Full wwPDB X-ray Structure Validation Report (i)

Nov 13, 2023 – 10:29 PM JST

PDB ID	:	5YA8
Title	:	Crystal structure of scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity complexed with myo-inositol
Authors	:	Fukano, K.; Shimizu, T.; Sasaki, Y.; Nakamura, A.; Yajima, S.
Deposited on	:	2017-08-31
Resolution	:	2.30 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 (i) 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 (i)) were used in the production of this report:

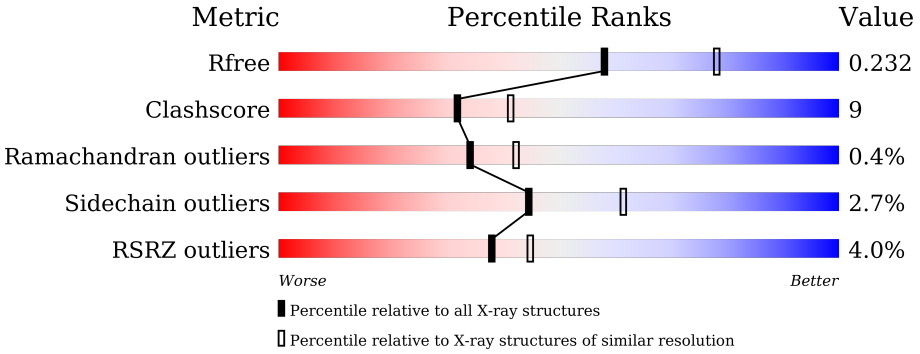
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	380	3% 83% 12%
1	B	380	3% 82% 13%
1	C	380	7% 82% 13%
1	D	380	2% 81% 14%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 11763 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 Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	366	2785	1744	510	513	18	0	0	0
1	B	366	2785	1744	510	513	18	0	0	0
1	C	366	2785	1744	510	513	18	0	0	0
1	D	365	2777	1738	509	512	18	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	72	SER	ASN	engineered mutation	UNP K7ZP76
A	373	LEU	-	expression tag	UNP K7ZP76
A	374	GLU	-	expression tag	UNP K7ZP76
A	375	HIS	-	expression tag	UNP K7ZP76
A	376	HIS	-	expression tag	UNP K7ZP76
A	377	HIS	-	expression tag	UNP K7ZP76
A	378	HIS	-	expression tag	UNP K7ZP76
A	379	HIS	-	expression tag	UNP K7ZP76
A	380	HIS	-	expression tag	UNP K7ZP76
B	72	SER	ASN	engineered mutation	UNP K7ZP76
B	373	LEU	-	expression tag	UNP K7ZP76
B	374	GLU	-	expression tag	UNP K7ZP76
B	375	HIS	-	expression tag	UNP K7ZP76
B	376	HIS	-	expression tag	UNP K7ZP76
B	377	HIS	-	expression tag	UNP K7ZP76
B	378	HIS	-	expression tag	UNP K7ZP76
B	379	HIS	-	expression tag	UNP K7ZP76
B	380	HIS	-	expression tag	UNP K7ZP76
C	72	SER	ASN	engineered mutation	UNP K7ZP76
C	373	LEU	-	expression tag	UNP K7ZP76

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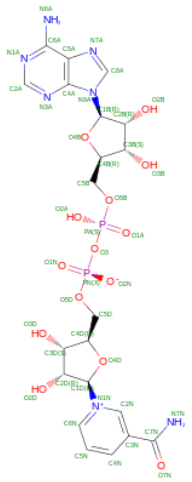
Chain	Residue	Modelled	Actual	Comment	Reference
C	374	GLU	-	expression tag	UNP K7ZP76
C	375	HIS	-	expression tag	UNP K7ZP76
C	376	HIS	-	expression tag	UNP K7ZP76
C	377	HIS	-	expression tag	UNP K7ZP76
C	378	HIS	-	expression tag	UNP K7ZP76
C	379	HIS	-	expression tag	UNP K7ZP76
C	380	HIS	-	expression tag	UNP K7ZP76
D	72	SER	ASN	engineered mutation	UNP K7ZP76
D	373	LEU	-	expression tag	UNP K7ZP76
D	374	GLU	-	expression tag	UNP K7ZP76
D	375	HIS	-	expression tag	UNP K7ZP76
D	376	HIS	-	expression tag	UNP K7ZP76
D	377	HIS	-	expression tag	UNP K7ZP76
D	378	HIS	-	expression tag	UNP K7ZP76
D	379	HIS	-	expression tag	UNP K7ZP76
D	380	HIS	-	expression tag	UNP K7ZP76

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

NAD



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

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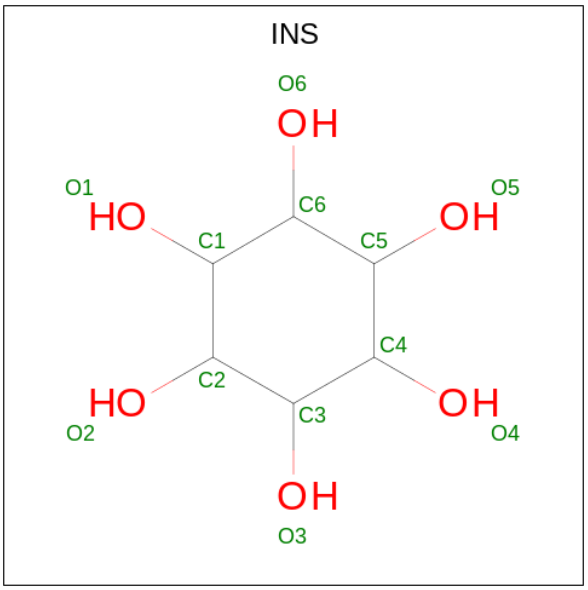
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	D	1	44	21	7	14	2	0	0

- Molecule 3 is 1,2,3,4,5,6-HEXAHYDROXY-CYCLOHEXANE (three-letter code: INS) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	12	6	6	0	0
3	B	1	12	6	6	0	0
3	D	1	12	6	6	0	0

- Molecule 4 is water.

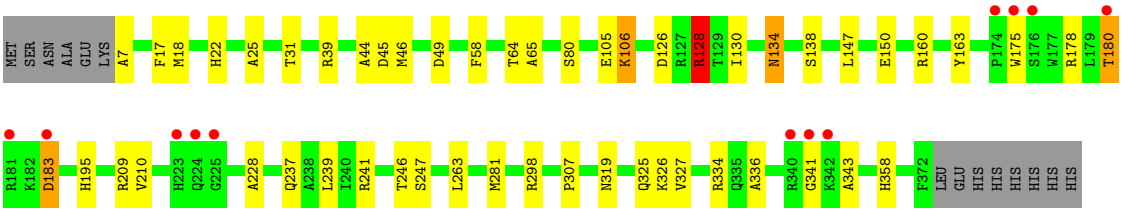
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	104	104	104	0	0
4	B	116	116	116	0	0
4	C	84	84	84	0	0
4	D	115	115	115	0	0

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.

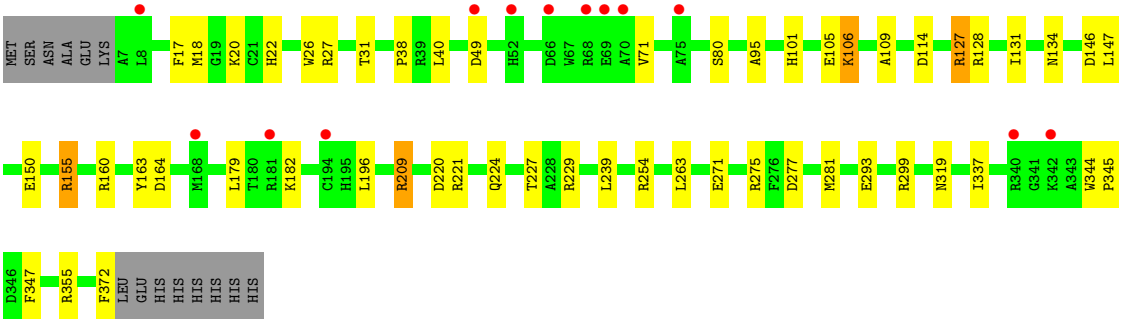
• Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity

Chain A:



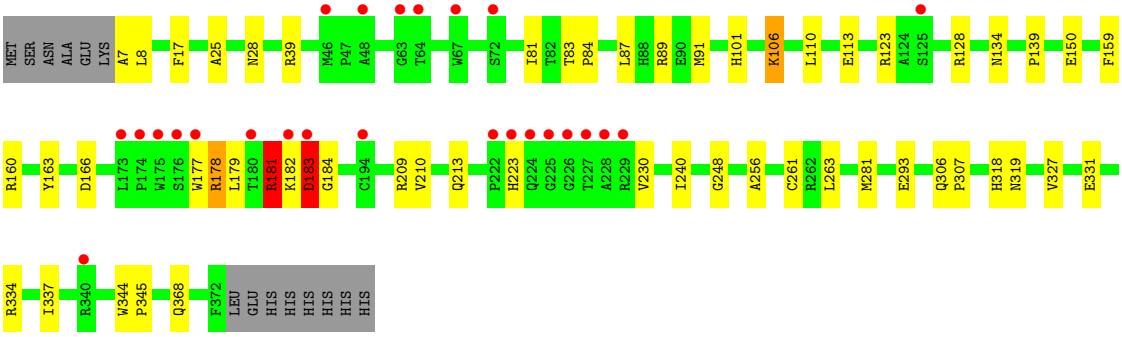
• Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity

Chain B:



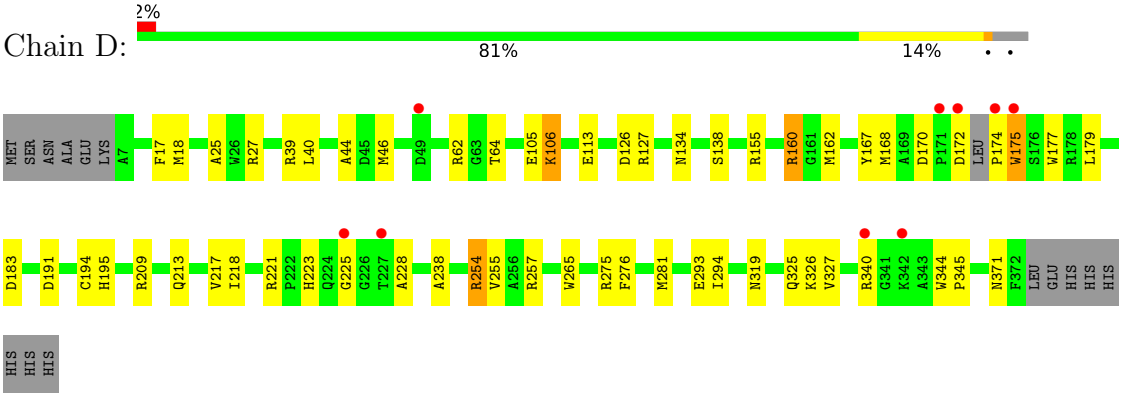
• Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity

Chain C:



• Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	86.21Å 120.15Å 140.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.30 – 2.30 49.29 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.0 (49.30-2.30) 98.0 (49.29-2.30)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.192 , 0.224 0.198 , 0.232	Depositor DCC
R_{free} test set	3208 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	24.9	Xtriage
Anisotropy	0.430	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 44.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11763	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.69	0/2851	0.87	9/3865 (0.2%)
1	B	0.70	0/2851	0.86	7/3865 (0.2%)
1	C	0.70	0/2851	0.83	2/3865 (0.1%)
1	D	0.70	0/2842	0.87	8/3850 (0.2%)
All	All	0.70	0/11395	0.86	26/15445 (0.2%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	257	ARG	NE-CZ-NH2	-8.51	116.05	120.30
1	A	128	ARG	NE-CZ-NH1	7.19	123.90	120.30
1	A	178	ARG	NE-CZ-NH1	7.09	123.85	120.30
1	D	275	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	B	209	ARG	NE-CZ-NH2	-6.17	117.22	120.30
1	B	277	ASP	CB-CG-OD1	6.16	123.85	118.30
1	A	178	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	D	160	ARG	NE-CZ-NH1	6.13	123.37	120.30
1	B	299	ARG	NE-CZ-NH1	6.10	123.35	120.30
1	D	275	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	A	128	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	C	160	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	D	257	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	B	155	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	A	160	ARG	NE-CZ-NH2	-5.73	117.43	120.30
1	C	181	ARG	NE-CZ-NH1	5.71	123.15	120.30
1	D	254	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	D	160	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	D	39	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	A	241	ARG	NE-CZ-NH2	-5.30	117.65	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	254	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	A	334	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	A	160	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	B	160	ARG	NE-CZ-NH2	-5.11	117.74	120.30
1	B	275	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	A	183	ASP	N-CA-C	-5.04	97.40	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2785	0	2702	46	0
1	B	2785	0	2702	41	1
1	C	2785	0	2702	51	1
1	D	2777	0	2691	58	0
2	A	44	0	26	9	0
2	B	44	0	26	6	0
2	C	44	0	26	3	0
2	D	44	0	26	12	0
3	A	12	0	12	4	0
3	B	12	0	12	5	0
3	D	12	0	12	5	0
4	A	104	0	0	15	0
4	B	116	0	0	12	0
4	C	84	0	0	21	0
4	D	115	0	0	16	0
All	All	11763	0	10937	199	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (199) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:HIS:HB3	4:A:578:HOH:O	1.33	1.29
4:A:537:HOH:O	1:D:294:ILE:HB	1.49	1.09
1:B:196:LEU:HA	4:B:558:HOH:O	1.56	1.05
1:C:181:ARG:HD3	4:C:606:HOH:O	1.56	1.04
1:C:178:ARG:HB3	4:C:649:HOH:O	1.59	1.03
1:B:196:LEU:HD12	4:B:558:HOH:O	1.60	1.01
1:A:58:PHE:HA	4:B:502:HOH:O	1.61	1.00
1:C:178:ARG:CB	4:C:649:HOH:O	2.13	0.96
1:D:191:ASP:HB3	4:D:518:HOH:O	1.67	0.95
2:B:401:NAD:C4N	3:B:402:INS:H4	1.99	0.93
1:D:46:MET:SD	4:D:507:HOH:O	2.27	0.92
2:C:500:NAD:H2B	4:C:638:HOH:O	1.69	0.91
1:D:113:GLU:HG3	4:D:596:HOH:O	1.71	0.90
2:A:401:NAD:C4N	3:A:402:INS:H4	2.03	0.88
1:B:155:ARG:HD3	4:B:512:HOH:O	1.75	0.87
1:B:372:PHE:C	4:B:505:HOH:O	2.16	0.82
1:C:182:LYS:O	1:C:184:GLY:N	2.13	0.82
1:C:223:HIS:HD2	4:C:678:HOH:O	1.63	0.81
1:D:162:MET:SD	4:D:611:HOH:O	2.40	0.80
1:B:101:HIS:HD2	1:B:128:ARG:H	1.31	0.79
1:D:276:PHE:CD2	4:D:506:HOH:O	2.37	0.77
1:A:263:LEU:HD11	4:A:578:HOH:O	1.87	0.75
2:D:401:NAD:C4N	3:D:402:INS:H4	2.17	0.74
2:D:401:NAD:C3N	3:D:402:INS:O4	2.37	0.73
1:C:306:GLN:HB3	4:C:603:HOH:O	1.90	0.72
1:C:83:THR:OG1	1:C:91:MET:CE	2.38	0.71
1:D:209:ARG:HG2	4:D:529:HOH:O	1.90	0.71
1:B:146:ASP:O	1:B:150:GLU:HG3	1.91	0.70
1:C:139:PRO:CA	4:C:603:HOH:O	2.40	0.70
2:B:401:NAD:C3N	3:B:402:INS:O4	2.39	0.70
2:B:401:NAD:C4N	3:B:402:INS:C4	2.70	0.70
1:D:175:TRP:HE3	1:D:221:ARG:HG3	1.58	0.69
1:D:276:PHE:CG	4:D:506:HOH:O	2.45	0.69
1:D:126:ASP:O	1:D:127:ARG:NH1	2.28	0.66
1:A:180:THR:O	1:A:183:ASP:O	2.13	0.66
1:B:106:LYS:HE2	3:B:402:INS:O4	1.96	0.65
1:C:83:THR:OG1	1:C:91:MET:HE1	1.95	0.65
1:C:139:PRO:HB3	4:C:603:HOH:O	1.96	0.64
1:A:147:LEU:HD21	1:D:293:GLU:HG2	1.80	0.64
1:C:139:PRO:HA	4:C:603:HOH:O	1.98	0.64
1:D:265:TRP:CZ2	4:D:506:HOH:O	2.50	0.64
1:A:263:LEU:CD1	4:A:578:HOH:O	2.45	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:183:ASP:HB2	4:C:656:HOH:O	1.98	0.63
1:C:182:LYS:C	1:C:184:GLY:H	1.99	0.63
1:A:307:PRO:HA	4:A:600:HOH:O	1.99	0.62
1:D:175:TRP:CE3	1:D:221:ARG:HG3	2.34	0.62
1:D:106:LYS:HD2	1:D:106:LYS:C	2.20	0.61
1:B:71:VAL:CG1	1:B:95:ALA:HA	2.30	0.61
1:A:175:TRP:CG	4:A:561:HOH:O	2.54	0.61
1:C:177:TRP:CE3	1:C:183:ASP:HB3	2.36	0.61
2:A:401:NAD:C3N	3:A:402:INS:O4	2.48	0.61
1:A:18:MET:HE1	1:A:134:ASN:OD1	2.02	0.60
1:C:83:THR:HB	1:C:84:PRO:HD2	1.83	0.60
1:B:31:THR:HG23	4:B:502:HOH:O	2.01	0.60
1:C:8:LEU:HD21	1:C:337:ILE:HG21	1.83	0.60
2:D:401:NAD:C6A	4:D:507:HOH:O	2.49	0.60
1:C:83:THR:OG1	1:C:91:MET:HE3	2.02	0.60
1:B:271:GLU:HG3	4:B:611:HOH:O	2.02	0.59
1:D:18:MET:HG2	2:D:401:NAD:C5N	2.32	0.59
1:D:209:ARG:HD2	1:D:371:ASN:OD1	2.02	0.59
2:A:401:NAD:C4N	3:A:402:INS:C4	2.81	0.58
1:C:178:ARG:HB2	4:C:649:HOH:O	1.93	0.58
2:D:401:NAD:C4N	3:D:402:INS:C4	2.82	0.57
1:A:18:MET:CE	2:A:401:NAD:N7N	2.68	0.57
1:A:175:TRP:CD1	4:A:561:HOH:O	2.53	0.57
2:B:401:NAD:C5N	3:B:402:INS:H4	2.34	0.57
1:A:298:ARG:CD	4:A:537:HOH:O	2.53	0.57
1:B:229:ARG:HD3	4:B:587:HOH:O	2.04	0.56
1:C:81:ILE:HG23	1:C:91:MET:HE3	1.87	0.56
1:A:18:MET:HE2	2:A:401:NAD:N7N	2.21	0.56
1:A:31:THR:O	1:B:20:LYS:HE3	2.06	0.56
1:A:138:SER:HB3	1:A:325:GLN:HE21	1.70	0.55
1:B:22:HIS:HE1	1:B:105:GLU:OE2	1.89	0.55
1:C:281:MET:HG2	1:D:319:ASN:HD22	1.71	0.55
1:A:106:LYS:HE2	3:A:402:INS:O4	2.07	0.54
1:C:139:PRO:CB	4:C:603:HOH:O	2.53	0.54
1:D:106:LYS:HD3	1:D:194:CYS:SG	2.47	0.54
1:A:281:MET:HG2	1:B:319:ASN:HD22	1.72	0.54
1:D:172:ASP:HA	1:D:223:HIS:HD2	1.73	0.53
1:A:18:MET:CE	2:A:401:NAD:C7N	2.87	0.53
1:A:22:HIS:HE1	1:A:105:GLU:OE2	1.92	0.53
1:C:166:ASP:HB3	1:C:256:ALA:O	2.09	0.53
2:D:401:NAD:O5D	2:D:401:NAD:H6N	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:500:NAD:C8A	4:C:638:HOH:O	2.57	0.53
1:D:168:MET:SD	1:D:179:LEU:HD21	2.48	0.52
1:D:265:TRP:CE2	4:D:506:HOH:O	2.62	0.52
1:C:344:TRP:CE3	1:C:345:PRO:HD3	2.45	0.52
1:A:25:ALA:HA	1:A:327:VAL:HG22	1.91	0.52
1:D:27:ARG:HG2	1:D:40:LEU:HD11	1.91	0.52
1:B:372:PHE:N	4:B:505:HOH:O	2.42	0.52
1:D:168:MET:HG3	1:D:254:ARG:HD2	1.91	0.52
1:A:18:MET:HE3	1:A:105:GLU:OE2	2.10	0.51
1:B:196:LEU:CD1	4:B:558:HOH:O	2.36	0.51
1:A:209:ARG:NE	4:A:501:HOH:O	2.32	0.51
1:A:7:ALA:O	1:A:39:ARG:NH1	2.44	0.51
1:C:318:HIS:NE2	3:D:402:INS:O2	2.40	0.51
1:B:182:LYS:HB2	4:B:541:HOH:O	2.10	0.50
1:D:265:TRP:NE1	4:D:506:HOH:O	2.44	0.50
1:B:71:VAL:HG13	1:B:95:ALA:HA	1.94	0.50
1:A:336:ALA:HA	1:A:341:GLY:O	2.13	0.49
1:C:209:ARG:HG2	1:C:210:VAL:N	2.27	0.49
1:D:167:TYR:CE1	1:D:168:MET:HG2	2.47	0.49
1:D:172:ASP:C	1:D:174:PRO:HG3	2.33	0.49
4:A:537:HOH:O	1:D:294:ILE:HD12	2.13	0.49
1:B:27:ARG:HD2	1:B:40:LEU:HD11	1.93	0.49
1:A:319:ASN:HD22	1:B:281:MET:HG2	1.77	0.49
1:A:126:ASP:O	1:A:128:ARG:NH2	2.45	0.48
1:C:101:HIS:CD2	1:C:337:ILE:HG12	2.48	0.48
2:C:500:NAD:H8A	4:C:638:HOH:O	2.13	0.48
1:C:209:ARG:CZ	4:C:618:HOH:O	2.61	0.48
1:D:217:VAL:HG21	1:D:255:VAL:HG21	1.95	0.48
1:A:228:ALA:CB	4:A:561:HOH:O	2.62	0.47
1:C:319:ASN:OD1	1:D:281:MET:HG2	2.14	0.47
1:D:25:ALA:HA	1:D:327:VAL:HG22	1.96	0.47
1:D:46:MET:CG	4:D:507:HOH:O	2.59	0.47
1:A:130:ILE:HD12	1:A:343:ALA:HB1	1.95	0.47
1:C:110:LEU:HD11	1:C:183:ASP:HA	1.95	0.47
1:B:105:GLU:OE1	2:B:401:NAD:H2N	2.15	0.47
1:D:170:ASP:CG	4:D:547:HOH:O	2.52	0.47
1:A:22:HIS:HD2	1:A:80:SER:OG	1.98	0.46
1:B:18:MET:HG3	2:B:401:NAD:C2N	2.45	0.46
1:B:22:HIS:HD2	1:B:80:SER:OG	1.99	0.46
1:C:106:LYS:O	1:C:106:LYS:HE3	2.15	0.46
1:C:123:ARG:HD2	4:C:674:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:VAL:HG12	1:C:240:ILE:HD12	1.97	0.46
1:C:87:LEU:HB3	1:C:91:MET:HE1	1.97	0.46
1:C:179:LEU:HD13	1:C:230:VAL:HG22	1.95	0.46
1:C:7:ALA:O	1:C:39:ARG:NH1	2.49	0.46
1:C:331:GLU:OE1	1:C:334:ARG:NH1	2.47	0.46
1:D:105:GLU:OE1	2:D:401:NAD:H2N	2.16	0.46
1:B:179:LEU:HD13	1:B:221:ARG:HD2	1.96	0.45
1:A:228:ALA:HB3	4:A:561:HOH:O	2.15	0.45
1:D:18:MET:HG2	2:D:401:NAD:C6N	2.47	0.45
1:A:106:LYS:HD2	1:A:106:LYS:C	2.37	0.45
1:C:306:GLN:CB	4:C:603:HOH:O	2.56	0.45
1:A:209:ARG:NH2	4:A:501:HOH:O	2.50	0.44
1:C:25:ALA:HA	1:C:327:VAL:HG22	1.99	0.44
1:D:138:SER:HB3	1:D:325:GLN:HE21	1.82	0.44
1:C:368:GLN:HG3	4:C:648:HOH:O	2.17	0.44
1:D:18:MET:CE	1:D:326:LYS:HE2	2.47	0.44
1:D:46:MET:HG3	4:D:507:HOH:O	2.18	0.44
1:B:101:HIS:CE1	1:B:337:ILE:HG23	2.53	0.44
1:C:307:PRO:N	4:C:603:HOH:O	2.51	0.44
1:D:218:ILE:HG21	1:D:221:ARG:HH21	1.82	0.44
1:A:44:ALA:HA	1:A:64:THR:O	2.17	0.44
1:A:298:ARG:HD2	4:A:537:HOH:O	2.15	0.44
1:D:18:MET:HG2	2:D:401:NAD:C4N	2.47	0.44
1:B:147:LEU:HD21	1:C:293:GLU:HG2	2.00	0.44
1:D:179:LEU:CD1	1:D:221:ARG:HD2	2.48	0.44
1:D:106:LYS:HE2	1:D:195:HIS:NE2	2.33	0.43
1:D:126:ASP:C	1:D:127:ARG:HH11	2.21	0.43
2:D:401:NAD:C7N	3:D:402:INS:O4	2.65	0.43
1:C:368:GLN:CD	4:C:648:HOH:O	2.56	0.43
1:A:18:MET:HE1	2:A:401:NAD:N7N	2.32	0.43
1:A:209:ARG:HD2	1:A:210:VAL:N	2.33	0.43
1:A:18:MET:HE2	2:A:401:NAD:C7N	2.48	0.43
1:B:127:ARG:C	1:B:128:ARG:HH11	2.21	0.43
1:D:106:LYS:HE3	2:D:401:NAD:C2N	2.49	0.43
1:A:106:LYS:HE2	1:A:195:HIS:NE2	2.34	0.43
1:D:168:MET:CE	1:D:179:LEU:HD21	2.49	0.43
1:B:127:ARG:HA	1:B:127:ARG:HE	1.83	0.43
1:D:155:ARG:HD2	4:D:545:HOH:O	2.19	0.43
1:C:179:LEU:HD23	4:C:649:HOH:O	2.19	0.42
1:A:45:ASP:O	1:A:65:ALA:HA	2.19	0.42
1:D:44:ALA:HA	1:D:64:THR:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:GLU:OE2	1:C:150:GLU:OE1	2.36	0.42
1:C:182:LYS:C	1:C:184:GLY:N	2.62	0.42
1:B:344:TRP:HA	1:B:345:PRO:HA	1.82	0.42
1:C:101:HIS:CD2	1:C:128:ARG:H	2.38	0.42
1:B:131:ILE:HG12	1:B:345:PRO:O	2.20	0.42
1:B:263:LEU:HD21	4:B:558:HOH:O	2.19	0.42
1:B:131:ILE:HD11	1:B:347:PHE:CD1	2.55	0.42
1:C:210:VAL:HG12	1:C:240:ILE:CD1	2.49	0.42
1:D:340:ARG:HB3	1:D:340:ARG:CZ	2.50	0.42
1:A:18:MET:HE1	2:A:401:NAD:C7N	2.50	0.41
1:A:18:MET:HE3	1:A:326:LYS:NZ	2.34	0.41
1:A:239:LEU:HD13	1:C:213:GLN:HG2	2.02	0.41
1:B:179:LEU:CD1	1:B:221:ARG:HD2	2.50	0.41
1:B:239:LEU:HD11	1:D:238:ALA:HA	2.02	0.41
1:C:91:MET:HE2	1:C:91:MET:HB2	1.75	0.41
1:C:159:PHE:O	1:C:248:GLY:HA2	2.19	0.41
1:B:26:TRP:HE3	1:B:38:PRO:HG2	1.84	0.41
1:D:175:TRP:CZ3	1:D:228:ALA:HB1	2.55	0.41
1:D:344:TRP:HA	1:D:345:PRO:HA	1.86	0.41
1:A:150:GLU:HG2	1:A:150:GLU:O	2.19	0.41
1:B:209:ARG:HH22	1:D:213:GLN:NE2	2.18	0.41
1:B:109:ALA:HB3	1:B:114:ASP:HB3	2.02	0.41
1:D:177:TRP:HB2	1:D:183:ASP:HB3	2.01	0.41
1:D:27:ARG:HG3	4:D:604:HOH:O	2.20	0.41
1:D:167:TYR:CD1	1:D:168:MET:HG2	2.56	0.41
1:B:164:ASP:OD2	1:D:160:ARG:NH2	2.54	0.41
1:A:358:HIS:HE1	4:A:523:HOH:O	2.04	0.41
1:C:223:HIS:O	1:C:223:HIS:CG	2.71	0.41
1:D:18:MET:HG3	2:D:401:NAD:C2N	2.51	0.40
1:D:27:ARG:CG	1:D:40:LEU:HD11	2.52	0.40
1:A:246:THR:HG22	1:A:247:SER:N	2.37	0.40
1:B:128:ARG:HH11	1:B:128:ARG:N	2.19	0.40
1:A:150:GLU:OE2	1:D:293:GLU:OE1	2.40	0.40
1:B:101:HIS:CD2	1:B:128:ARG:H	2.23	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:ARG:NH1	1:C:113:GLU:OE2[1_655]	2.18	0.02

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	364/380 (96%)	351 (96%)	13 (4%)	0	100	100
1	B	364/380 (96%)	352 (97%)	10 (3%)	2 (0%)	29	35
1	C	364/380 (96%)	346 (95%)	16 (4%)	2 (0%)	29	35
1	D	361/380 (95%)	348 (96%)	11 (3%)	2 (1%)	25	31
All	All	1453/1520 (96%)	1397 (96%)	50 (3%)	6 (0%)	34	42

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	227	THR
1	B	224	GLN
1	C	183	ASP
1	D	175	TRP
1	D	225	GLY
1	C	261	CYS

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	275/288 (96%)	266 (97%)	9 (3%)	38	53
1	B	275/288 (96%)	268 (98%)	7 (2%)	47	65
1	C	275/288 (96%)	265 (96%)	10 (4%)	35	49
1	D	274/288 (95%)	270 (98%)	4 (2%)	65	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1099/1152 (95%)	1069 (97%)	30 (3%)	44 61

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	PHE
1	A	46	MET
1	A	49	ASP
1	A	106	LYS
1	A	128	ARG
1	A	134	ASN
1	A	163	TYR
1	A	180	THR
1	A	237	GLN
1	B	17	PHE
1	B	49	ASP
1	B	106	LYS
1	B	127	ARG
1	B	134	ASN
1	B	163	TYR
1	B	220	ASP
1	C	17	PHE
1	C	28	ASN
1	C	89	ARG
1	C	106	LYS
1	C	134	ASN
1	C	163	TYR
1	C	178	ARG
1	C	181	ARG
1	C	183	ASP
1	C	263	LEU
1	D	17	PHE
1	D	62	ARG
1	D	106	LYS
1	D	134	ASN

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	22	HIS
1	A	28	ASN

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Mol	Chain	Res	Type
1	A	237	GLN
1	A	306	GLN
1	A	319	ASN
1	A	325	GLN
1	A	358	HIS
1	B	22	HIS
1	B	28	ASN
1	B	101	HIS
1	B	306	GLN
1	B	319	ASN
1	B	358	HIS
1	C	28	ASN
1	C	101	HIS
1	C	223	HIS
1	C	306	GLN
1	D	28	ASN
1	D	213	GLN
1	D	223	HIS
1	D	278	GLN
1	D	319	ASN
1	D	325	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	NAD	C	500	-	42,48,48	1.04	2 (4%)	50,73,73	1.53	9 (18%)
2	NAD	B	401	-	42,48,48	1.37	5 (11%)	50,73,73	1.57	8 (16%)
2	NAD	D	401	-	42,48,48	1.28	5 (11%)	50,73,73	1.34	7 (14%)
2	NAD	A	401	-	42,48,48	1.21	4 (9%)	50,73,73	1.33	6 (12%)
3	INS	B	402	-	12,12,12	0.86	0	18,18,18	2.61	8 (44%)
3	INS	A	402	-	12,12,12	1.02	0	18,18,18	2.37	6 (33%)
3	INS	D	402	-	12,12,12	0.86	0	18,18,18	2.29	5 (27%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	C	500	-	-	4/26/62/62	0/5/5/5
2	NAD	B	401	-	-	6/26/62/62	0/5/5/5
2	NAD	D	401	-	-	5/26/62/62	0/5/5/5
2	NAD	A	401	-	-	4/26/62/62	0/5/5/5
3	INS	B	402	-	-	-	0/1/1/1
3	INS	A	402	-	-	-	0/1/1/1
3	INS	D	402	-	-	-	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAD	C2N-N1N	-3.98	1.30	1.35
2	D	401	NAD	C2N-N1N	-3.82	1.30	1.35
2	D	401	NAD	C4N-C3N	3.44	1.45	1.39
2	C	500	NAD	O4B-C1B	3.22	1.45	1.41
2	B	401	NAD	C2A-N3A	3.02	1.37	1.32
2	A	401	NAD	C3N-C7N	2.79	1.54	1.50
2	B	401	NAD	O4D-C1D	2.71	1.44	1.41
2	A	401	NAD	O4D-C1D	2.67	1.44	1.41
2	A	401	NAD	C5A-C4A	2.58	1.47	1.40
2	D	401	NAD	C2A-N3A	2.32	1.35	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAD	C7N-N7N	2.21	1.37	1.33
2	D	401	NAD	C5A-C4A	2.20	1.46	1.40
2	B	401	NAD	C4A-N3A	2.14	1.38	1.35
2	B	401	NAD	C5A-C4A	2.11	1.46	1.40
2	C	500	NAD	C2A-N3A	2.08	1.35	1.32
2	D	401	NAD	C4A-N3A	-2.06	1.32	1.35

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	402	INS	C3-C2-C1	6.25	121.74	110.82
3	A	402	INS	C3-C2-C1	5.97	121.24	110.82
3	B	402	INS	C6-C5-C4	-5.19	101.76	110.82
3	D	402	INS	C3-C2-C1	4.71	119.04	110.82
3	A	402	INS	O6-C6-C5	-4.62	99.68	110.35
3	D	402	INS	C6-C5-C4	-4.43	103.10	110.82
2	C	500	NAD	PN-O3-PA	-4.27	118.19	132.83
2	B	401	NAD	C3B-C2B-C1B	-4.22	94.63	100.98
2	B	401	NAD	N3A-C2A-N1A	-4.10	122.27	128.68
3	B	402	INS	O6-C6-C5	-4.01	101.08	110.35
3	D	402	INS	O6-C6-C5	-3.95	101.21	110.35
2	A	401	NAD	N3A-C2A-N1A	-3.86	122.65	128.68
2	A	401	NAD	PN-O3-PA	-3.85	119.63	132.83
3	D	402	INS	C6-C1-C2	3.80	117.46	110.82
2	C	500	NAD	N3A-C2A-N1A	-3.73	122.84	128.68
2	B	401	NAD	C1B-N9A-C4A	-3.64	120.25	126.64
2	D	401	NAD	PN-O3-PA	-3.18	121.90	132.83
2	D	401	NAD	C6N-C5N-C4N	-3.15	114.86	119.44
3	A	402	INS	C6-C5-C4	-3.14	105.35	110.82
2	B	401	NAD	PN-O3-PA	-3.01	122.51	132.83
2	C	500	NAD	C1B-N9A-C4A	-2.90	121.55	126.64
3	B	402	INS	O2-C2-C1	-2.84	103.77	110.35
3	A	402	INS	O2-C2-C3	-2.80	103.89	110.35
3	B	402	INS	O5-C5-C4	2.74	116.68	110.35
3	A	402	INS	O3-C3-C4	-2.74	104.03	110.35
2	D	401	NAD	N3A-C2A-N1A	-2.68	124.48	128.68
2	D	401	NAD	N6A-C6A-N1A	2.66	124.10	118.57
2	A	401	NAD	C4A-C5A-N7A	-2.65	106.64	109.40
2	C	500	NAD	C3B-C2B-C1B	-2.61	97.05	100.98
2	A	401	NAD	O2A-PA-O1A	2.60	125.07	112.24
2	C	500	NAD	C3N-C2N-N1N	2.58	122.95	120.43
2	C	500	NAD	C3N-C7N-N7N	2.54	120.80	117.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	401	NAD	C3N-C2N-N1N	2.49	122.86	120.43
3	D	402	INS	O5-C5-C4	2.47	116.06	110.35
2	A	401	NAD	O4B-C1B-C2B	-2.39	103.43	106.93
2	D	401	NAD	C2N-N1N-C1D	-2.36	113.88	119.14
2	B	401	NAD	N6A-C6A-N1A	2.30	123.34	118.57
3	B	402	INS	O2-C2-C3	-2.30	105.04	110.35
2	C	500	NAD	O4B-C1B-C2B	-2.28	103.59	106.93
2	C	500	NAD	O3D-C3D-C4D	-2.24	104.59	111.05
2	A	401	NAD	C1B-N9A-C4A	-2.23	122.73	126.64
2	D	401	NAD	O4B-C1B-C2B	-2.20	103.71	106.93
3	B	402	INS	O3-C3-C4	-2.19	105.29	110.35
2	B	401	NAD	O2A-PA-O1A	2.16	122.90	112.24
2	B	401	NAD	C6N-N1N-C2N	-2.09	120.07	121.97
3	A	402	INS	C5-C6-C1	-2.08	107.19	110.82
2	D	401	NAD	C3B-C2B-C1B	-2.04	97.90	100.98
3	B	402	INS	C6-C1-C2	2.02	114.35	110.82
2	C	500	NAD	C5N-C6N-N1N	-2.01	117.53	120.40

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAD	O4D-C1D-N1N-C2N
2	A	401	NAD	O4D-C1D-N1N-C6N
2	A	401	NAD	C2D-C1D-N1N-C2N
2	B	401	NAD	O4D-C1D-N1N-C2N
2	B	401	NAD	O4D-C1D-N1N-C6N
2	B	401	NAD	C2D-C1D-N1N-C2N
2	C	500	NAD	O4D-C1D-N1N-C2N
2	C	500	NAD	O4D-C1D-N1N-C6N
2	C	500	NAD	C2D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	C2D-C1D-N1N-C2N
2	D	401	NAD	C2D-C1D-N1N-C6N
2	B	401	NAD	PN-O3-PA-O1A
2	C	500	NAD	O4B-C4B-C5B-O5B
2	A	401	NAD	O4B-C4B-C5B-O5B
2	B	401	NAD	O4B-C4B-C5B-O5B
2	D	401	NAD	O4B-C4B-C5B-O5B
2	B	401	NAD	PN-O3-PA-O2A

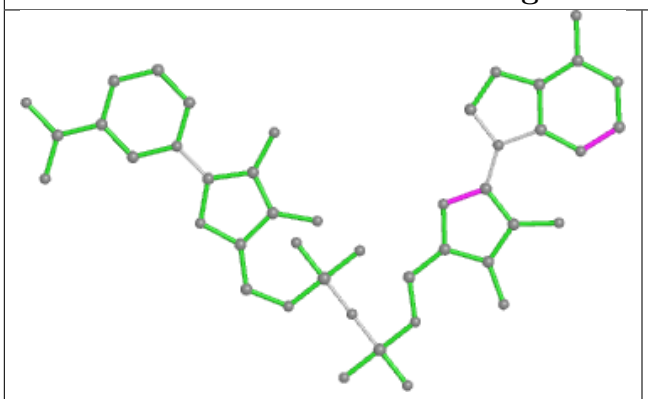
There are no ring outliers.

7 monomers are involved in 33 short contacts:

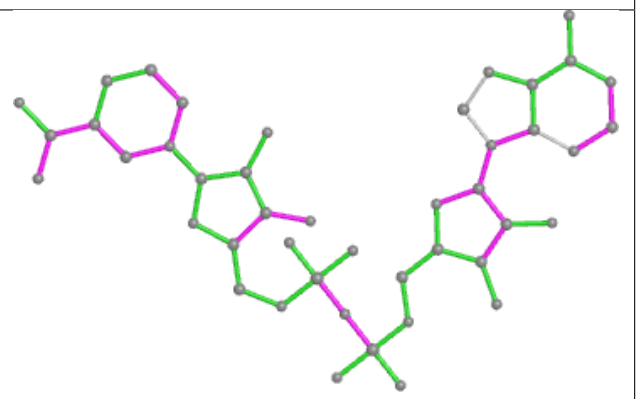
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	500	NAD	3	0
2	B	401	NAD	6	0
2	D	401	NAD	12	0
2	A	401	NAD	9	0
3	B	402	INS	5	0
3	A	402	INS	4	0
3	D	402	INS	5	0

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

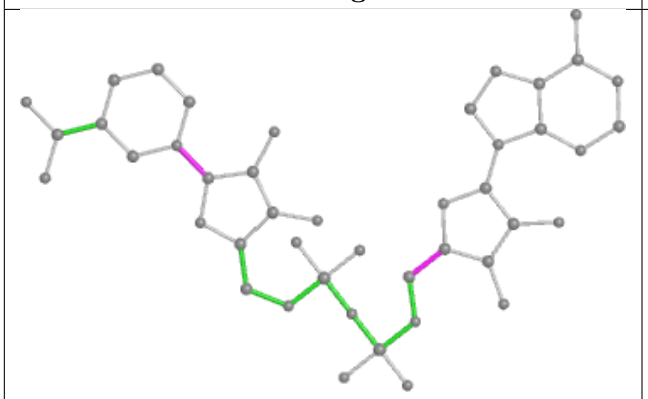
Ligand NAD C 500



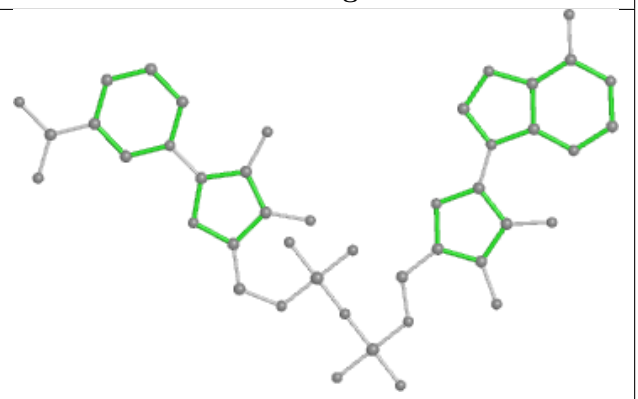
Bond lengths



Bond angles

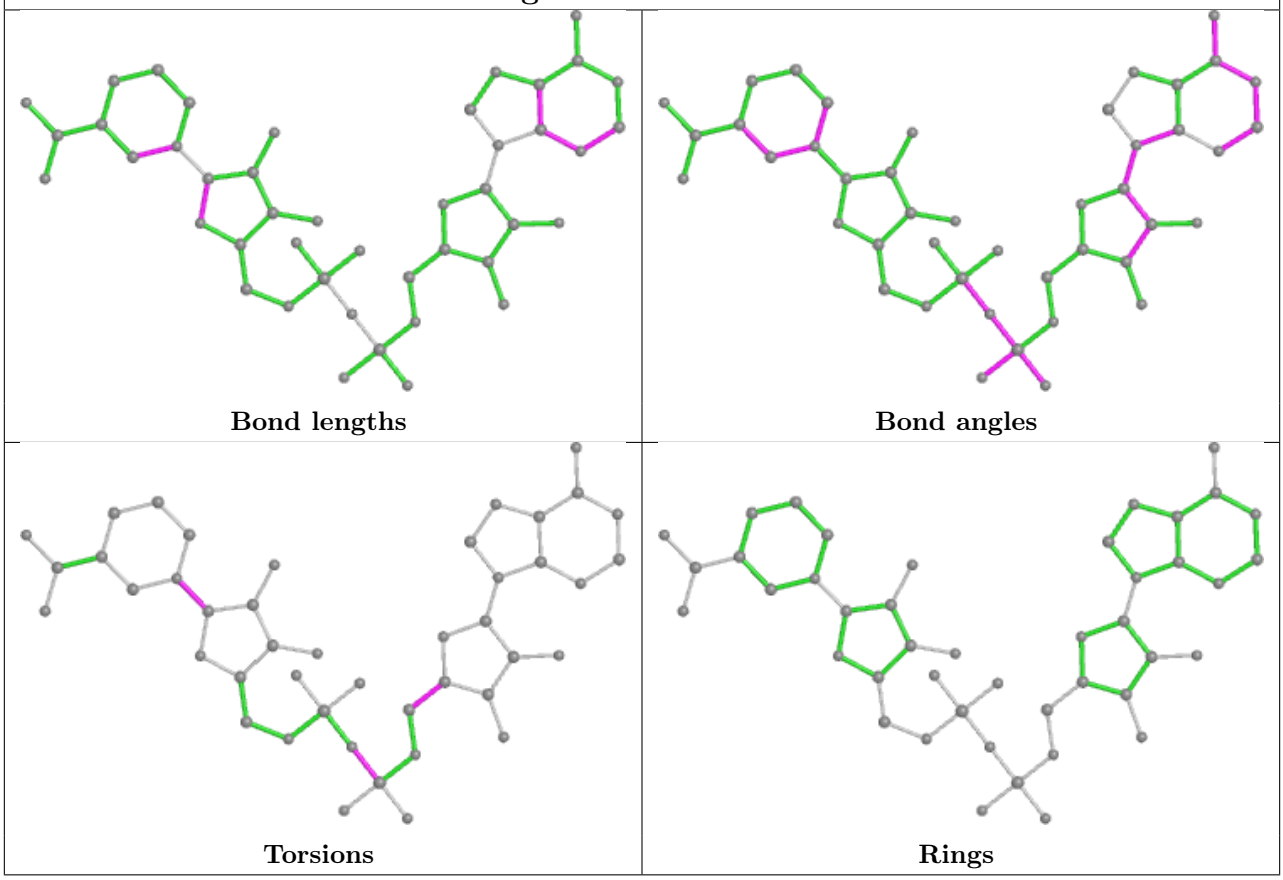


Torsions

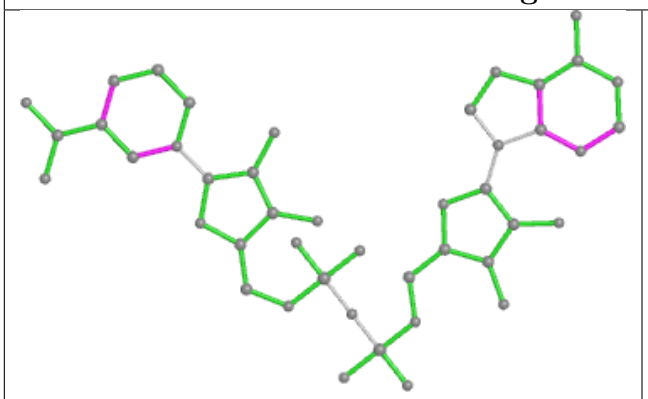


Rings

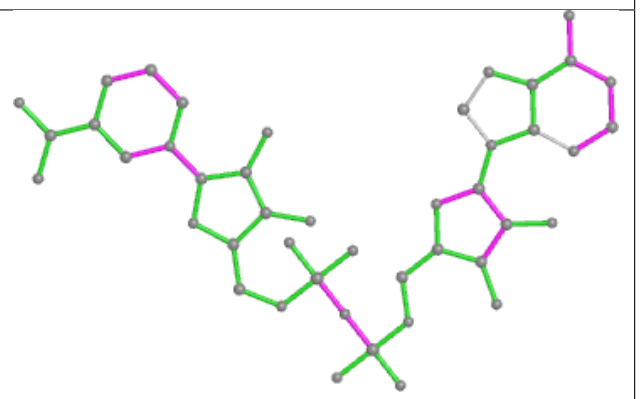
Ligand NAD B 401



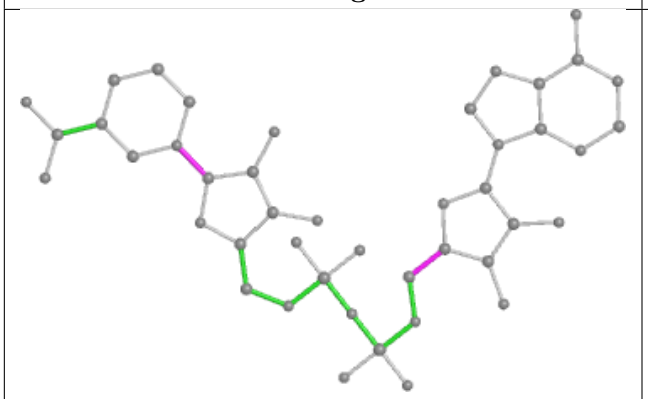
Ligand NAD D 401



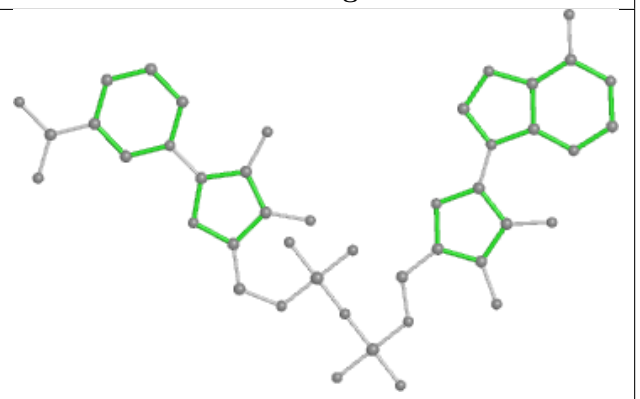
Bond lengths



Bond angles

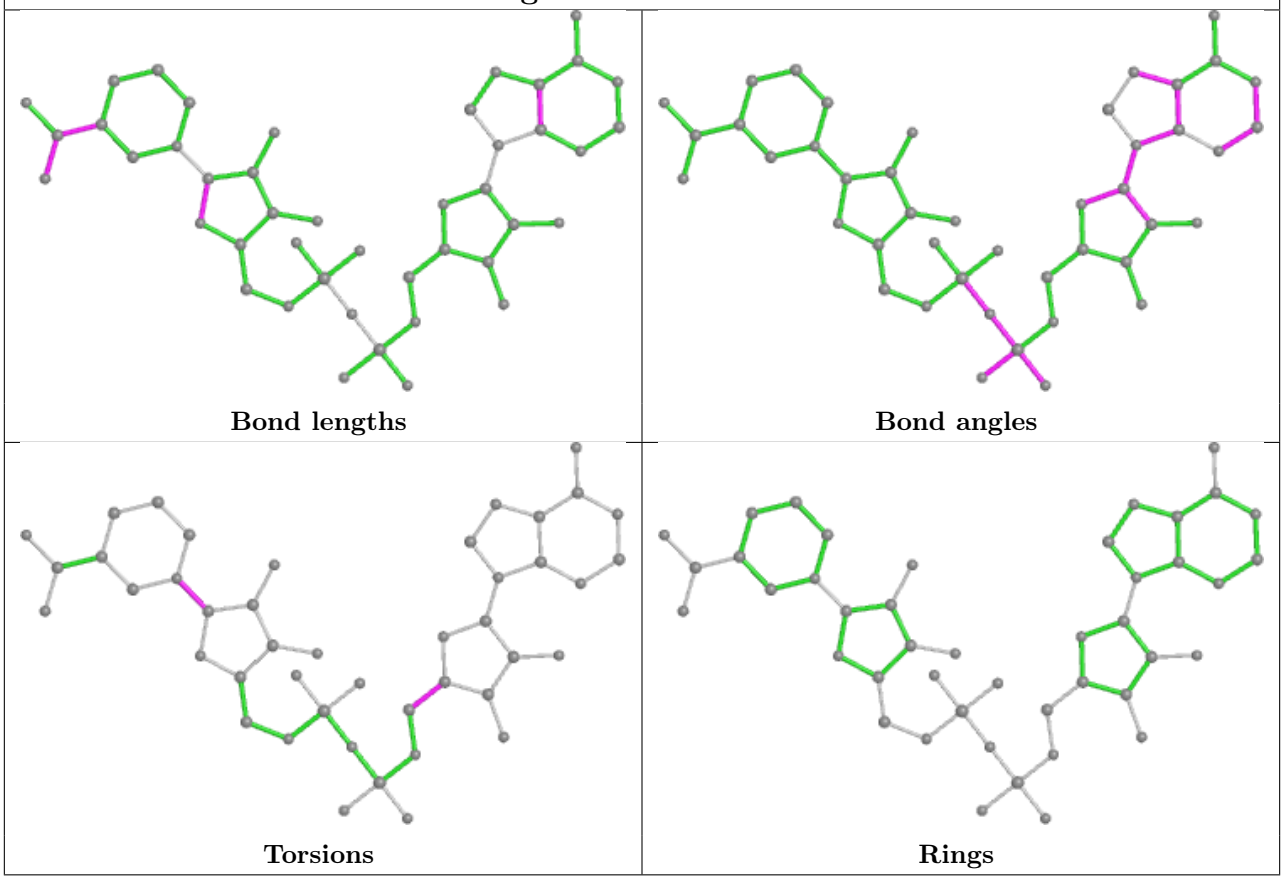


Torsions



Rings

Ligand NAD A 401



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	366/380 (96%)	-0.15	12 (3%)	46	53	16, 25, 46, 87	0
1	B	366/380 (96%)	-0.14	13 (3%)	42	49	14, 24, 45, 66	4 (1%)
1	C	366/380 (96%)	0.15	25 (6%)	17	22	15, 26, 58, 103	2 (0%)
1	D	365/380 (96%)	-0.27	9 (2%)	57	64	13, 24, 41, 71	1 (0%)
All	All	1463/1520 (96%)	-0.10	59 (4%)	38	45	13, 25, 48, 103	7 (0%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	224	GLN	9.0
1	C	183	ASP	6.4
1	C	223	HIS	5.4
1	C	177	TRP	5.1
1	C	175	TRP	4.9
1	C	227	THR	4.5
1	C	229	ARG	4.3
1	D	174	PRO	4.3
1	C	225	GLY	4.3
1	A	176	SER	4.3
1	A	175	TRP	3.9
1	C	228	ALA	3.9
1	C	174	PRO	3.8
1	C	222	PRO	3.4
1	D	340	ARG	3.2
1	C	180	THR	3.1
1	C	125	SER	3.1
1	B	75	ALA	3.1
1	B	70	ALA	3.1
1	D	172	ASP	3.1
1	C	226	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	225	GLY	2.9
1	A	224	GLN	2.8
1	D	225	GLY	2.8
1	A	181	ARG	2.8
1	C	46	MET	2.8
1	A	340	ARG	2.7
1	A	174	PRO	2.7
1	B	340	ARG	2.6
1	C	176	SER	2.6
1	C	194	CYS	2.6
1	A	183	ASP	2.5
1	D	175	TRP	2.5
1	A	223	HIS	2.5
1	B	69	GLU	2.5
1	B	168	MET	2.5
1	B	342	LYS	2.5
1	C	72	SER	2.5
1	C	48	ALA	2.4
1	B	52	HIS	2.4
1	D	171	PRO	2.4
1	B	8	LEU	2.3
1	C	182	LYS	2.3
1	B	194	CYS	2.3
1	B	66	ASP	2.3
1	B	49	ASP	2.3
1	D	227	THR	2.3
1	A	342	LYS	2.3
1	B	181	ARG	2.2
1	A	180	THR	2.1
1	A	341	GLY	2.1
1	D	49	ASP	2.1
1	C	63	GLY	2.1
1	C	340	ARG	2.1
1	D	342	LYS	2.1
1	C	64	THR	2.1
1	B	68	ARG	2.0
1	C	67	TRP	2.0
1	C	173	LEU	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

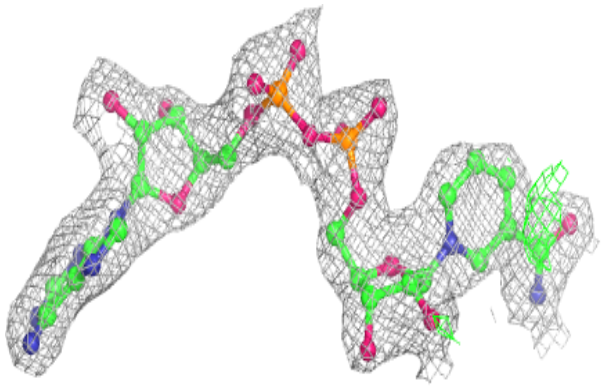
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

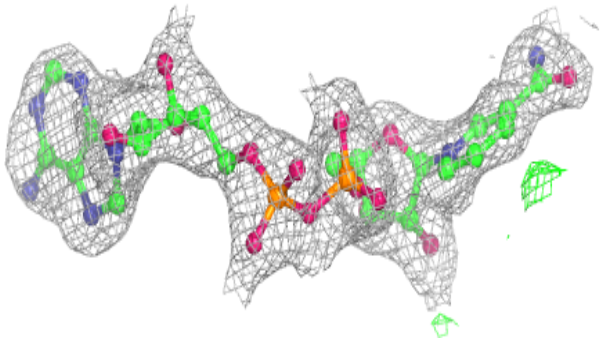
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	INS	B	402	12/12	0.89	0.14	23,23,23,26	0
3	INS	A	402	12/12	0.91	0.21	23,24,26,26	0
3	INS	D	402	12/12	0.91	0.16	22,23,24,24	0
2	NAD	C	500	44/44	0.94	0.17	26,30,36,38	0
2	NAD	B	401	44/44	0.94	0.14	23,28,32,35	0
2	NAD	D	401	44/44	0.95	0.15	18,25,33,35	0
2	NAD	A	401	44/44	0.95	0.14	23,27,32,34	0

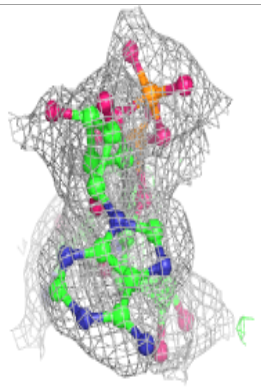
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD C 500:

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

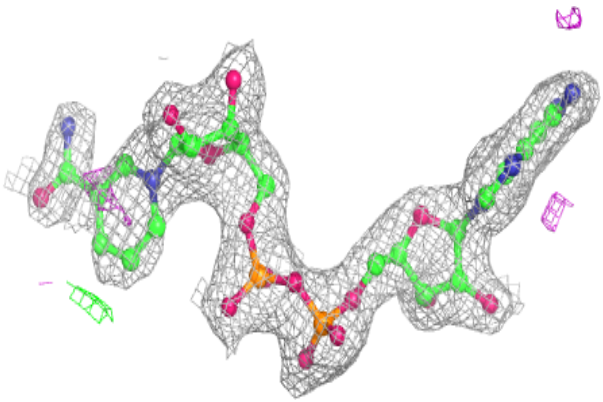


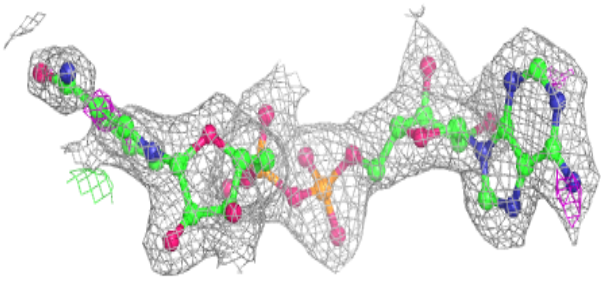


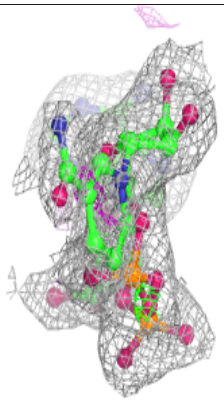


Electron density around NAD B 401:

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

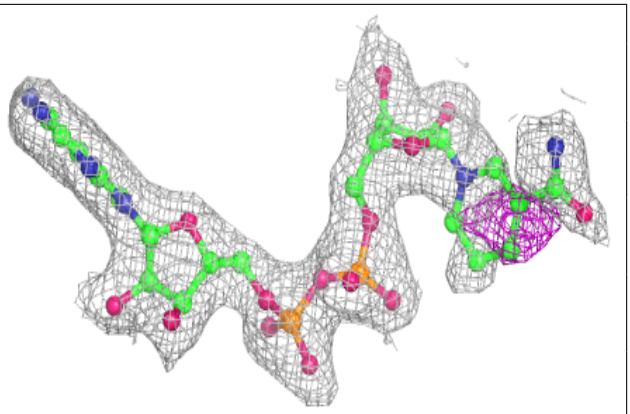


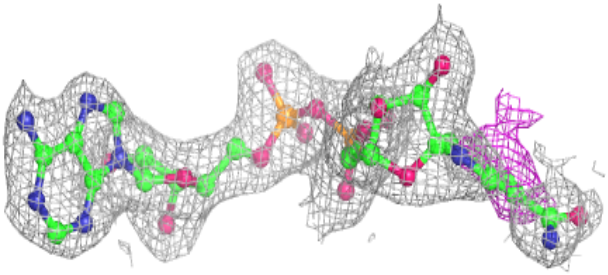


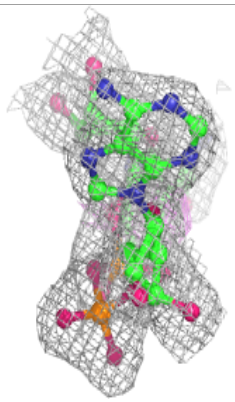


Electron density around NAD D 401:

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

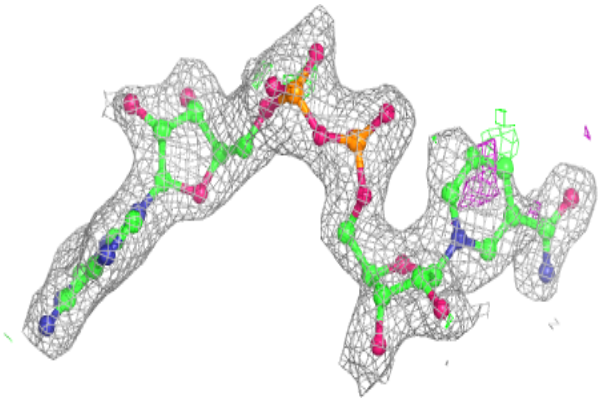


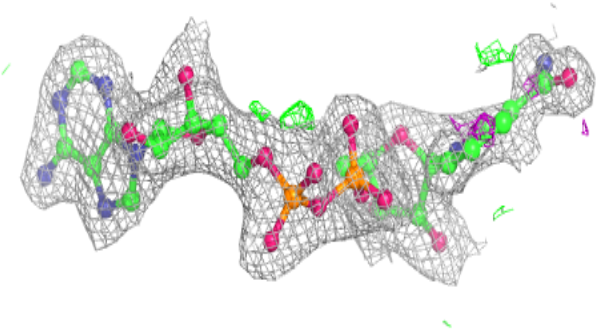


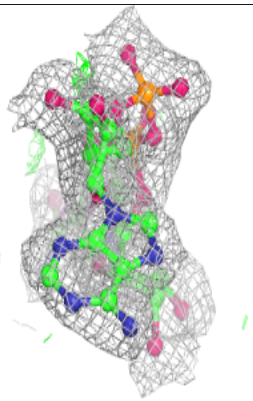


Electron density around NAD A 401:

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







6.5 Other polymers (i)

There are no such residues in this entry.