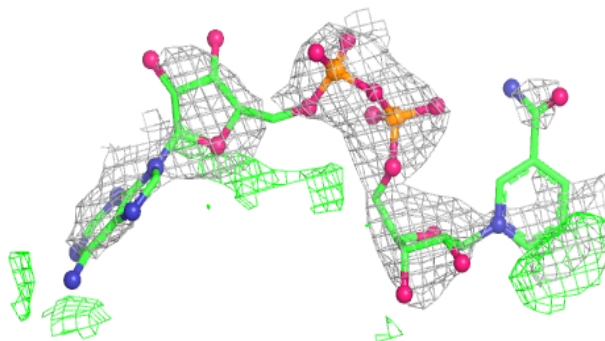
**Electron density around NAD K 401:**

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

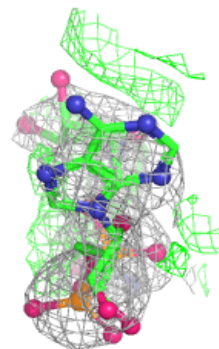
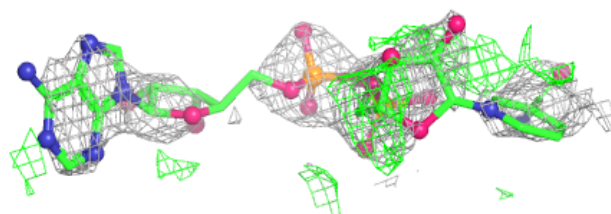
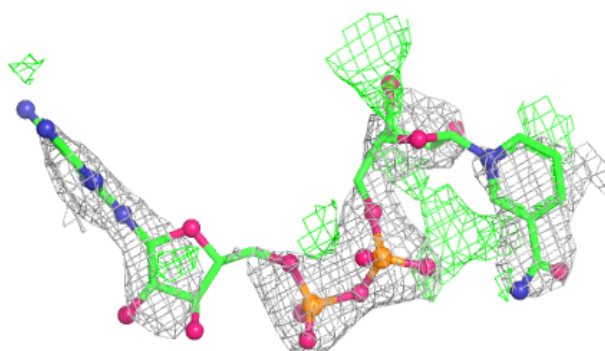


Electron density around NAD D 401:

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

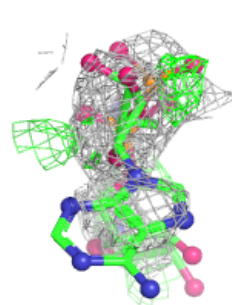
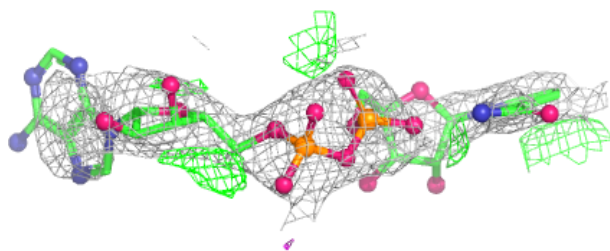
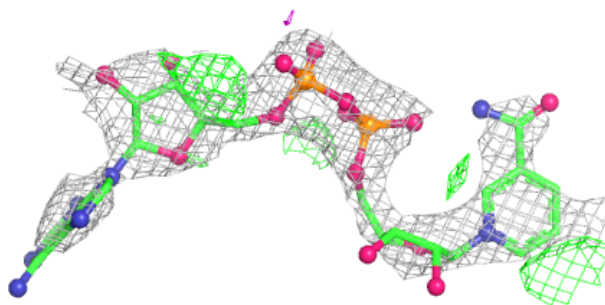
**Electron density around NAD P 401:**

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

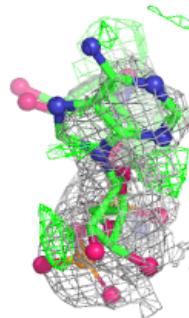
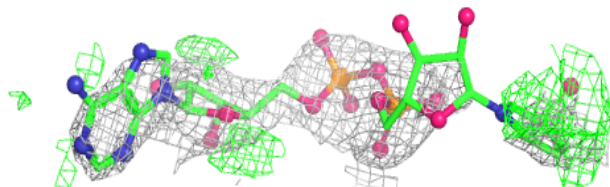
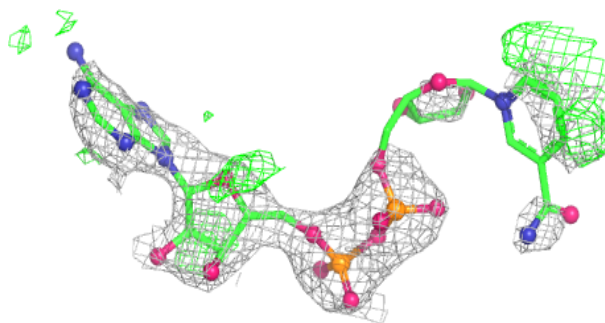


Electron density around NAD E 401:

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

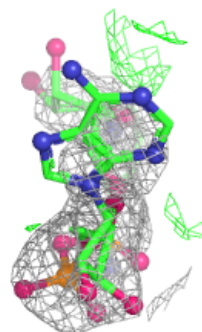
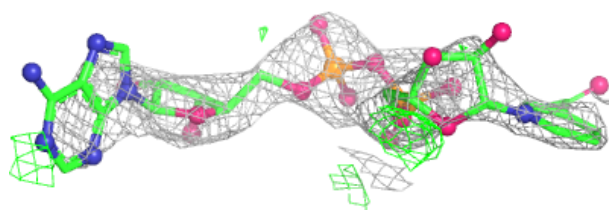
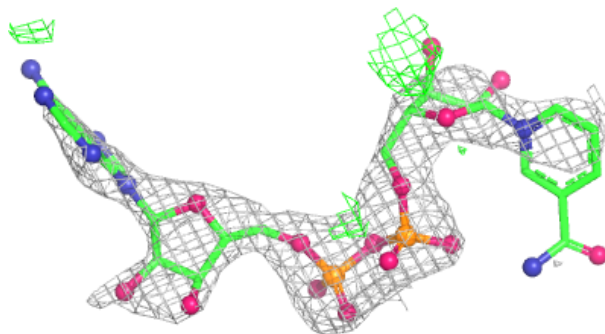
**Electron density around NAD O 401:**

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



Electron density around NAD 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.