Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2023 – 01:23 AM EDT

PDB ID	:	3LC1
Title	:	Crystal Structure of H178N mutant of Glyceraldehyde-3-phosphate-dehydrogenase 1 (GAPDH 1) from Staphylococcus aureus MRSA252 complexed with NAD at 2.0 angstrom resolution.
Authors	:	Mukherjee, S.; Dutta, D.; Saha, B.; Das, A.K.
Deposited on	:	2010-01-09
Resolution	:	2.00 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:

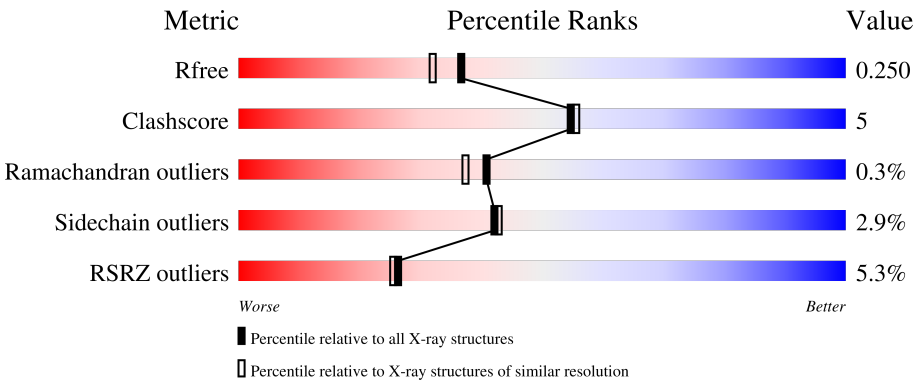
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	O	336	4% 85% 14% ••
1	P	336	4% 91% 8% •
1	Q	336	8% 87% 12% ••
1	R	336	4% 93% 5% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10756 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 1.

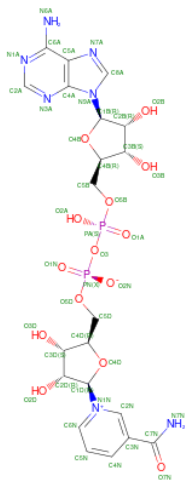
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	P	336	2555	1590	440	515	10	0	2	0
1	R	336	2552	1587	438	517	10	0	1	0
1	O	334	2527	1572	434	512	9	0	0	0
1	Q	334	2527	1572	434	512	9	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	178	ASN	HIS	engineered mutation	UNP Q6GIL8
R	178	ASN	HIS	engineered mutation	UNP Q6GIL8
O	178	ASN	HIS	engineered mutation	UNP Q6GIL8
Q	178	ASN	HIS	engineered mutation	UNP Q6GIL8

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

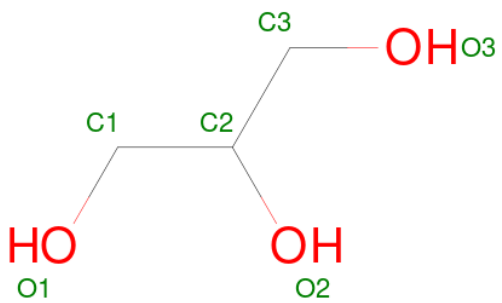
NAD



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

GOL



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	P	1	6	3	3	0	0

- Molecule 4 is water.

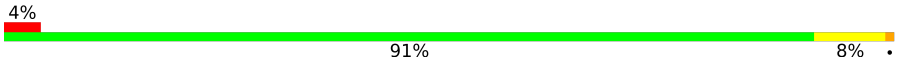
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	P	100	100	100	0	0
			Total	O		
4	R	111	111	111	0	0
			Total	O		
4	O	118	118	118	0	0
			Total	O		
4	Q	84	84	84	0	0

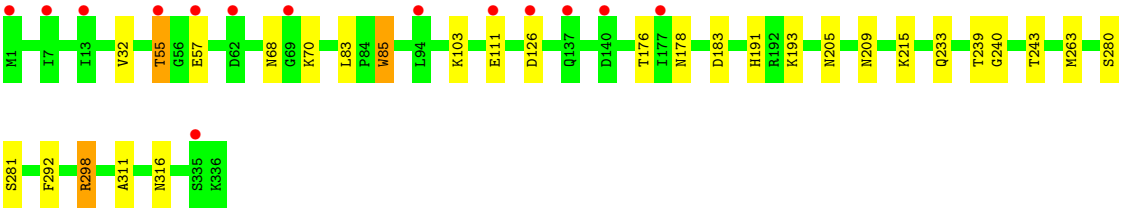
3 Residue-property plots (i)

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 1

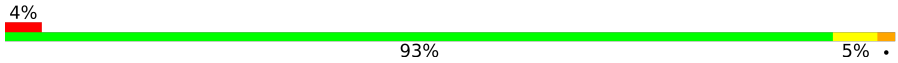
Chain P:

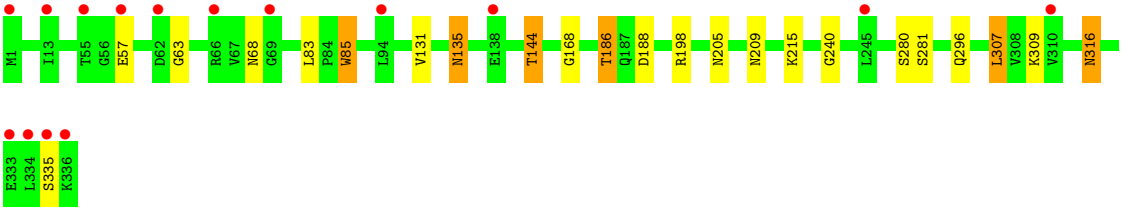




• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

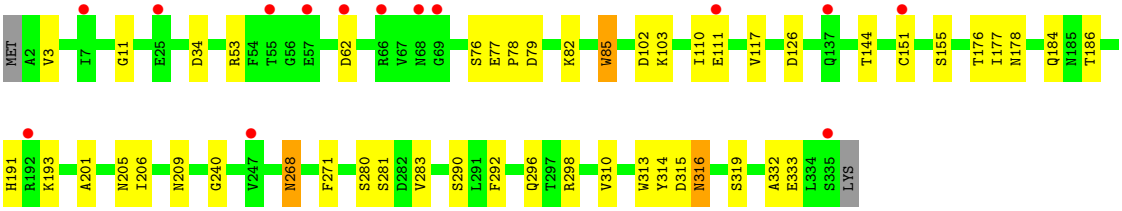
Chain R:





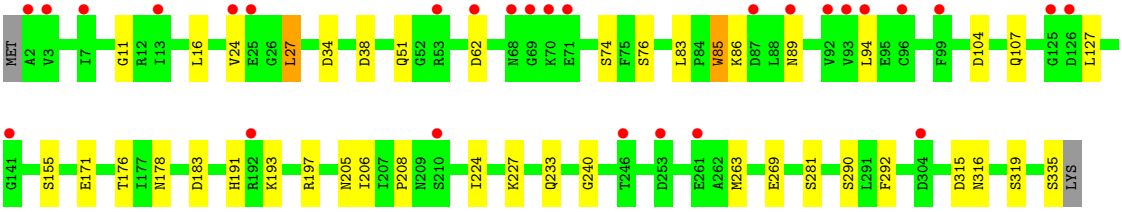
• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

Chain O:



• Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

Chain Q:


MET A2 V3 I7 G11 R12 I13 L16 V24 E25 G26 L27 D34 D38 Q51 G52 R53 D62 N68 G69 K70 E71 S74 F75 S76 L83 P84 W85 K86 D87 L88 N89 V92 V93 L94 E95 C96 F99 D104 Q107 G125 D126 L127
G141 S155 E171 T176 I177 N178 D183 H191 R192 K193 R197 N205 I206 I207 P208 N209 S210 I224 K227 Q233 G240 T246 D253 E261 A262 M263 E269 S281 S290 L291 F292 D304 D315 N316 S319 S335 LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	68.85Å 105.94Å 91.47Å 90.00° 108.34° 90.00°	Depositor
Resolution (Å)	29.06 – 2.00 29.06 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.2 (29.06-2.00) 95.3 (29.06-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.56 (at 2.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.197 , 0.242 0.209 , 0.250	Depositor DCC
R_{free} test set	4003 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	30.8	Xtriage
Anisotropy	0.660	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 37.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	10756	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: GOL, 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	O	0.54	0/2560	0.64	0/3469
1	P	0.54	0/2594	0.63	0/3512
1	Q	0.52	0/2560	0.62	0/3469
1	R	0.56	0/2585	0.66	1/3501 (0.0%)
All	All	0.54	0/10299	0.64	1/13951 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	R	188	ASP	CB-CG-OD1	6.23	123.91	118.30

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	O	2527	0	2504	41	0
1	P	2555	0	2547	30	0
1	Q	2527	0	2504	26	0
1	R	2552	0	2532	25	0
2	O	44	0	26	1	0
2	P	44	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	Q	44	0	26	1	0
2	R	44	0	26	0	0
3	P	6	0	8	0	0
4	O	118	0	0	5	0
4	P	100	0	0	2	0
4	Q	84	0	0	2	0
4	R	111	0	0	6	0
All	All	10756	0	10199	106	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (106) 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:P:298[A]:ARG:HG3	1:P:298[A]:ARG:NH1	1.52	1.11
1:P:298[A]:ARG:HH11	1:P:298[A]:ARG:CG	1.64	1.07
1:R:144:THR:HG21	4:R:411:HOH:O	1.63	0.99
1:R:63:GLY:HA3	4:R:444:HOH:O	1.72	0.89
1:O:111:GLU:HG2	4:O:364:HOH:O	1.72	0.88
4:P:432:HOH:O	1:O:298:ARG:HD3	1.74	0.88
1:R:131:VAL:H	1:R:135:ASN:HD21	1.32	0.75
1:P:298[A]:ARG:HG3	1:P:298[A]:ARG:HH11	0.71	0.74
1:P:298[A]:ARG:NH1	1:P:298[A]:ARG:CG	2.35	0.71
1:P:183:ASP:O	1:P:191:HIS:HE1	1.75	0.69
1:Q:191:HIS:CD2	1:Q:193:LYS:H	2.13	0.66
1:R:131:VAL:H	1:R:135:ASN:ND2	1.96	0.64
1:Q:183:ASP:O	1:Q:191:HIS:HE1	1.81	0.64
1:Q:176:THR:HG23	1:Q:178:ASN:ND2	2.15	0.61
1:P:233:GLN:HG3	1:O:177:ILE:HD13	1.84	0.59
1:Q:38:ASP:OD1	1:Q:74:SER:OG	2.18	0.58
1:Q:224:ILE:HA	1:Q:227:LYS:HD2	1.84	0.58
1:R:198:ARG:NH1	4:R:390:HOH:O	2.38	0.57
1:Q:83:LEU:HD13	1:Q:85:TRP:CZ2	2.40	0.56
1:P:176:THR:HG23	1:P:178:ASN:ND2	2.20	0.56
1:P:57:GLU:H	1:P:68:ASN:ND2	2.03	0.55
1:P:191:HIS:CD2	1:P:193:LYS:H	2.24	0.55
1:R:186:THR:HG21	1:O:184:GLN:O	2.06	0.55
1:Q:240:GLY:H	1:Q:316:ASN:HD21	1.54	0.54
1:Q:24:VAL:HG11	1:Q:27:LEU:HD22	1.89	0.53
1:P:263:MET:HG3	1:P:292:PHE:CZ	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:268:ASN:ND2	1:O:271:PHE:H	2.06	0.53
1:R:186:THR:HG23	4:R:438:HOH:O	2.08	0.53
1:O:77:GLU:HG2	1:O:82:LYS:HB2	1.91	0.52
1:P:205:ASN:HD22	1:O:281:SER:H	1.58	0.51
1:P:240:GLY:H	1:P:316:ASN:HD21	1.58	0.51
1:O:103:LYS:HD2	1:O:126:ASP:O	2.11	0.51
1:P:83:LEU:HD13	1:P:85:TRP:CZ2	2.47	0.50
1:Q:240:GLY:H	1:Q:316:ASN:ND2	2.09	0.50
1:Q:104:ASP:O	1:Q:107:GLN:HB3	2.13	0.49
1:R:309:LYS:HE2	4:R:412:HOH:O	2.12	0.49
1:Q:269:GLU:HG2	4:Q:414:HOH:O	2.13	0.49
1:P:243:THR:O	1:P:311:ALA:HA	2.13	0.49
1:O:155:SER:HA	1:O:290:SER:HB2	1.94	0.49
1:O:240:GLY:H	1:O:316:ASN:ND2	2.11	0.48
1:Q:197:ARG:HD3	1:Q:208:PRO:O	2.13	0.48
1:P:263:MET:HG3	1:P:292:PHE:CE1	2.48	0.48
1:P:176:THR:HG23	1:P:178:ASN:HD21	1.79	0.48
1:O:268:ASN:HD22	1:O:268:ASN:C	2.17	0.48
1:O:176:THR:HG23	1:O:178:ASN:ND2	2.29	0.47
1:R:83:LEU:HD13	1:R:85:TRP:CZ2	2.50	0.47
1:P:281:SER:CB	1:O:205:ASN:HD21	2.28	0.47
1:R:281:SER:H	1:Q:205:ASN:HD22	1.63	0.47
1:O:3:VAL:HG21	1:O:332:ALA:HB1	1.97	0.47
1:O:240:GLY:H	1:O:316:ASN:HD21	1.63	0.46
1:O:292:PHE:HE1	1:O:310:VAL:HB	1.80	0.46
1:P:55:THR:HG22	4:P:371:HOH:O	2.15	0.46
1:O:144:THR:HG21	4:O:393:HOH:O	2.15	0.46
1:O:144:THR:HG22	4:O:433:HOH:O	2.14	0.46
1:Q:176:THR:CG2	1:Q:178:ASN:HD21	2.28	0.46
1:O:283:VAL:HG11	1:O:313:TRP:HB3	1.96	0.46
1:P:233:GLN:HG3	1:O:177:ILE:CD1	2.45	0.45
1:P:240:GLY:H	1:P:316:ASN:ND2	2.13	0.45
1:O:34:ASP:O	1:O:76:SER:HA	2.16	0.45
1:P:233:GLN:HE22	1:O:296:GLN:HE22	1.64	0.45
1:R:240:GLY:H	1:R:316:ASN:ND2	2.14	0.45
1:R:135:ASN:HD22	1:R:135:ASN:H	1.65	0.45
1:P:316:ASN:ND2	1:P:316:ASN:H	2.16	0.44
1:R:240:GLY:H	1:R:316:ASN:HD21	1.63	0.44
1:P:103:LYS:HD2	1:P:126:ASP:O	2.16	0.44
1:Q:176:THR:HG23	1:Q:178:ASN:HD21	1.82	0.44
1:P:233:GLN:HE22	1:O:296:GLN:NE2	2.14	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:215:LYS:HB3	1:R:215:LYS:HE2	1.66	0.44
1:R:280:SER:HB3	1:Q:206:ILE:HB	2.00	0.44
1:O:268:ASN:HD22	1:O:271:PHE:H	1.65	0.44
1:O:280:SER:O	1:O:283:VAL:HG22	2.18	0.44
1:Q:263:MET:HG3	1:Q:292:PHE:CZ	2.53	0.44
1:R:205:ASN:HD22	1:Q:281:SER:H	1.65	0.43
1:Q:155:SER:HA	1:Q:290:SER:HB2	2.00	0.43
1:R:85:TRP:CE3	1:R:85:TRP:HA	2.54	0.43
1:R:135:ASN:HD22	1:R:135:ASN:N	2.17	0.43
1:O:316:ASN:ND2	1:O:316:ASN:H	2.17	0.43
1:O:315:ASP:O	1:O:319:SER:HB2	2.18	0.43
1:R:57:GLU:H	1:R:68:ASN:ND2	2.16	0.43
1:R:316:ASN:ND2	1:R:316:ASN:H	2.17	0.43
1:O:191:HIS:CE1	1:O:193:LYS:HB2	2.54	0.43
1:R:85:TRP:HA	1:R:85:TRP:HE3	1.84	0.42
1:O:79:ASP:OD2	1:O:82:LYS:HE2	2.19	0.42
1:Q:191:HIS:HD2	1:Q:193:LYS:H	1.64	0.42
1:P:281:SER:H	1:O:205:ASN:ND2	2.17	0.42
1:R:68:ASN:HB2	4:R:370:HOH:O	2.19	0.42
1:O:11:GLY:HA3	2:O:0:NAD:O5B	2.20	0.42
1:O:110:ILE:HD11	1:O:117:VAL:HG23	2.02	0.42
1:Q:315:ASP:O	1:Q:319:SER:HB2	2.19	0.42
1:P:205:ASN:ND2	1:O:281:SER:H	2.18	0.41
1:O:53:ARG:HD3	4:Q:420:HOH:O	2.20	0.41
1:Q:16:LEU:HD21	1:Q:51:GLN:HG3	2.01	0.41
1:O:85:TRP:HA	1:O:85:TRP:CE3	2.55	0.41
1:O:193:LYS:HE2	4:O:389:HOH:O	2.20	0.41
1:Q:34:ASP:O	1:Q:76:SER:HA	2.20	0.41
1:P:281:SER:H	1:O:205:ASN:HD22	1.67	0.41
1:O:85:TRP:HA	1:O:85:TRP:HE3	1.84	0.41
1:P:85:TRP:HA	1:P:85:TRP:CE3	2.56	0.41
1:R:131:VAL:N	1:R:135:ASN:HD21	2.10	0.41
1:Q:11:GLY:HA3	2:Q:0:NAD:O5B	2.20	0.41
1:P:85:TRP:HA	1:P:85:TRP:HE3	1.85	0.41
1:P:280:SER:HB3	1:O:206:ILE:HB	2.02	0.41
1:O:151:CYS:SG	4:O:450:HOH:O	2.62	0.40
1:R:296:GLN:NE2	1:Q:233:GLN:HE22	2.20	0.40
1:R:307:LEU:HD22	1:Q:171:GLU:HG3	2.04	0.40
1:O:77:GLU:HA	1:O:78:PRO:HD2	1.98	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	O	332/336 (99%)	318 (96%)	13 (4%)	1 (0%)	41	37
1	P	336/336 (100%)	321 (96%)	15 (4%)	0	100	100
1	Q	332/336 (99%)	317 (96%)	14 (4%)	1 (0%)	41	37
1	R	335/336 (100%)	321 (96%)	12 (4%)	2 (1%)	25	19
All	All	1335/1344 (99%)	1277 (96%)	54 (4%)	4 (0%)	41	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	Q	62	ASP
1	R	335	SER
1	O	201	ALA
1	R	168	GLY

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	O	272/274 (99%)	263 (97%)	9 (3%)	38	37
1	P	276/274 (101%)	266 (96%)	10 (4%)	35	34
1	Q	272/274 (99%)	265 (97%)	7 (3%)	46	48
1	R	275/274 (100%)	268 (98%)	7 (2%)	47	49
All	All	1095/1096 (100%)	1062 (97%)	33 (3%)	42	41

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

Mol	Chain	Res	Type
1	P	32	VAL
1	P	55	THR
1	P	70	LYS
1	P	85	TRP
1	P	111	GLU
1	P	209	ASN
1	P	215	LYS
1	P	239	THR
1	P	298[A]	ARG
1	P	298[B]	ARG
1	R	85	TRP
1	R	135	ASN
1	R	144	THR
1	R	186	THR
1	R	209	ASN
1	R	307	LEU
1	R	316	ASN
1	O	62	ASP
1	O	85	TRP
1	O	102	ASP
1	O	186	THR
1	O	209	ASN
1	O	268	ASN
1	O	314	TYR
1	O	316	ASN
1	O	333	GLU
1	Q	27	LEU
1	Q	85	TRP
1	Q	86	LYS
1	Q	89	ASN
1	Q	94	LEU
1	Q	127	LEU
1	Q	335	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	P	68	ASN
1	P	178	ASN
1	P	191	HIS
1	P	205	ASN

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Mol	Chain	Res	Type
1	P	252	GLN
1	P	316	ASN
1	R	68	ASN
1	R	135	ASN
1	R	178	ASN
1	R	205	ASN
1	R	265	ASN
1	R	296	GLN
1	R	316	ASN
1	O	178	ASN
1	O	205	ASN
1	O	265	ASN
1	O	268	ASN
1	O	296	GLN
1	O	316	ASN
1	Q	178	ASN
1	Q	191	HIS
1	Q	205	ASN
1	Q	265	ASN
1	Q	296	GLN
1	Q	316	ASN

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)

5 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	R	0	-	42,48,48	1.70	4 (9%)	50,73,73	1.19	4 (8%)
2	NAD	O	0	-	42,48,48	1.66	4 (9%)	50,73,73	1.33	4 (8%)
2	NAD	P	0	-	42,48,48	1.68	4 (9%)	50,73,73	1.35	5 (10%)
2	NAD	Q	0	-	42,48,48	1.65	4 (9%)	50,73,73	1.28	5 (10%)
3	GOL	P	337	-	5,5,5	0.29	0	5,5,5	0.34	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	R	0	-	-	6/26/62/62	0/5/5/5
2	NAD	O	0	-	-	6/26/62/62	0/5/5/5
2	NAD	P	0	-	-	6/26/62/62	0/5/5/5
2	NAD	Q	0	-	-	5/26/62/62	0/5/5/5
3	GOL	P	337	-	-	2/4/4/4	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	0	NAD	O7N-C7N	8.30	1.40	1.24
2	O	0	NAD	O7N-C7N	8.20	1.39	1.24
2	R	0	NAD	O7N-C7N	8.06	1.39	1.24
2	P	0	NAD	O7N-C7N	8.00	1.39	1.24
2	P	0	NAD	C2A-N3A	4.04	1.38	1.32
2	Q	0	NAD	C2A-N3A	3.92	1.38	1.32
2	O	0	NAD	C2A-N3A	3.87	1.38	1.32
2	R	0	NAD	C2A-N3A	3.54	1.37	1.32
2	R	0	NAD	C2N-N1N	3.31	1.39	1.35
2	P	0	NAD	C2N-N1N	3.23	1.38	1.35
2	O	0	NAD	C2N-N1N	2.63	1.38	1.35
2	R	0	NAD	C2A-N1A	2.55	1.38	1.33
2	Q	0	NAD	C2N-N1N	2.53	1.38	1.35
2	O	0	NAD	C2A-N1A	2.52	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	0	NAD	C2A-N1A	2.41	1.38	1.33
2	Q	0	NAD	C2A-N1A	2.27	1.38	1.33

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	P	0	NAD	N3A-C2A-N1A	-5.81	119.60	128.68
2	O	0	NAD	N3A-C2A-N1A	-5.53	120.04	128.68
2	R	0	NAD	N3A-C2A-N1A	-5.48	120.11	128.68
2	Q	0	NAD	N3A-C2A-N1A	-5.07	120.76	128.68
2	P	0	NAD	C3N-C7N-N7N	3.99	122.53	117.75
2	O	0	NAD	C3N-C7N-N7N	3.72	122.22	117.75
2	Q	0	NAD	PN-O3-PA	-2.64	123.78	132.83
2	Q	0	NAD	C3N-C7N-N7N	2.61	120.88	117.75
2	O	0	NAD	O7N-C7N-C3N	-2.60	116.52	119.63
2	O	0	NAD	PN-O3-PA	-2.38	124.67	132.83
2	Q	0	NAD	O7N-C7N-C3N	-2.27	116.92	119.63
2	P	0	NAD	O7N-C7N-C3N	-2.22	116.98	119.63
2	R	0	NAD	C3N-C7N-N7N	2.19	120.38	117.75
2	R	0	NAD	C1B-N9A-C4A	-2.15	122.87	126.64
2	P	0	NAD	PN-O3-PA	-2.09	125.65	132.83
2	R	0	NAD	O7N-C7N-N7N	-2.08	119.62	122.58
2	Q	0	NAD	C4A-C5A-N7A	-2.07	107.24	109.40
2	P	0	NAD	C1B-N9A-C4A	-2.01	123.10	126.64

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	P	0	NAD	O4D-C1D-N1N-C2N
2	P	0	NAD	O4D-C1D-N1N-C6N
2	R	0	NAD	O4D-C1D-N1N-C2N
2	R	0	NAD	O4D-C1D-N1N-C6N
2	R	0	NAD	C2D-C1D-N1N-C2N
2	R	0	NAD	C2D-C1D-N1N-C6N
2	O	0	NAD	O4D-C1D-N1N-C2N
2	O	0	NAD	O4D-C1D-N1N-C6N
2	O	0	NAD	C2D-C1D-N1N-C2N
2	O	0	NAD	C2D-C1D-N1N-C6N
2	Q	0	NAD	O4D-C1D-N1N-C2N
2	Q	0	NAD	O4D-C1D-N1N-C6N
2	Q	0	NAD	C2D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
2	Q	0	NAD	C2D-C1D-N1N-C6N
3	P	337	GOL	C1-C2-C3-O3
3	P	337	GOL	O2-C2-C3-O3
2	O	0	NAD	O4B-C4B-C5B-O5B
2	P	0	NAD	O4B-C4B-C5B-O5B
2	R	0	NAD	O4B-C4B-C5B-O5B
2	O	0	NAD	C3B-C4B-C5B-O5B
2	Q	0	NAD	O4B-C4B-C5B-O5B
2	P	0	NAD	C2D-C1D-N1N-C2N
2	P	0	NAD	C2D-C1D-N1N-C6N
2	P	0	NAD	C3B-C4B-C5B-O5B
2	R	0	NAD	C3B-C4B-C5B-O5B

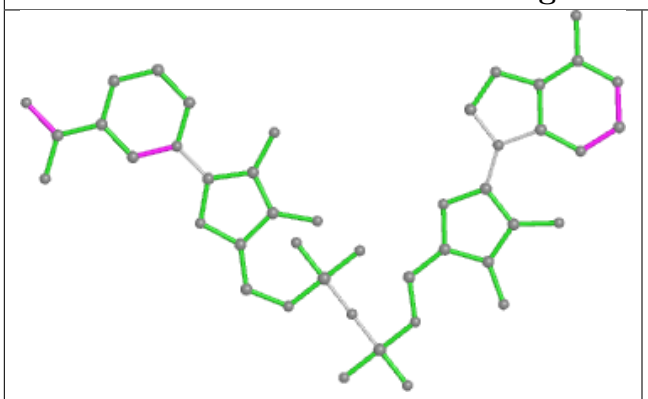
There are no ring outliers.

2 monomers are involved in 2 short contacts:

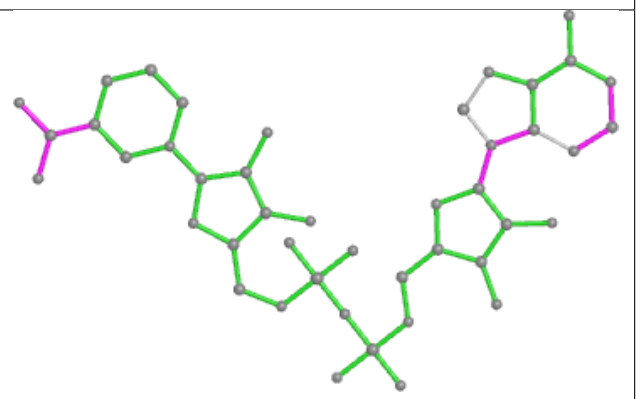
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	O	0	NAD	1	0
2	Q	0	NAD	1	0

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

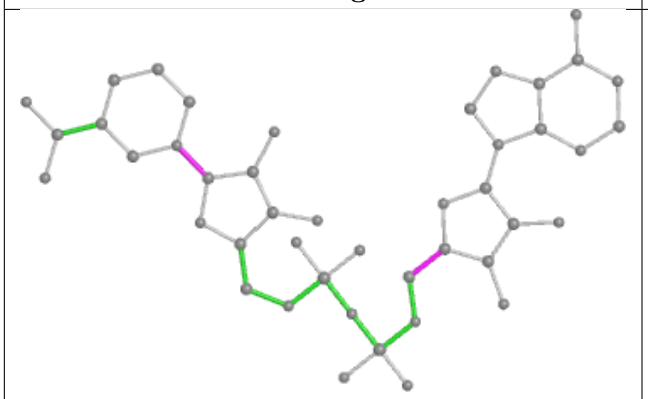
Ligand NAD R 0



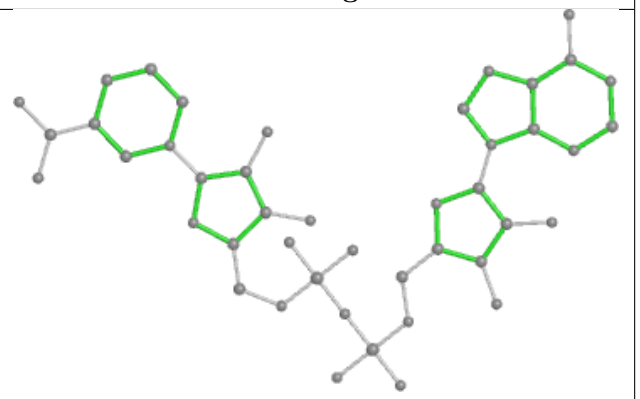
Bond lengths



Bond angles

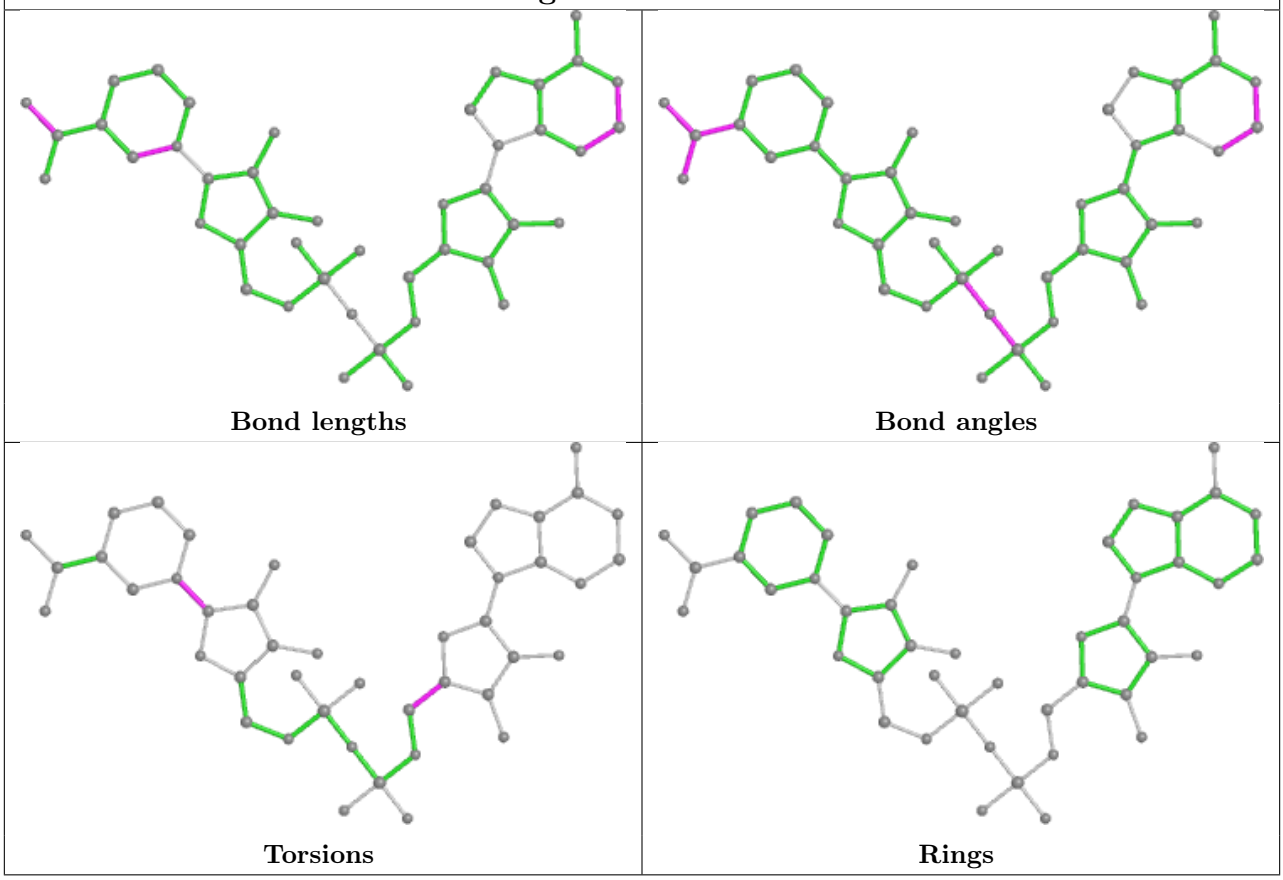


Torsions

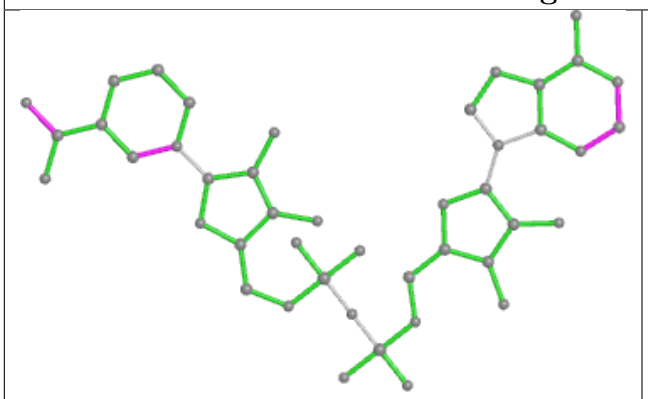


Rings

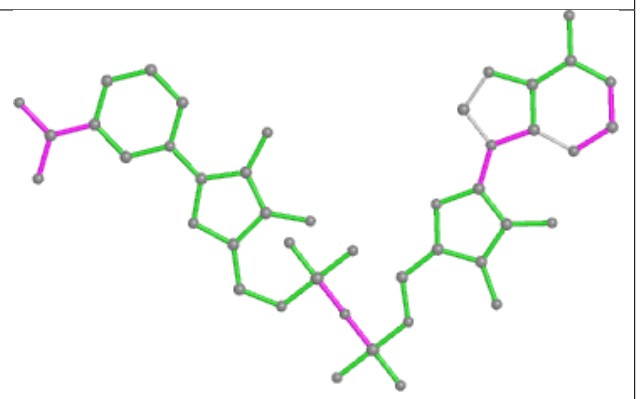
Ligand NAD O 0



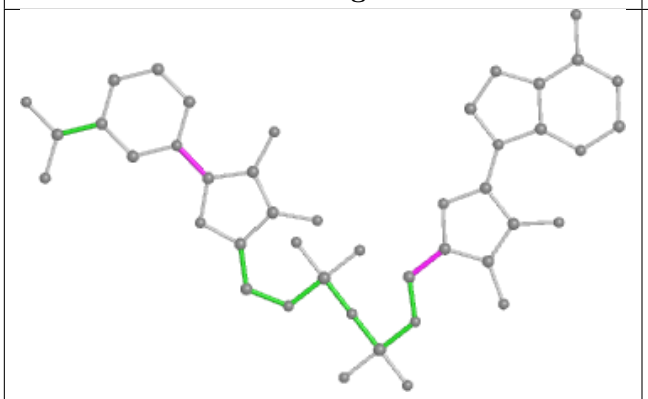
Ligand NAD P 0



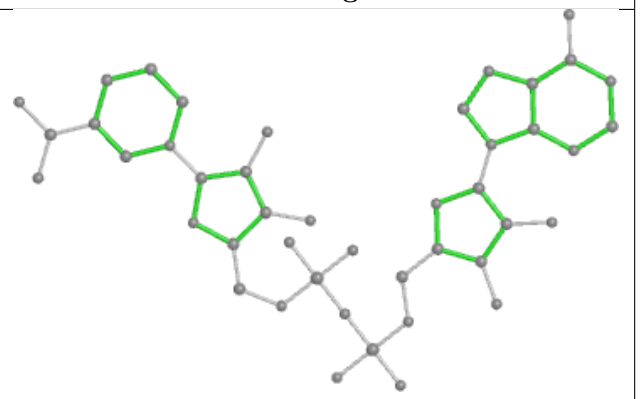
Bond lengths



Bond angles

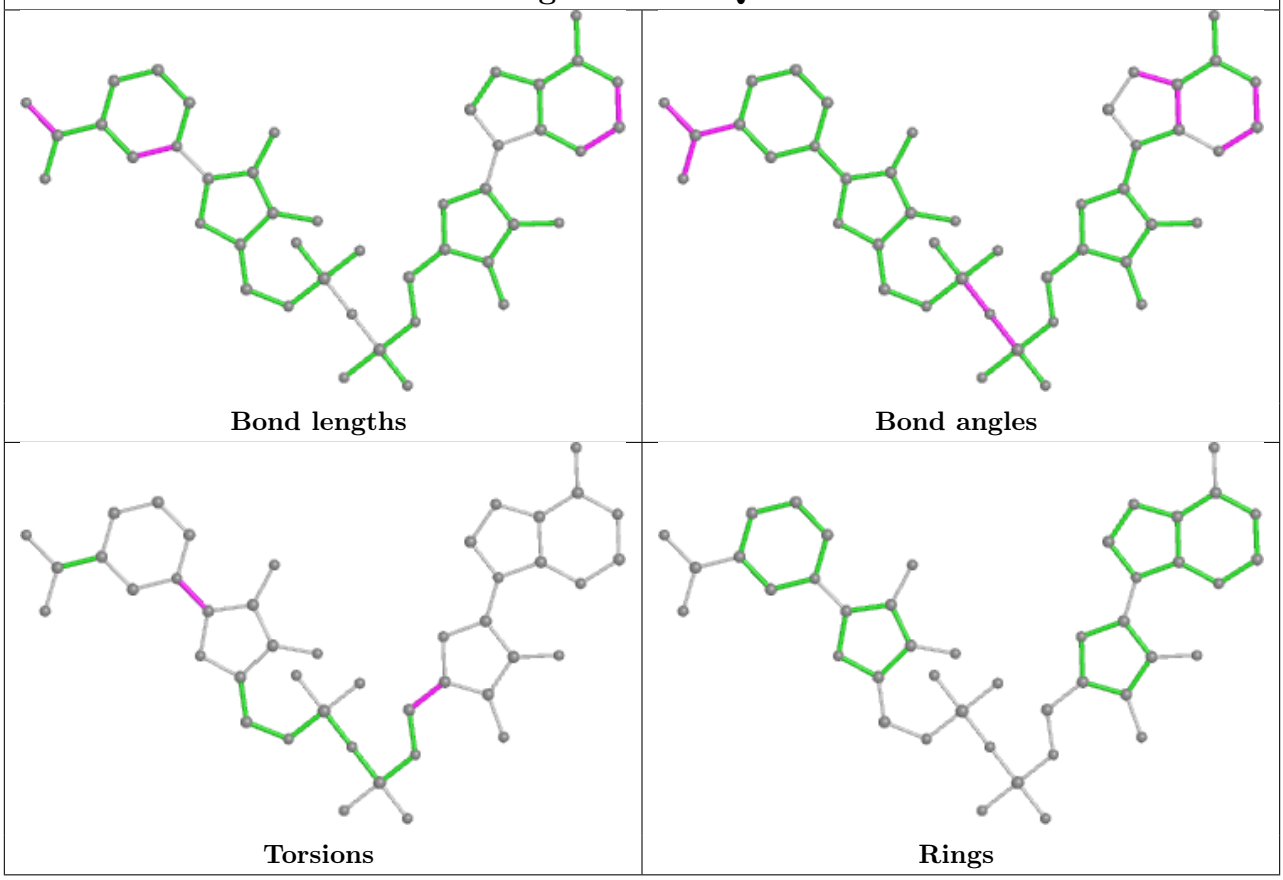


Torsions



Rings

Ligand NAD Q 0



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	O	334/336 (99%)	0.21	14 (4%)	36	35	20, 30, 45, 56	0
1	P	336/336 (100%)	0.28	14 (4%)	36	35	20, 31, 47, 64	0
1	Q	334/336 (99%)	0.51	28 (8%)	11	10	21, 34, 54, 63	0
1	R	336/336 (100%)	0.22	15 (4%)	33	32	21, 30, 44, 67	0
All	All	1340/1344 (99%)	0.30	71 (5%)	26	25	20, 31, 49, 67	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	R	335	SER	6.7
1	R	1	MET	5.8
1	R	336	LYS	4.8
1	Q	99	PHE	4.4
1	P	1	MET	4.0
1	O	62	ASP	3.7
1	Q	62	ASP	3.6
1	R	334	LEU	3.5
1	P	335	SER	3.5
1	Q	304	ASP	3.3
1	O	335	SER	3.3
1	Q	126	ASP	3.3
1	R	66	ARG	3.3
1	Q	2	ALA	3.3
1	O	55	THR	3.2
1	Q	94	LEU	3.2
1	Q	89	ASN	3.2
1	P	126	ASP	3.2
1	Q	25	GLU	3.1
1	Q	210	SER	3.1
1	P	94	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	R	62	ASP	2.8
1	Q	69	GLY	2.8
1	P	55	THR	2.7
1	P	13	ILE	2.7
1	O	151	CYS	2.7
1	O	68	ASN	2.7
1	Q	68	ASN	2.7
1	P	62	ASP	2.6
1	O	57	GLU	2.6
1	O	69	GLY	2.6
1	Q	253	ASP	2.6
1	R	55	THR	2.6
1	R	245	LEU	2.6
1	O	66	ARG	2.5
1	P	7	ILE	2.5
1	R	333	GLU	2.5
1	Q	261	GLU	2.5
1	Q	70	LYS	2.5
1	O	25	GLU	2.4
1	Q	125	GLY	2.4
1	P	69	GLY	2.4
1	P	57	GLU	2.4
1	P	111	GLU	2.4
1	Q	13	ILE	2.4
1	R	69	GLY	2.4
1	Q	71	GLU	2.3
1	R	310	VAL	2.3
1	R	57	GLU	2.3
1	P	137	GLN	2.3
1	R	13	ILE	2.3
1	Q	7	ILE	2.3
1	O	247	VAL	2.3
1	P	140	ASP	2.3
1	Q	87	ASP	2.3
1	Q	96	CYS	2.3
1	Q	53	ARG	2.2
1	Q	92	VAL	2.2
1	Q	3	VAL	2.2
1	Q	24	VAL	2.2
1	R	94	LEU	2.2
1	Q	141	GLY	2.2
1	Q	192	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	O	7	ILE	2.1
1	O	137	GLN	2.1
1	P	177	ILE	2.1
1	O	111	GLU	2.1
1	Q	246	THR	2.1
1	R	138	GLU	2.0
1	Q	93	VAL	2.0
1	O	192	ARG	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)

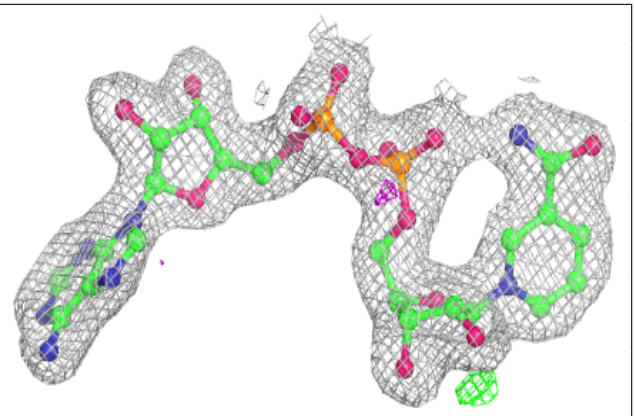
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

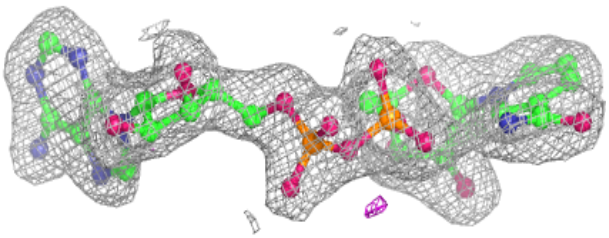
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	P	337	6/6	0.53	0.34	54,55,56,57	0
2	NAD	O	0	44/44	0.93	0.13	23,26,28,30	0
2	NAD	Q	0	44/44	0.95	0.10	21,27,33,33	0
2	NAD	P	0	44/44	0.95	0.12	21,28,33,35	0
2	NAD	R	0	44/44	0.96	0.11	20,25,26,28	0

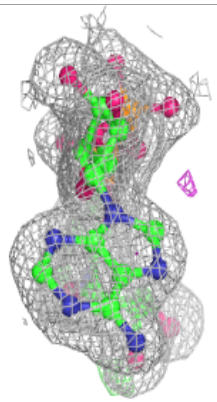
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD O 0:

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

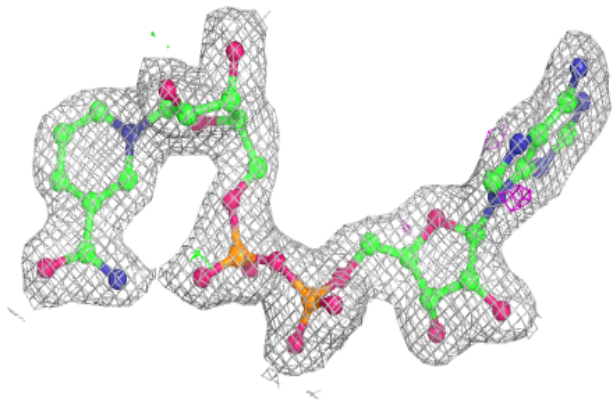


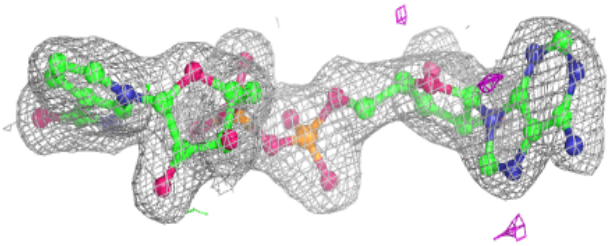


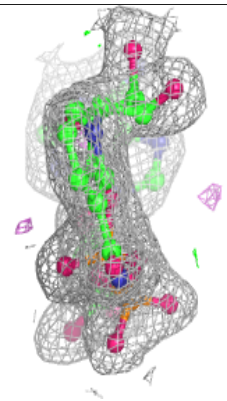


Electron density around NAD Q 0:

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

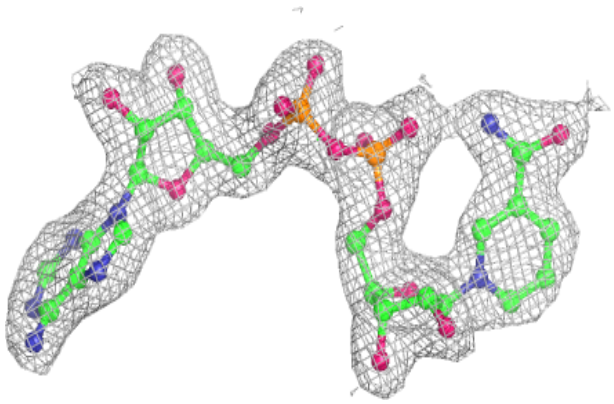


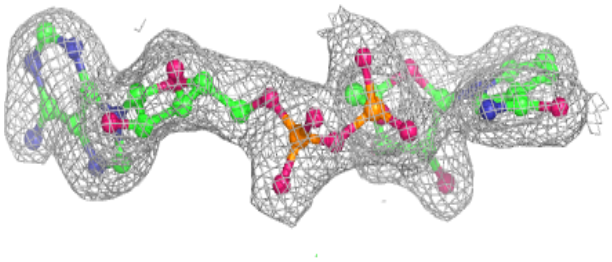


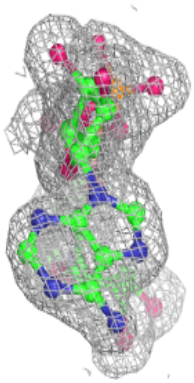


Electron density around NAD P 0:

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

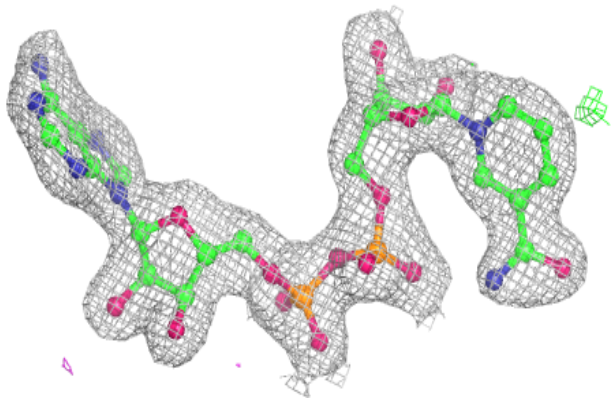


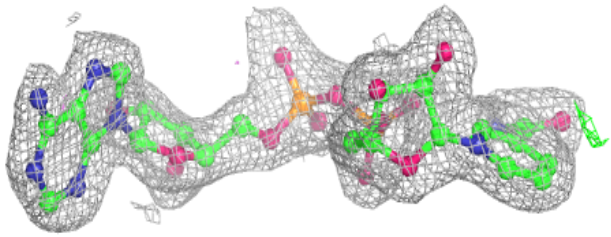


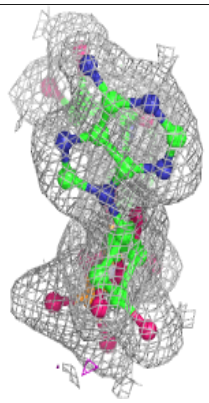


Electron density around NAD R 0:

$2mF_o$-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.