



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 12:05 pm BST

PDB ID : 1LW7  
Title : NADR PROTEIN FROM HAEMOPHILUS INFLUENZAE  
Authors : Singh, S.K.; Kurnasov, O.V.; Chen, B.; Robinson, H.; Grishin, N.V.; Osterman, A.L.; Zhang, H.  
Deposited on : 2002-05-30  
Resolution : 2.90 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

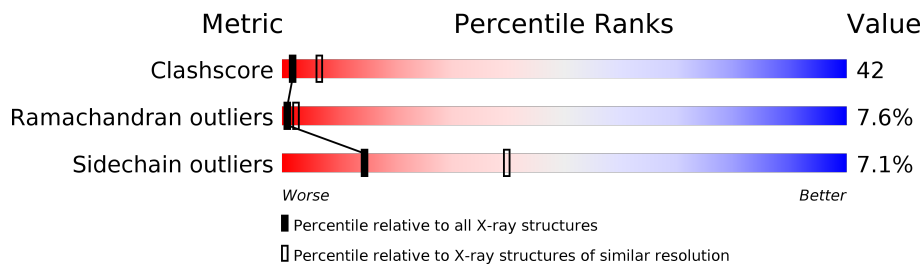
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	365	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2955 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.

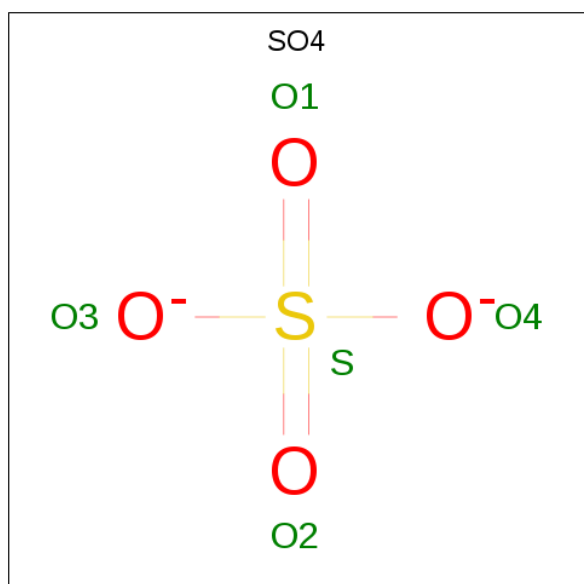
- Molecule 1 is a protein called TRANSCRIPTIONAL REGULATOR NADR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	344	2842	1849	469	515	3	6	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

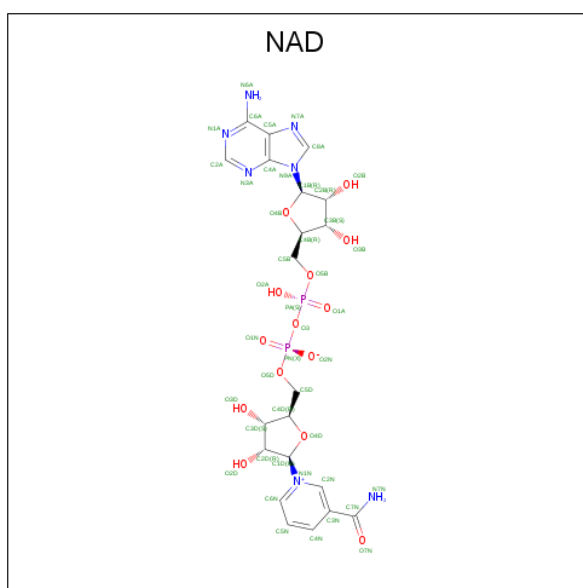
Chain	Residue	Modelled	Actual	Comment	Reference
A	77	MSE	MET	MODIFIED RESIDUE	UNP P44308
A	108	MSE	MET	MODIFIED RESIDUE	UNP P44308
A	111	MSE	MET	MODIFIED RESIDUE	UNP P44308
A	121	MSE	MET	MODIFIED RESIDUE	UNP P44308
A	282	MSE	MET	MODIFIED RESIDUE	UNP P44308
A	328	MSE	MET	MODIFIED RESIDUE	UNP P44308

- Molecule 2 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ).



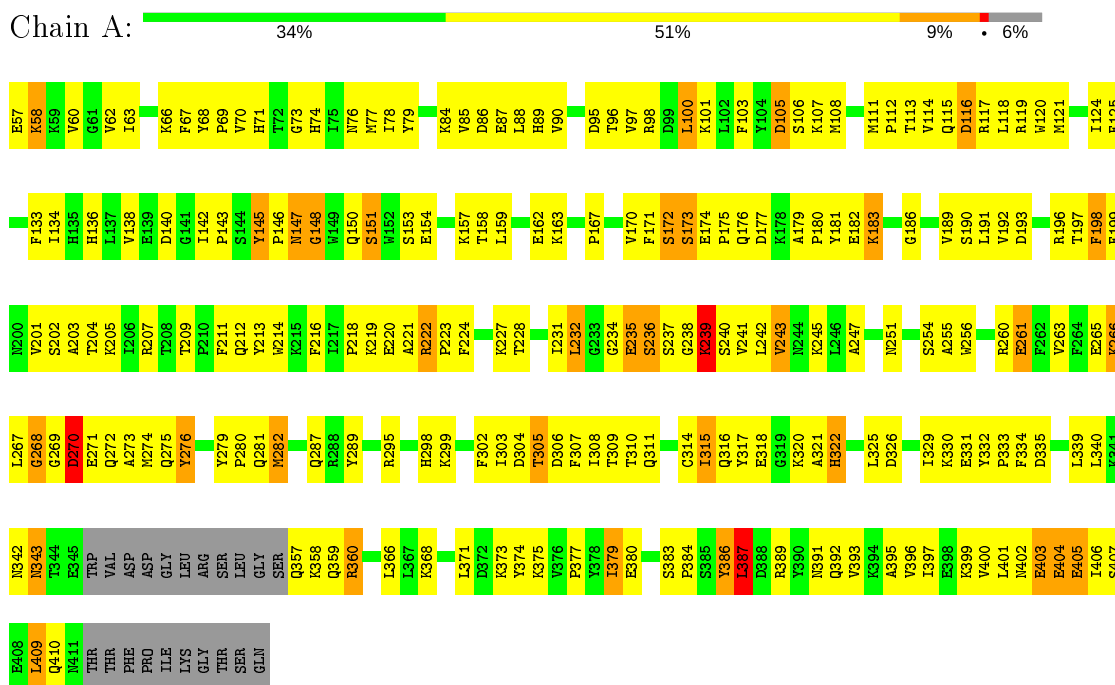
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

Note EDS was not executed.

- Molecule 1: TRANSCRIPTIONAL REGULATOR NADR



## 4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.87Å 106.87Å 174.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.90	Depositor
% Data completeness (in resolution range)	91.5 (30.00-2.90)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.236 , 0.298	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	2955	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

## 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: SO4, 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	0.53	1/2914 (0.0%)	0.74	3/3929 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	282	MSE	CG-SE	-5.04	1.78	1.95

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	268	GLY	N-CA-C	-6.43	97.03	113.10
1	A	266	LYS	N-CA-C	-6.00	94.79	111.00
1	A	173	SER	N-CA-C	-5.93	95.00	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	2842	0	2794	245	0
2	A	25	0	0	0	1
3	A	88	0	52	3	0
All	All	2955	0	2846	245	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 42.

All (245) 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:71:HIS:H	1:A:74:HIS:HD2	1.04	0.93
1:A:373:LYS:HG2	1:A:374:TYR:H	1.35	0.92
1:A:219:LYS:HA	1:A:222:ARG:HD2	1.56	0.88
1:A:240:SER:O	1:A:243:VAL:HG12	1.73	0.88
1:A:340:LEU:HD13	1:A:389:ARG:HD3	1.59	0.84
1:A:396:VAL:HG22	1:A:409:LEU:HD13	1.60	0.83
1:A:409:LEU:HG	1:A:410:GLN:N	1.94	0.83
1:A:71:HIS:H	1:A:74:HIS:CD2	1.96	0.82
1:A:263:VAL:HA	1:A:274:MET:HE3	1.61	0.82
1:A:409:LEU:HG	1:A:410:GLN:H	1.43	0.81
1:A:209:THR:HG21	1:A:212:GLN:NE2	1.95	0.80
1:A:373:LYS:HG2	1:A:374:TYR:N	1.97	0.79
1:A:211:PHE:HB3	1:A:299:LYS:HZ3	1.47	0.78
1:A:231:ILE:O	1:A:239:LYS:HD3	1.84	0.77
1:A:113:THR:HG22	1:A:114:VAL:H	1.48	0.77
1:A:173:SER:HB2	3:A:601:NAD:O2B	1.85	0.76
1:A:85:VAL:HG12	1:A:87:GLU:H	1.48	0.76
1:A:266:LYS:C	1:A:268:GLY:H	1.88	0.75
1:A:232:LEU:HD23	1:A:232:LEU:N	2.05	0.71
1:A:70:VAL:HG21	1:A:121:MSE:SE	2.41	0.71
1:A:406:ILE:HG23	1:A:409:LEU:HD12	1.73	0.70
1:A:179:ALA:HB3	1:A:180:PRO:HD3	1.72	0.70
1:A:227:LYS:HA	1:A:335:ASP:OD2	1.89	0.70
1:A:113:THR:HG22	1:A:114:VAL:N	2.07	0.70
1:A:358:LYS:HB3	1:A:360:ARG:HD2	1.71	0.70
1:A:67:PHE:HB3	1:A:70:VAL:HG22	1.71	0.70
1:A:214:TRP:CE2	1:A:298:HIS:O	2.44	0.70
1:A:115:GLN:HE21	1:A:119:ARG:HH12	1.41	0.69
1:A:150:GLN:O	1:A:154:GLU:HG2	1.93	0.69
1:A:60:VAL:HG12	1:A:167:PRO:HA	1.74	0.69
1:A:379:ILE:HD11	1:A:396:VAL:HG21	1.76	0.68
1:A:170:VAL:CG1	1:A:189:VAL:HG22	2.25	0.67
1:A:311:GLN:HE22	1:A:321:ALA:CA	2.07	0.67
1:A:57:GLU:O	1:A:58:LYS:HB3	1.94	0.67
1:A:158:THR:HG22	1:A:162:GLU:OE1	1.95	0.66
1:A:231:ILE:HD12	1:A:243:VAL:HG23	1.78	0.66
1:A:63:ILE:HD13	1:A:77:MSE:HE2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:TYR:HB3	1:A:146:PRO:HD3	1.77	0.65
1:A:280:PRO:HB2	1:A:281:GLN:NE2	2.11	0.65
1:A:97:VAL:HG12	1:A:101:LYS:HE3	1.78	0.65
1:A:266:LYS:HD2	1:A:270:ASP:O	1.97	0.65
1:A:66:LYS:HG2	1:A:98:ARG:HH12	1.60	0.65
1:A:305:THR:HG23	1:A:309:THR:HB	1.78	0.64
1:A:375:LYS:O	1:A:377:PRO:HD3	1.98	0.64
1:A:266:LYS:O	1:A:268:GLY:N	2.31	0.64
1:A:209:THR:HG21	1:A:212:GLN:HE21	1.62	0.64
1:A:235:GLU:O	1:A:237:SER:N	2.31	0.63
1:A:236:SER:C	1:A:238:GLY:H	2.00	0.63
1:A:402:ASN:ND2	1:A:403:GLU:HG3	2.13	0.63
1:A:379:ILE:HD12	1:A:409:LEU:HB2	1.81	0.62
1:A:85:VAL:HG12	1:A:86:ASP:N	2.14	0.62
1:A:240:SER:HB3	1:A:260:ARG:CZ	2.30	0.62
1:A:387:LEU:CD1	1:A:391:ASN:HD21	2.13	0.62
1:A:228:THR:HB	1:A:303:ILE:CD1	2.30	0.61
1:A:95:ASP:HB3	1:A:98:ARG:HB3	1.81	0.61
1:A:203:ALA:O	1:A:207:ARG:HB2	2.00	0.61
1:A:179:ALA:O	1:A:183:LYS:N	2.31	0.61
1:A:266:LYS:C	1:A:268:GLY:N	2.54	0.61
1:A:402:ASN:O	1:A:403:GLU:O	2.19	0.61
1:A:175:PRO:O	1:A:177:ASP:N	2.33	0.60
1:A:101:LYS:O	1:A:105:ASP:HB2	2.02	0.60
1:A:153:SER:O	1:A:157:LYS:HG3	2.02	0.60
1:A:211:PHE:HB3	1:A:299:LYS:NZ	2.16	0.59
1:A:387:LEU:HD12	1:A:391:ASN:HD21	1.67	0.59
1:A:316:GLN:HG2	1:A:317:TYR:CD1	2.37	0.59
1:A:170:VAL:HG12	1:A:189:VAL:HG22	1.84	0.59
1:A:311:GLN:HE22	1:A:321:ALA:C	2.06	0.59
1:A:316:GLN:HG2	1:A:317:TYR:CE1	2.37	0.59
1:A:115:GLN:HE21	1:A:119:ARG:NH1	2.00	0.59
1:A:373:LYS:O	1:A:374:TYR:HB2	2.03	0.59
1:A:62:VAL:HG22	1:A:167:PRO:CB	2.33	0.58
1:A:325:LEU:O	1:A:329:ILE:HG12	2.03	0.58
1:A:205:LYS:HB3	1:A:213:TYR:CE2	2.39	0.57
1:A:147:ASN:HD22	1:A:148:GLY:N	2.02	0.57
1:A:177:ASP:C	1:A:180:PRO:HD2	2.25	0.57
1:A:76:ASN:ND2	1:A:199:PHE:HE2	2.03	0.57
1:A:62:VAL:HG12	1:A:63:ILE:H	1.70	0.57
1:A:68:TYR:HE1	1:A:108:MSE:HE1	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:THR:O	1:A:100:LEU:HB2	2.05	0.57
1:A:392:GLN:O	1:A:395:ALA:HB3	2.05	0.57
1:A:358:LYS:HG2	1:A:359:GLN:H	1.70	0.56
1:A:311:GLN:NE2	1:A:321:ALA:O	2.39	0.56
1:A:311:GLN:HE22	1:A:321:ALA:HA	1.69	0.56
1:A:266:LYS:O	1:A:267:LEU:HB2	2.06	0.55
1:A:62:VAL:CG2	1:A:167:PRO:CB	2.85	0.55
1:A:88:LEU:HD23	1:A:89:HIS:N	2.22	0.55
1:A:175:PRO:C	1:A:177:ASP:H	2.10	0.55
1:A:303:ILE:HD13	1:A:334:PHE:CE2	2.41	0.55
1:A:228:THR:HB	1:A:303:ILE:HD11	1.88	0.54
1:A:204:THR:HA	1:A:207:ARG:NH1	2.22	0.54
1:A:303:ILE:HD13	1:A:334:PHE:CD2	2.42	0.54
1:A:71:HIS:HD2	1:A:73:GLY:CA	2.20	0.54
1:A:228:THR:CG2	1:A:303:ILE:HD11	2.38	0.54
1:A:62:VAL:HG22	1:A:167:PRO:CG	2.38	0.54
1:A:373:LYS:C	1:A:375:LYS:H	2.11	0.54
1:A:115:GLN:NE2	1:A:119:ARG:HH22	2.06	0.53
1:A:189:VAL:HG12	1:A:190:SER:N	2.23	0.53
1:A:112:PRO:HG3	1:A:224:PHE:CD1	2.43	0.53
1:A:389:ARG:O	1:A:393:VAL:HG23	2.08	0.53
1:A:103:PHE:HD1	1:A:108:MSE:HE3	1.73	0.53
1:A:191:LEU:H	1:A:191:LEU:HD23	1.74	0.53
1:A:279:TYR:HB2	1:A:280:PRO:HD3	1.91	0.53
1:A:193:ASP:HB2	1:A:198:PHE:HE1	1.74	0.53
1:A:120:TRP:O	1:A:124:ILE:HG13	2.08	0.53
1:A:182:GLU:O	1:A:186:GLY:HA2	2.08	0.53
1:A:266:LYS:HE3	1:A:271:GLU:O	2.08	0.53
1:A:371:LEU:O	1:A:373:LYS:O	2.27	0.53
1:A:198:PHE:N	1:A:198:PHE:CD1	2.77	0.52
1:A:201:VAL:HG22	1:A:202:SER:H	1.74	0.52
1:A:315:ILE:HA	1:A:320:LYS:O	2.10	0.52
1:A:331:GLU:HA	1:A:331:GLU:OE1	2.09	0.52
1:A:402:ASN:HD22	1:A:403:GLU:HG3	1.75	0.51
1:A:241:VAL:HG13	1:A:242:LEU:N	2.25	0.51
1:A:88:LEU:HD22	1:A:90:VAL:HG23	1.93	0.51
1:A:263:VAL:HG22	1:A:274:MET:HE1	1.92	0.51
1:A:211:PHE:CD1	1:A:211:PHE:N	2.79	0.50
1:A:85:VAL:HG12	1:A:86:ASP:H	1.76	0.50
1:A:145:TYR:O	1:A:146:PRO:C	2.50	0.50
1:A:214:TRP:NE1	1:A:298:HIS:O	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:SER:OG	1:A:108:MSE:HG3	2.12	0.50
1:A:171:PHE:O	1:A:172:SER:HB2	2.12	0.50
1:A:234:GLY:C	1:A:236:SER:H	2.13	0.49
1:A:193:ASP:HB2	1:A:198:PHE:CE1	2.46	0.49
1:A:138:VAL:HG23	1:A:138:VAL:O	2.13	0.49
1:A:201:VAL:HG22	1:A:202:SER:N	2.27	0.49
1:A:219:LYS:HE3	1:A:295:ARG:HH21	1.78	0.49
1:A:406:ILE:CG2	1:A:409:LEU:HA	2.42	0.49
1:A:71:HIS:CD2	1:A:74:HIS:H	2.30	0.49
1:A:254:SER:HA	1:A:302:PHE:O	2.12	0.49
1:A:298:HIS:O	1:A:299:LYS:HG3	2.12	0.49
1:A:331:GLU:HB3	1:A:332:TYR:CD1	2.47	0.49
1:A:67:PHE:HB3	1:A:70:VAL:CG2	2.40	0.49
1:A:404:GLU:O	1:A:405:GLU:C	2.52	0.49
1:A:199:PHE:HA	1:A:216:PHE:HE1	1.77	0.48
1:A:331:GLU:O	1:A:333:PRO:HD3	2.13	0.48
1:A:219:LYS:HA	1:A:222:ARG:CD	2.35	0.48
1:A:60:VAL:HG22	1:A:87:GLU:HB3	1.94	0.48
1:A:113:THR:HB	1:A:116:ASP:OD1	2.13	0.48
1:A:133:PHE:C	1:A:134:ILE:HG13	2.33	0.48
1:A:211:PHE:CB	1:A:299:LYS:HZ3	2.23	0.48
1:A:159:LEU:O	1:A:163:LYS:HB2	2.14	0.48
1:A:383:SER:HA	1:A:384:PRO:HD3	1.68	0.48
1:A:326:ASP:OD1	1:A:374:TYR:OH	2.29	0.47
1:A:62:VAL:HG12	1:A:63:ILE:N	2.29	0.47
1:A:181:TYR:O	1:A:186:GLY:N	2.48	0.47
1:A:68:TYR:CE1	1:A:108:MSE:HE1	2.49	0.47
1:A:221:ALA:O	1:A:223:PRO:N	2.47	0.47
1:A:396:VAL:HG22	1:A:409:LEU:CD1	2.40	0.47
1:A:236:SER:C	1:A:238:GLY:N	2.67	0.47
1:A:221:ALA:O	1:A:222:ARG:C	2.52	0.47
1:A:142:ILE:HG23	1:A:151:SER:HB2	1.97	0.47
1:A:114:VAL:HG11	1:A:136:HIS:CE1	2.49	0.47
1:A:220:GLU:N	1:A:220:GLU:OE1	2.48	0.47
1:A:71:HIS:N	1:A:74:HIS:HD2	1.89	0.47
1:A:117:ARG:HG2	1:A:120:TRP:CZ3	2.50	0.46
1:A:193:ASP:OD1	1:A:197:THR:N	2.42	0.46
1:A:387:LEU:O	1:A:391:ASN:ND2	2.48	0.46
1:A:211:PHE:HD1	1:A:211:PHE:N	2.14	0.46
1:A:219:LYS:CA	1:A:222:ARG:HD2	2.36	0.46
1:A:63:ILE:HD13	1:A:77:MSE:CE	2.42	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:MET:HE1	1:A:282:MSE:HE1	1.98	0.46
1:A:211:PHE:O	1:A:299:LYS:NZ	2.44	0.46
1:A:311:GLN:HE21	1:A:315:ILE:HG13	1.80	0.46
1:A:57:GLU:O	1:A:58:LYS:CB	2.63	0.46
1:A:198:PHE:O	1:A:199:PHE:HB2	2.15	0.45
1:A:399:LYS:HG2	1:A:403:GLU:OE1	2.16	0.45
1:A:196:ARG:HD3	1:A:196:ARG:N	2.31	0.45
1:A:265:GLU:C	1:A:266:LYS:O	2.50	0.45
1:A:379:ILE:CG1	1:A:379:ILE:O	2.65	0.45
1:A:113:THR:CG2	1:A:114:VAL:N	2.77	0.45
1:A:307:PHE:HB2	1:A:329:ILE:HD13	1.97	0.45
1:A:263:VAL:HG22	1:A:274:MET:CE	2.46	0.45
1:A:95:ASP:HA	1:A:140:ASP:OD1	2.16	0.45
1:A:366:LEU:O	1:A:366:LEU:HD12	2.16	0.45
1:A:321:ALA:O	1:A:322:HIS:CB	2.65	0.45
1:A:193:ASP:OD2	1:A:198:PHE:HD1	1.99	0.45
1:A:234:GLY:C	1:A:236:SER:N	2.71	0.45
1:A:308:ILE:HG12	1:A:329:ILE:HD12	1.99	0.45
1:A:235:GLU:C	1:A:237:SER:N	2.70	0.44
1:A:173:SER:O	1:A:174:GLU:HB2	2.18	0.44
1:A:228:THR:CB	1:A:303:ILE:HD11	2.48	0.44
1:A:107:LYS:HD3	1:A:107:LYS:HA	1.77	0.44
1:A:172:SER:OG	1:A:173:SER:N	2.51	0.44
1:A:113:THR:CG2	1:A:114:VAL:H	2.24	0.44
1:A:228:THR:HG22	1:A:303:ILE:HD11	1.99	0.44
1:A:238:GLY:O	1:A:241:VAL:CG1	2.66	0.44
1:A:409:LEU:CG	1:A:410:GLN:H	2.22	0.44
1:A:177:ASP:O	1:A:180:PRO:HD2	2.18	0.44
1:A:397:ILE:O	1:A:400:VAL:N	2.49	0.44
1:A:84:LYS:N	1:A:84:LYS:HD2	2.32	0.44
1:A:71:HIS:HD2	1:A:73:GLY:HA3	1.82	0.43
1:A:147:ASN:HD22	1:A:148:GLY:H	1.67	0.43
1:A:311:GLN:NE2	1:A:321:ALA:HA	2.32	0.43
1:A:326:ASP:OD2	1:A:330:LYS:HE3	2.18	0.43
1:A:173:SER:HB2	3:A:601:NAD:HO2A	1.80	0.43
1:A:406:ILE:HG23	1:A:409:LEU:HA	2.01	0.43
1:A:307:PHE:HE2	1:A:332:TYR:HB2	1.83	0.43
1:A:189:VAL:CG1	1:A:190:SER:N	2.82	0.43
1:A:238:GLY:O	1:A:241:VAL:HG12	2.19	0.43
1:A:310:THR:HG22	1:A:325:LEU:CD1	2.48	0.43
1:A:308:ILE:CG1	1:A:329:ILE:HD12	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:386:TYR:O	1:A:387:LEU:C	2.57	0.43
1:A:174:GLU:OE2	1:A:196:ARG:NH2	2.52	0.43
1:A:62:VAL:HG22	1:A:167:PRO:HG2	2.00	0.43
1:A:239:LYS:CB	1:A:239:LYS:NZ	2.82	0.42
1:A:270:ASP:HB3	1:A:273:ALA:HB2	2.01	0.42
1:A:272:GLN:HA	1:A:275:GLN:NE2	2.34	0.42
1:A:373:LYS:C	1:A:375:LYS:N	2.72	0.42
1:A:339:LEU:HB3	1:A:380:GLU:HA	2.01	0.42
1:A:310:THR:HG22	1:A:325:LEU:HD13	2.01	0.42
1:A:71:HIS:CE1	1:A:74:HIS:NE2	2.87	0.42
1:A:71:HIS:CD2	1:A:73:GLY:HA3	2.54	0.42
1:A:112:PRO:HG3	1:A:224:PHE:CG	2.54	0.42
1:A:255:ALA:HB1	1:A:289:TYR:HB3	2.00	0.42
1:A:148:GLY:O	1:A:151:SER:N	2.52	0.42
1:A:173:SER:O	1:A:174:GLU:CB	2.68	0.42
1:A:368:LYS:HE2	1:A:380:GLU:OE1	2.20	0.42
1:A:85:VAL:CG1	1:A:86:ASP:N	2.82	0.42
1:A:281:GLN:N	1:A:281:GLN:CD	2.73	0.41
1:A:305:THR:HG23	1:A:306:ASP:N	2.35	0.41
1:A:74:HIS:O	1:A:78:ILE:HG13	2.19	0.41
1:A:263:VAL:HG12	1:A:267:LEU:HD12	2.01	0.41
1:A:175:PRO:C	1:A:177:ASP:N	2.69	0.41
1:A:306:ASP:OD1	1:A:308:ILE:HB	2.20	0.41
1:A:276:TYR:HB2	1:A:318:GLU:OE1	2.19	0.41
1:A:247:ALA:HA	1:A:302:PHE:CE2	2.56	0.41
1:A:342:ASN:CB	1:A:389:ARG:HH22	2.33	0.41
1:A:76:ASN:ND2	1:A:199:PHE:CE2	2.86	0.41
1:A:111:MSE:HA	1:A:112:PRO:HD3	1.91	0.41
1:A:118:LEU:HD11	1:A:136:HIS:HB2	2.03	0.41
1:A:403:GLU:HG3	1:A:403:GLU:H	1.73	0.41
1:A:279:TYR:HB3	1:A:325:LEU:HD11	2.02	0.41
1:A:325:LEU:HA	1:A:325:LEU:HD23	1.92	0.40
1:A:307:PHE:CB	1:A:329:ILE:HD13	2.51	0.40
1:A:332:TYR:CD1	1:A:332:TYR:N	2.88	0.40
1:A:143:PRO:O	3:A:601:NAD:N7N	2.52	0.40
1:A:192:VAL:O	1:A:193:ASP:C	2.60	0.40
1:A:261:GLU:HG3	1:A:261:GLU:H	1.53	0.40
1:A:79:TYR:CZ	1:A:125:PHE:CE1	3.09	0.40
1:A:221:ALA:C	1:A:223:PRO:HD2	2.42	0.40
1:A:303:ILE:HG22	1:A:305:THR:O	2.21	0.40
1:A:90:VAL:O	1:A:134:ILE:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:VAL:O	1:A:218:PRO:CD	2.70	0.40
1:A:58:LYS:HG2	1:A:58:LYS:O	2.20	0.40
1:A:71:HIS:HD2	1:A:73:GLY:N	2.19	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)
2:A:603:SO4:O4	2:A:603:SO4:O4[12_555]	1.75	0.45

## 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	340/365 (93%)	271 (80%)	43 (13%)	26 (8%)	<a href="#">1</a> <a href="#">2</a>

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	TYR
1	A	176	GLN
1	A	236	SER
1	A	403	GLU
1	A	405	GLU
1	A	58	LYS
1	A	256	TRP
1	A	270	ASP
1	A	304	ASP
1	A	343	ASN
1	A	407	SER
1	A	409	LEU
1	A	269	GLY

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Mol	Chain	Res	Type
1	A	276	TYR
1	A	314	CYS
1	A	404	GLU
1	A	235	GLU
1	A	322	HIS
1	A	386	TYR
1	A	172	SER
1	A	239	LYS
1	A	315	ILE
1	A	183	LYS
1	A	387	LEU
1	A	222	ARG
1	A	148	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	A	309/325 (95%)	287 (93%)	22 (7%)	14 40

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	PRO
1	A	100	LEU
1	A	105	ASP
1	A	116	ASP
1	A	147	ASN
1	A	151	SER
1	A	198	PHE
1	A	232	LEU
1	A	239	LYS
1	A	243	VAL
1	A	245	LYS
1	A	251	ASN
1	A	261	GLU

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Mol	Chain	Res	Type
1	A	270	ASP
1	A	287	GLN
1	A	305	THR
1	A	343	ASN
1	A	357	GLN
1	A	360	ARG
1	A	379	ILE
1	A	387	LEU
1	A	401	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	HIS
1	A	74	HIS
1	A	115	GLN
1	A	123	GLN
1	A	147	ASN
1	A	176	GLN
1	A	212	GLN
1	A	275	GLN
1	A	281	GLN
1	A	287	GLN
1	A	296	HIS
1	A	311	GLN
1	A	343	ASN
1	A	357	GLN
1	A	364	GLN
1	A	391	ASN
1	A	402	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 carbohydrates 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	SO4	A	603	-	4,4,4	0.20	0	6,6,6	0.28	0
2	SO4	A	607	-	4,4,4	0.23	0	6,6,6	0.28	0
2	SO4	A	606	-	4,4,4	0.27	0	6,6,6	0.22	0
2	SO4	A	501	-	4,4,4	0.29	0	6,6,6	0.24	0
3	NAD	A	605	-	42,48,48	1.90	10 (23%)	50,73,73	1.79	11 (22%)
3	NAD	A	601	-	42,48,48	2.16	10 (23%)	50,73,73	1.77	9 (18%)
2	SO4	A	604	-	4,4,4	0.21	0	6,6,6	0.21	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
3	NAD	A	601	-	-	5/26/62/62	0/5/5/5
3	NAD	A	605	-	-	9/26/62/62	0/5/5/5

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	NAD	C5N-C4N	6.43	1.52	1.38
3	A	601	NAD	C2N-N1N	6.14	1.42	1.35
3	A	605	NAD	C5N-C4N	5.83	1.51	1.38
3	A	601	NAD	C4N-C3N	5.23	1.48	1.39
3	A	601	NAD	C2N-C3N	5.21	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	605	NAD	C2N-N1N	4.40	1.40	1.35
3	A	605	NAD	C4N-C3N	4.26	1.46	1.39
3	A	605	NAD	C2N-C3N	3.94	1.45	1.39
3	A	601	NAD	C6N-N1N	3.80	1.44	1.35
3	A	605	NAD	O4B-C1B	-3.56	1.36	1.41
3	A	605	NAD	C6N-C5N	-3.06	1.31	1.38
3	A	601	NAD	C2A-N1A	2.92	1.39	1.33
3	A	605	NAD	C6N-N1N	2.92	1.42	1.35
3	A	605	NAD	C7N-N7N	2.84	1.38	1.33
3	A	601	NAD	O4D-C1D	2.64	1.44	1.41
3	A	601	NAD	C6N-C5N	-2.50	1.33	1.38
3	A	601	NAD	C7N-N7N	2.33	1.37	1.33
3	A	601	NAD	O4B-C1B	-2.17	1.38	1.41
3	A	605	NAD	C2A-N1A	2.10	1.37	1.33
3	A	605	NAD	C2D-C3D	-2.05	1.47	1.53

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	NAD	C3N-C7N-N7N	6.07	125.03	117.75
3	A	601	NAD	O7N-C7N-C3N	-5.77	112.73	119.63
3	A	605	NAD	O7N-C7N-C3N	-5.54	113.01	119.63
3	A	605	NAD	C3N-C7N-N7N	5.44	124.28	117.75
3	A	601	NAD	C6N-N1N-C2N	-4.31	118.04	121.97
3	A	605	NAD	C6N-N1N-C2N	-4.26	118.09	121.97
3	A	601	NAD	O4B-C1B-C2B	-3.78	101.40	106.93
3	A	605	NAD	C5N-C4N-C3N	-3.27	116.48	120.34
3	A	605	NAD	O5B-C5B-C4B	-3.16	98.10	108.99
3	A	605	NAD	PN-O3-PA	-3.08	122.27	132.83
3	A	605	NAD	C5A-C6A-N6A	3.04	124.97	120.35
3	A	601	NAD	C5N-C4N-C3N	-2.97	116.83	120.34
3	A	601	NAD	C5N-C6N-N1N	2.69	124.26	120.40
3	A	605	NAD	C3B-C2B-C1B	2.66	104.99	100.98
3	A	605	NAD	C5N-C6N-N1N	2.66	124.22	120.40
3	A	601	NAD	C3B-C2B-C1B	2.37	104.55	100.98
3	A	601	NAD	C4A-C5A-N7A	2.29	111.78	109.40
3	A	601	NAD	C5A-C6A-N6A	2.27	123.81	120.35
3	A	605	NAD	C4A-C5A-N7A	2.09	111.58	109.40
3	A	605	NAD	O4B-C1B-C2B	-2.04	103.95	106.93

There are no chirality outliers.

All (14) torsion outliers are listed below:

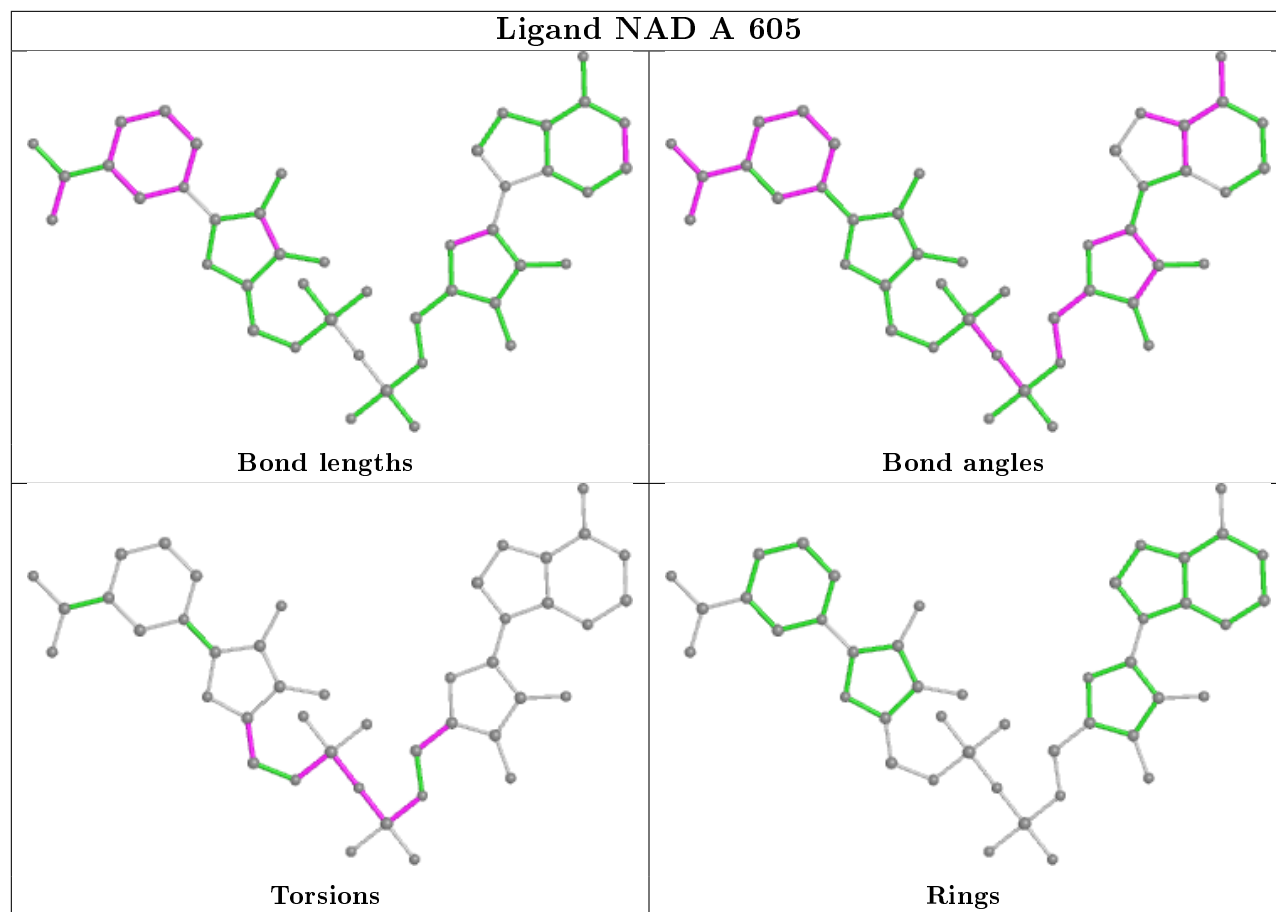
Mol	Chain	Res	Type	Atoms
3	A	605	NAD	C5B-O5B-PA-O1A
3	A	605	NAD	PN-O3-PA-O5B
3	A	605	NAD	C5D-O5D-PN-O3
3	A	605	NAD	C5D-O5D-PN-O1N
3	A	605	NAD	O4D-C4D-C5D-O5D
3	A	605	NAD	C3D-C4D-C5D-O5D
3	A	601	NAD	C5D-O5D-PN-O3
3	A	601	NAD	O4D-C4D-C5D-O5D
3	A	605	NAD	PA-O3-PN-O1N
3	A	601	NAD	C5D-O5D-PN-O1N
3	A	605	NAD	PA-O3-PN-O2N
3	A	601	NAD	PN-O3-PA-O1A
3	A	601	NAD	PN-O3-PA-O5B
3	A	605	NAD	C3B-C4B-C5B-O5B

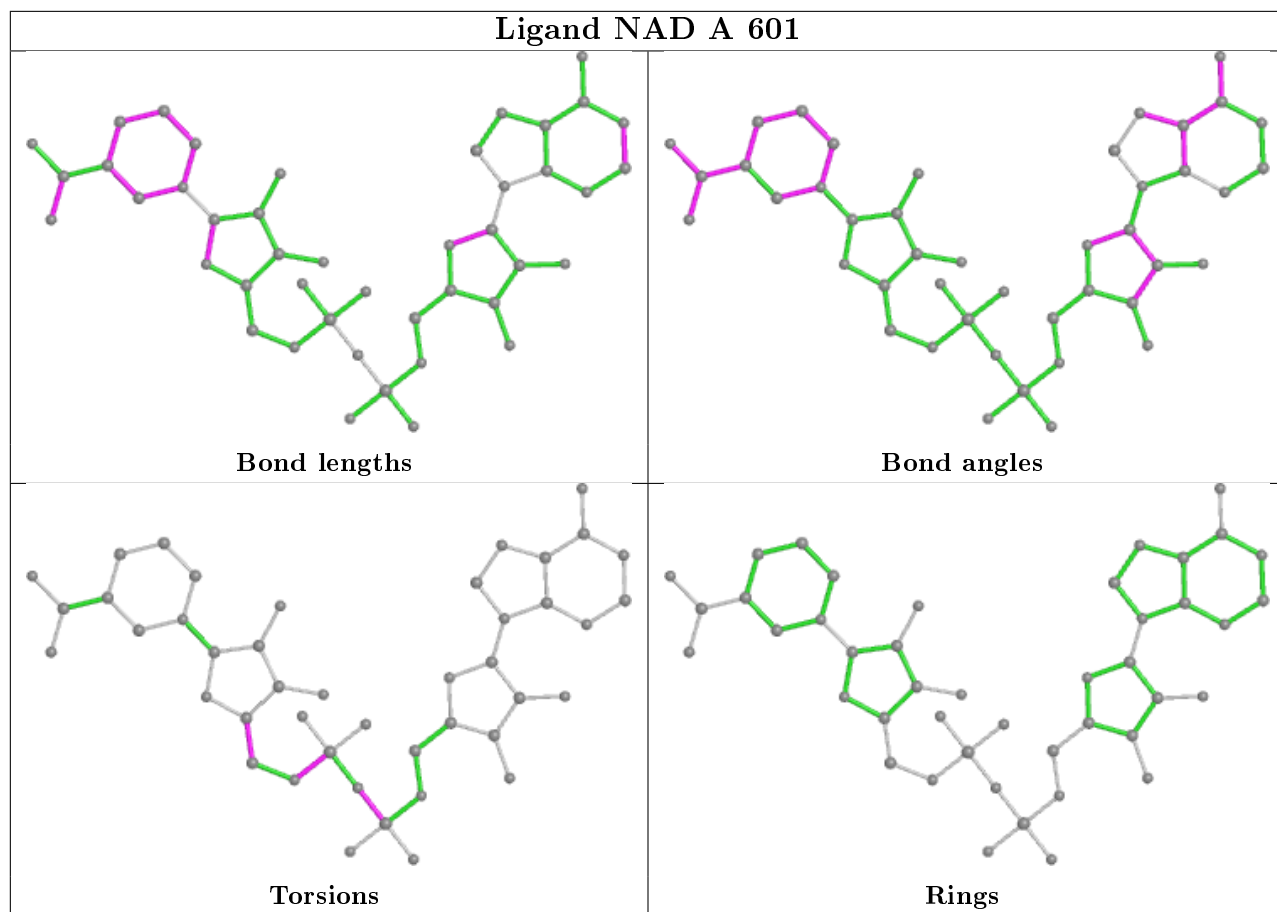
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	603	SO4	0	1
3	A	601	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

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.