



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:35 AM UTC

PDB ID : 1DC6 / pdb_00001dc6
Title : STRUCTURAL ANALYSIS OF GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE FROM ESCHERICHIA COLI: DIRECT EVIDENCE FOR SUBSTRATE BINDING AND COFACTOR-INDUCED CONFORMATIONAL CHANGES.
Authors : Yun, M.; Park, C.-G.; Kim, J.-Y.; Park, H.-W.
Deposited on : 1999-11-04
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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtrriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

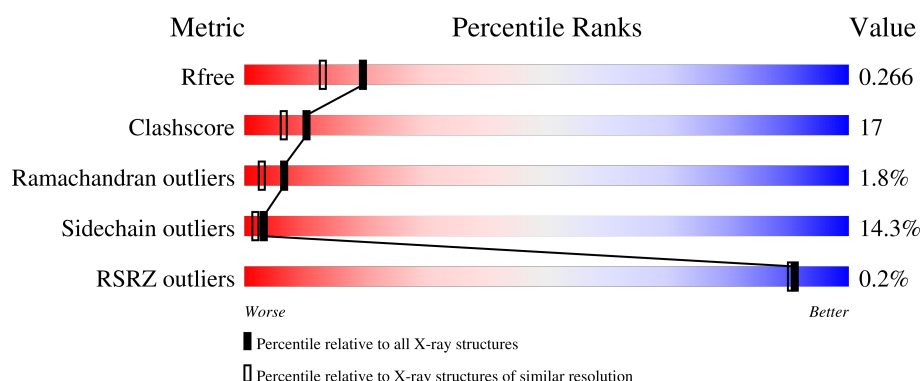
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	330	 65% 28% 7%
1	B	330	 55% 36% 8% .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5321 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2488	1563	431	484	10			
1	B	330	Total	C	N	O	S	0	0	0
			2488	1563	431	484	10			

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



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

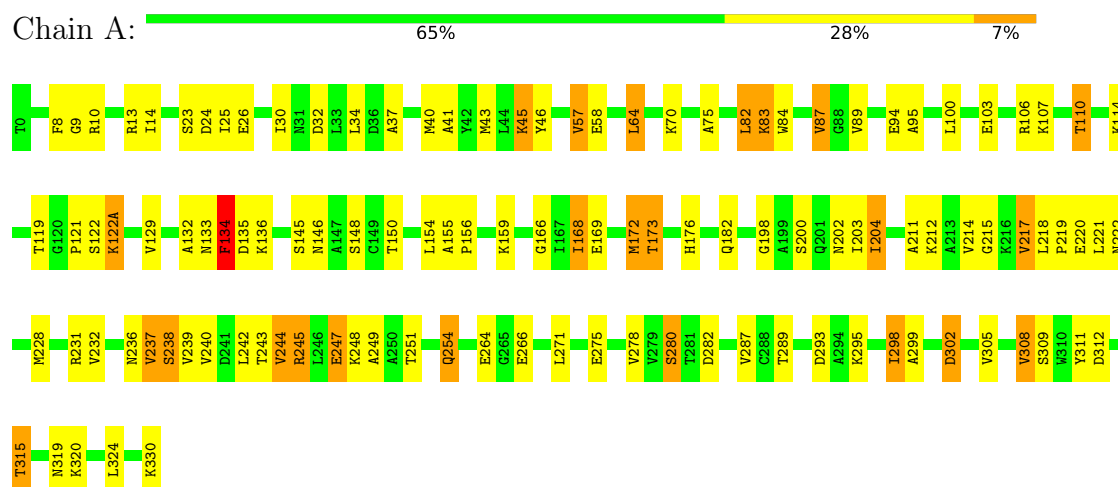
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	134	Total 134	O 134	0	0
3	B	123	Total 123	O 123	0	0

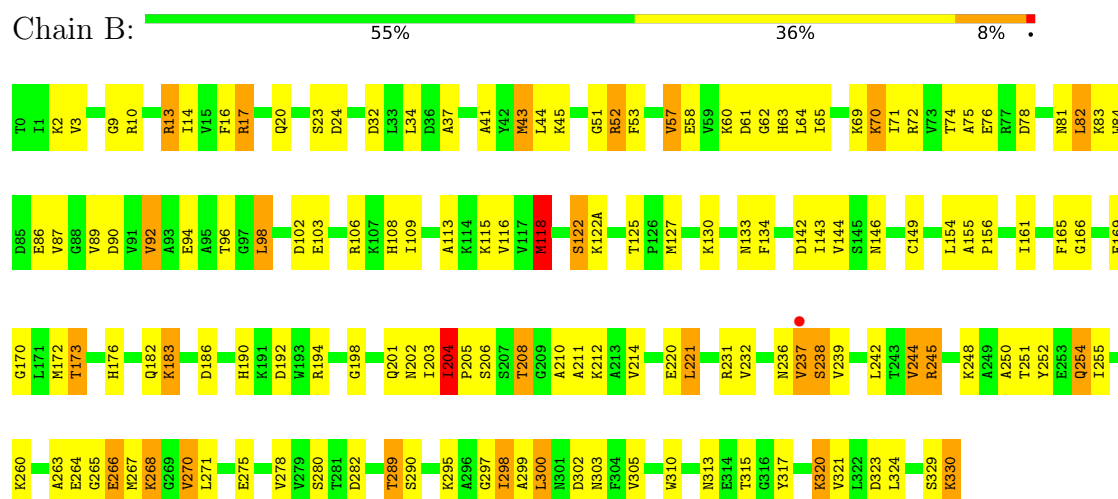
3 Residue-property plots

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

• Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE



• Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.45Å 134.47Å 67.34Å 90.00° 107.89° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.00) 81.9 (20.00-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 2.01Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.224 , 0.291 0.204 , 0.266	Depositor DCC
R_{free} test set	2245 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.599	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 74.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5321	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
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	0.55	1/2527 (0.0%)	0.97	10/3421 (0.3%)
1	B	0.55	3/2527 (0.1%)	1.04	13/3421 (0.4%)
All	All	0.55	4/5054 (0.1%)	1.00	23/6842 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	43	MET	SD-CE	-7.29	1.61	1.79
1	A	43	MET	SD-CE	-6.25	1.64	1.79
1	B	127	MET	SD-CE	-6.07	1.64	1.79
1	B	118	MET	SD-CE	-5.59	1.65	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	203	ILE	N-CA-C	-10.11	92.41	107.37
1	B	204	ILE	N-CA-C	8.57	116.77	107.60
1	A	203	ILE	N-CA-C	-8.15	95.36	107.51
1	A	204	ILE	N-CA-C	7.82	115.97	107.60
1	B	202	ASN	N-CA-C	7.21	121.25	109.85
1	A	202	ASN	N-CA-C	6.83	119.80	109.41
1	B	315	THR	N-CA-C	6.71	118.67	111.36
1	B	82	LEU	N-CA-C	6.46	120.89	112.89
1	B	109	ILE	N-CA-C	-6.22	104.21	111.00
1	A	82	LEU	N-CA-C	5.95	120.38	113.18
1	A	280	SER	N-CA-C	5.88	118.16	111.11
1	B	133	ASN	N-CA-C	5.83	118.34	110.88
1	B	278	VAL	N-CA-C	5.78	116.74	110.21
1	A	168	ILE	N-CA-C	-5.69	105.07	110.42
1	A	134	PHE	N-CA-C	-5.68	105.23	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	270	VAL	N-CA-C	-5.59	106.75	111.56
1	B	244	VAL	N-CA-C	5.57	118.14	108.90
1	B	232	VAL	CA-C-N	5.47	126.68	119.84
1	B	232	VAL	C-N-CA	5.47	126.68	119.84
1	B	266	GLU	N-CA-C	5.43	117.91	111.33
1	A	278	VAL	N-CA-C	5.32	118.07	110.09
1	A	232	VAL	CA-C-N	5.07	126.18	119.84
1	A	232	VAL	C-N-CA	5.07	126.18	119.84

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	2488	0	2506	70	0
1	B	2488	0	2506	105	0
2	A	44	0	26	2	0
2	B	44	0	26	2	0
3	A	134	0	0	4	0
3	B	123	0	0	6	1
All	All	5321	0	5064	173	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 17.

All (173) 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:154:LEU:HD22	1:A:172:MET:HE3	1.28	1.10
1:B:169:GLU:OE2	1:B:245:ARG:HD3	1.77	0.84
1:B:251:THR:H	1:B:254:GLN:NE2	1.82	0.78
1:A:289:THR:HG23	1:A:320:LYS:HE2	1.66	0.76
1:B:41:ALA:O	1:B:45:LYS:HG3	1.85	0.76
1:B:237:VAL:HG23	1:B:280:SER:HB2	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:LEU:HD13	1:B:84:TRP:CZ2	2.22	0.75
1:A:275:GLU:O	1:A:295:LYS:HD2	1.86	0.74
1:A:41:ALA:O	1:A:45:LYS:HG3	1.87	0.74
1:A:242:LEU:HD11	1:A:244:VAL:HG13	1.69	0.73
1:A:154:LEU:HD11	1:A:242:LEU:HD22	1.71	0.72
1:A:176:HIS:HA	1:A:238:SER:HB3	1.71	0.71
1:B:299:ALA:CB	1:B:305:VAL:HG12	2.21	0.70
1:B:236:ASN:O	1:B:237:VAL:HG22	1.91	0.70
1:B:172:MET:HG2	1:B:173:THR:N	2.07	0.69
1:A:182:GLN:HB3	3:A:2041:HOH:O	1.91	0.68
1:A:9:GLY:HA3	2:A:351:NAD:O5B	1.94	0.67
1:A:146:ASN:HD22	1:A:324:LEU:HD22	1.58	0.67
1:B:263:ALA:O	1:B:268:LYS:HG2	1.95	0.66
1:B:72:ARG:HD3	1:B:74:THR:HG23	1.78	0.65
1:B:41:ALA:HB2	1:B:64:LEU:CD2	2.25	0.65
1:B:298:ILE:HD12	1:B:299:ALA:H	1.62	0.65
1:A:168:ILE:HD11	1:A:247:GLU:HG3	1.78	0.64
1:A:172:MET:HG2	1:A:173:THR:N	2.12	0.64
1:B:122:SER:HB3	1:B:125:THR:HB	1.80	0.63
1:A:32:ASP:O	1:A:75:ALA:HA	1.97	0.63
1:A:155:ALA:HB3	1:A:156:PRO:HD3	1.80	0.63
1:B:92:VAL:HG13	1:B:116:VAL:HG22	1.81	0.63
1:B:251:THR:H	1:B:254:GLN:HE21	1.46	0.63
1:B:176:HIS:HA	1:B:238:SER:HB3	1.79	0.63
1:A:134:PHE:HB2	3:A:2021:HOH:O	1.99	0.62
1:B:32:ASP:O	1:B:75:ALA:HA	1.99	0.62
1:A:154:LEU:CD2	1:A:172:MET:HE3	2.18	0.62
1:B:237:VAL:CG2	1:B:280:SER:HB2	2.30	0.62
1:B:76:GLU:HG2	1:B:81:ASN:HB2	1.82	0.61
1:B:115:LYS:HG2	1:B:142:ASP:HA	1.80	0.61
1:B:182:GLN:OE1	1:B:231:ARG:HD2	1.99	0.61
1:A:10:ARG:O	1:A:14:ILE:HG12	2.01	0.60
1:A:106:ARG:O	1:A:110:THR:HG22	2.02	0.60
1:A:293:ASP:HB3	1:A:308:VAL:HG22	1.83	0.60
1:B:130:LYS:HG2	1:B:270:VAL:CG2	2.31	0.60
1:A:107:LYS:HA	1:A:110:THR:HG23	1.82	0.59
1:A:236:ASN:O	1:A:237:VAL:HG22	2.02	0.59
1:B:155:ALA:HB3	1:B:156:PRO:HD3	1.83	0.59
1:A:299:ALA:CB	1:A:305:VAL:HG12	2.32	0.59
1:A:251:THR:H	1:A:254:GLN:NE2	2.03	0.57
1:B:165:PHE:O	1:B:248:LYS:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:LYS:HA	1:A:110:THR:CG2	2.35	0.57
1:B:146:ASN:HD22	1:B:324:LEU:HD22	1.69	0.56
1:A:299:ALA:HB1	1:A:305:VAL:HG12	1.88	0.56
1:B:92:VAL:CG1	1:B:116:VAL:HG22	2.35	0.56
1:B:60:LYS:C	1:B:62:GLY:H	2.13	0.55
1:B:130:LYS:HG2	1:B:270:VAL:HG22	1.88	0.55
1:A:168:ILE:HG21	1:A:245:ARG:HH12	1.71	0.55
1:B:313:ASN:O	2:B:352:NAD:H4N	2.07	0.55
1:B:9:GLY:O	1:B:13:ARG:HB2	2.08	0.54
1:B:237:VAL:HG21	1:B:280:SER:O	2.07	0.54
1:B:58:GLU:CG	1:B:65:ILE:HB	2.38	0.53
1:B:208:THR:HG22	1:B:210:ALA:H	1.72	0.53
1:B:41:ALA:HB2	1:B:64:LEU:HD21	1.89	0.53
1:A:168:ILE:HD11	1:A:247:GLU:CG	2.39	0.53
1:B:92:VAL:HG11	1:B:108:HIS:HB3	1.90	0.53
1:B:204:ILE:HB	1:B:231:ARG:HB2	1.91	0.52
1:B:83:LYS:HB3	1:B:86:GLU:HG2	1.91	0.52
1:B:94:GLU:HB3	1:B:118:MET:HA	1.91	0.52
1:B:106:ARG:HA	1:B:143:ILE:HD11	1.93	0.51
1:A:251:THR:H	1:A:254:GLN:HE21	1.59	0.51
1:B:300:LEU:HD12	1:B:300:LEU:O	2.10	0.51
1:A:82:LEU:O	1:A:83:LYS:C	2.53	0.51
1:A:168:ILE:HG21	1:A:245:ARG:NH1	2.26	0.51
1:A:298:ILE:HG12	1:A:299:ALA:N	2.25	0.51
1:B:299:ALA:HA	1:B:305:VAL:HG12	1.92	0.51
1:A:100:LEU:HB2	1:A:122(A):LYS:HG3	1.92	0.51
1:B:250:ALA:HA	1:B:254:GLN:NE2	2.26	0.51
1:B:267:MET:HA	1:B:267:MET:HE2	1.93	0.50
1:A:121:PRO:HG3	1:A:148:SER:HB3	1.93	0.50
1:B:298:ILE:HD12	1:B:299:ALA:N	2.26	0.50
1:B:58:GLU:HG3	1:B:65:ILE:HB	1.93	0.50
1:B:211:ALA:HB3	3:B:2058:HOH:O	2.12	0.50
1:B:84:TRP:HB3	1:B:89:VAL:HB	1.94	0.50
1:A:219:PRO:O	1:A:222:ASN:HB2	2.12	0.49
1:A:150:THR:HG22	1:A:172:MET:HE1	1.93	0.49
1:A:45:LYS:HG2	1:A:57:VAL:HG11	1.93	0.49
1:B:329:SER:O	1:B:330:LYS:CB	2.59	0.49
1:A:25:ILE:CG2	1:A:26:GLU:N	2.76	0.49
1:B:45:LYS:HG2	1:B:57:VAL:HG11	1.95	0.49
1:B:118:MET:HG3	1:B:144:VAL:O	2.13	0.49
1:A:159:LYS:HB2	1:A:218:LEU:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:GLY:HA3	1:B:244:VAL:HG12	1.96	0.48
1:A:46:TYR:HB3	1:B:282:ASP:OD1	2.14	0.48
1:B:250:ALA:HA	1:B:254:GLN:HE21	1.77	0.48
1:B:44:LEU:O	1:B:53:PHE:HB2	2.13	0.48
1:A:95:ALA:O	2:A:351:NAD:H52N	2.14	0.48
1:B:60:LYS:O	1:B:62:GLY:N	2.47	0.48
1:B:299:ALA:CA	1:B:305:VAL:HG12	2.43	0.47
1:A:169:GLU:OE2	1:A:245:ARG:NH1	2.47	0.47
1:A:251:THR:HA	1:A:302:ASP:OD2	2.14	0.47
1:B:329:SER:O	1:B:330:LYS:HB2	2.13	0.47
1:B:252:TYR:OH	1:B:297:GLY:HA3	2.14	0.47
1:B:76:GLU:CG	1:B:81:ASN:HB2	2.42	0.47
1:B:9:GLY:HA3	2:B:352:NAD:O5B	2.15	0.47
1:A:8:PHE:CE1	1:A:30:ILE:HD13	2.50	0.47
1:A:154:LEU:HD23	1:A:214:VAL:HG21	1.97	0.47
1:A:215:GLY:HA3	1:A:222:ASN:HA	1.95	0.47
1:A:282:ASP:HB3	1:B:52:ARG:HH11	1.78	0.47
1:B:63:HIS:CD2	1:B:70:LYS:HZ1	2.33	0.47
1:B:183:LYS:NZ	3:B:2100:HOH:O	2.47	0.47
1:B:251:THR:N	1:B:254:GLN:HE21	2.12	0.47
1:B:60:LYS:C	1:B:62:GLY:N	2.74	0.46
1:A:87:VAL:HG12	1:A:89:VAL:HG23	1.97	0.46
1:B:63:HIS:CD2	1:B:70:LYS:NZ	2.83	0.46
1:B:169:GLU:HG3	1:B:245:ARG:HG3	1.97	0.46
1:B:17:ARG:NH2	1:B:51:GLY:O	2.43	0.46
1:A:30:ILE:HG22	1:A:40:MET:HE1	1.98	0.46
1:A:242:LEU:HD12	1:A:243:THR:N	2.31	0.46
1:B:72:ARG:HD2	3:B:2077:HOH:O	2.15	0.46
1:B:24:ASP:OD2	1:B:24:ASP:N	2.48	0.45
1:B:245:ARG:HA	1:B:303:ASN:O	2.16	0.45
1:B:37:ALA:O	1:B:64:LEU:HD21	2.16	0.45
1:A:95:ALA:HA	1:A:119:THR:OG1	2.16	0.45
1:B:2:LYS:O	1:B:90:ASP:HB2	2.16	0.45
1:A:237:VAL:HG21	1:A:280:SER:O	2.16	0.45
1:B:201:GLN:HG3	3:B:2084:HOH:O	2.17	0.45
1:B:266:GLU:H	1:B:266:GLU:CD	2.24	0.45
1:A:220:GLU:HG2	1:A:221:LEU:N	2.32	0.45
1:B:122:SER:CB	1:B:125:THR:HB	2.47	0.45
1:B:45:LYS:HG2	1:B:57:VAL:CG1	2.47	0.45
1:B:16:PHE:O	1:B:20:GLN:HG2	2.17	0.45
1:B:154:LEU:HD11	1:B:242:LEU:HD22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:VAL:CG2	1:A:280:SER:HB2	2.48	0.44
1:B:210:ALA:O	1:B:214:VAL:HG23	2.17	0.44
1:A:82:LEU:HD13	1:A:84:TRP:CZ2	2.53	0.44
1:A:135:ASP:OD2	1:A:136:LYS:HD3	2.18	0.44
1:A:169:GLU:CD	1:A:245:ARG:HH11	2.26	0.44
1:B:96:THR:OG1	1:B:98:LEU:HB2	2.17	0.44
1:A:129:VAL:HB	1:A:132:ALA:HB3	1.99	0.44
1:A:295:LYS:NZ	3:A:2150:HOH:O	2.51	0.44
1:A:37:ALA:O	1:A:64:LEU:HD21	2.18	0.44
1:A:287:VAL:HG23	1:A:319:ASN:ND2	2.34	0.43
1:B:289:THR:HG23	1:B:320:LYS:HD3	2.01	0.43
1:A:24:ASP:OD2	1:A:24:ASP:N	2.51	0.43
1:A:133:ASN:CG	1:A:217:VAL:HG22	2.44	0.43
1:B:317:TYR:O	1:B:321:VAL:HG23	2.18	0.43
1:A:211:ALA:HB3	3:A:2103:HOH:O	2.19	0.43
1:B:220:GLU:HG2	1:B:221:LEU:N	2.33	0.43
1:B:194:ARG:HH11	1:B:205:PRO:HG2	1.84	0.42
1:A:215:GLY:O	1:A:219:PRO:HA	2.19	0.42
1:B:130:LYS:HG2	1:B:270:VAL:HG21	1.99	0.42
1:B:295:LYS:HB2	3:B:2016:HOH:O	2.19	0.42
1:B:275:GLU:O	1:B:295:LYS:HE3	2.20	0.42
1:B:10:ARG:O	1:B:14:ILE:HG12	2.19	0.42
1:B:86:GLU:OE2	1:B:86:GLU:HA	2.19	0.42
1:B:251:THR:O	1:B:255:ILE:HG13	2.20	0.42
1:B:320:LYS:O	1:B:323:ASP:HB2	2.20	0.42
1:B:102:ASP:O	1:B:106:ARG:HB2	2.19	0.42
1:B:41:ALA:HB2	1:B:64:LEU:HD22	2.00	0.42
1:B:170:GLY:CA	1:B:244:VAL:HG12	2.49	0.42
1:B:239:VAL:HB	1:B:310:TRP:CE3	2.55	0.42
1:A:168:ILE:CD1	1:A:247:GLU:HG3	2.48	0.41
1:B:220:GLU:HG2	1:B:221:LEU:H	1.85	0.41
1:B:108:HIS:O	1:B:113:ALA:HB3	2.20	0.41
1:B:271:LEU:HD12	1:B:290:SER:O	2.21	0.41
1:A:248:LYS:O	1:A:248:LYS:HG2	2.19	0.41
1:B:78:ASP:C	1:B:78:ASP:OD2	2.61	0.41
1:B:87:VAL:HG23	1:B:89:VAL:HG23	2.02	0.41
1:A:237:VAL:HG12	1:A:312:ASP:HA	2.03	0.41
1:A:239:VAL:HG23	1:A:309:SER:O	2.21	0.41
1:A:176:HIS:HB3	1:A:231:ARG:HD3	2.03	0.41
1:B:260:LYS:O	1:B:264:GLU:HG3	2.21	0.41
1:B:280:SER:HB2	3:B:2108:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:GLY:O	1:B:268:LYS:HB2	2.21	0.40
1:A:315:THR:HG22	1:A:319:ASN:ND2	2.36	0.40
1:B:45:LYS:HB2	1:B:45:LYS:HE2	1.84	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)
3:B:2146:HOH:O	3:B:2146:HOH:O[2_455]	1.28	0.92

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	328/330 (99%)	300 (92%)	23 (7%)	5 (2%)	8	4
1	B	328/330 (99%)	305 (93%)	16 (5%)	7 (2%)	5	2
All	All	656/660 (99%)	605 (92%)	39 (6%)	12 (2%)	6	3

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	GLY
1	B	192	ASP
1	A	237	VAL
1	A	302	ASP
1	B	61	ASP
1	B	190	HIS
1	B	198	GLY
1	B	237	VAL
1	A	249	ALA
1	B	166	GLY

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Mol	Chain	Res	Type
1	B	186	ASP
1	A	198	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	265/265 (100%)	226 (85%)	39 (15%)	3	2
1	B	265/265 (100%)	228 (86%)	37 (14%)	3	2
All	All	530/530 (100%)	454 (86%)	76 (14%)	3	2

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	23	SER
1	A	34	LEU
1	A	45	LYS
1	A	57	VAL
1	A	58	GLU
1	A	64	LEU
1	A	70	LYS
1	A	83	LYS
1	A	87	VAL
1	A	94	GLU
1	A	103	GLU
1	A	110	THR
1	A	114	LYS
1	A	122	SER
1	A	122(A)	LYS
1	A	134	PHE
1	A	145	SER
1	A	172	MET
1	A	173	THR
1	A	200	SER

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Mol	Chain	Res	Type
1	A	204	ILE
1	A	212	LYS
1	A	217	VAL
1	A	228	MET
1	A	238	SER
1	A	240	VAL
1	A	244	VAL
1	A	245	ARG
1	A	247	GLU
1	A	254	GLN
1	A	264	GLU
1	A	266	GLU
1	A	271	LEU
1	A	298	ILE
1	A	308	VAL
1	A	311	TYR
1	A	315	THR
1	A	330	LYS
1	B	3	VAL
1	B	13	ARG
1	B	17	ARG
1	B	23	SER
1	B	34	LEU
1	B	43	MET
1	B	52	ARG
1	B	57	VAL
1	B	69	LYS
1	B	70	LYS
1	B	71	ILE
1	B	92	VAL
1	B	98	LEU
1	B	103	GLU
1	B	118	MET
1	B	122	SER
1	B	122(A)	LYS
1	B	134	PHE
1	B	149	CYS
1	B	161	ILE
1	B	173	THR
1	B	183	LYS
1	B	204	ILE
1	B	206	SER

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Mol	Chain	Res	Type
1	B	208	THR
1	B	212	LYS
1	B	221	LEU
1	B	238	SER
1	B	245	ARG
1	B	254	GLN
1	B	268	LYS
1	B	289	THR
1	B	298	ILE
1	B	300	LEU
1	B	302	ASP
1	B	320	LYS
1	B	330	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	ASN
1	A	201	GLN
1	A	254	GLN
1	A	319	ASN
1	B	63	HIS
1	B	124	ASN
1	B	146	ASN
1	B	164	ASN
1	B	176	HIS
1	B	201	GLN
1	B	222	ASN
1	B	236	ASN
1	B	254	GLN
1	B	284	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	B	352	-	46,48,48	2.35	9 (19%)	64,73,73	1.75	10 (15%)
2	NAD	A	351	-	46,48,48	2.47	10 (21%)	64,73,73	1.71	9 (14%)

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	B	352	-	-	12/30/62/62	0/5/5/5
2	NAD	A	351	-	-	11/30/62/62	0/5/5/5

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	352	NAD	C4N-C3N	7.93	1.51	1.39
2	A	351	NAD	C4N-C3N	7.93	1.51	1.39
2	A	351	NAD	C2N-C3N	7.44	1.50	1.39
2	B	352	NAD	C2N-C3N	6.71	1.49	1.39
2	A	351	NAD	C5N-C4N	6.35	1.49	1.38
2	B	352	NAD	C5N-C4N	6.32	1.49	1.38
2	A	351	NAD	C2N-N1N	6.17	1.41	1.35
2	B	352	NAD	C2N-N1N	6.06	1.41	1.35
2	A	351	NAD	PA-O3	-4.66	1.54	1.59
2	A	351	NAD	PN-O3	3.64	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	352	NAD	PA-O3	-3.64	1.55	1.59
2	B	352	NAD	C6N-C5N	-3.37	1.31	1.38
2	A	351	NAD	C6N-N1N	3.32	1.42	1.35
2	B	352	NAD	C6N-N1N	3.16	1.42	1.35
2	A	351	NAD	C6N-C5N	-2.82	1.32	1.38
2	B	352	NAD	PN-O3	2.69	1.62	1.59
2	A	351	NAD	C2A-N1A	2.65	1.38	1.33
2	B	352	NAD	C2A-N1A	2.38	1.38	1.33
2	A	351	NAD	C3N-C7N	-2.16	1.47	1.50

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	351	NAD	C5N-C4N-C3N	-7.78	112.71	120.36
2	B	352	NAD	C5N-C4N-C3N	-7.53	112.96	120.36
2	B	352	NAD	C3N-C7N-N7N	5.07	123.98	117.74
2	A	351	NAD	C6N-C5N-C4N	4.65	126.15	119.45
2	B	352	NAD	C6N-C5N-C4N	4.58	126.05	119.45
2	A	351	NAD	C3N-C7N-N7N	4.51	123.29	117.74
2	B	352	NAD	C6N-N1N-C2N	-4.11	118.38	121.88
2	A	351	NAD	C6N-N1N-C2N	-3.98	118.49	121.88
2	B	352	NAD	C4D-O4D-C1D	3.47	113.10	109.92
2	B	352	NAD	O4D-C4D-C5D	-3.13	99.32	109.33
2	A	351	NAD	O4D-C4D-C5D	-3.05	99.56	109.33
2	A	351	NAD	O4B-C1B-C2B	-2.71	100.83	106.62
2	A	351	NAD	C4D-O4D-C1D	2.51	112.22	109.92
2	B	352	NAD	O4B-C1B-C2B	-2.51	101.25	106.62
2	B	352	NAD	O7N-C7N-N7N	-2.43	119.11	122.62
2	B	352	NAD	O7N-C7N-C3N	-2.20	116.91	119.60
2	A	351	NAD	O7N-C7N-C3N	-2.13	117.00	119.60
2	B	352	NAD	O3-PN-O1N	-2.07	104.48	110.70
2	A	351	NAD	O7N-C7N-N7N	-2.01	119.71	122.62

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	351	NAD	C5D-O5D-PN-O3
2	A	351	NAD	C5D-O5D-PN-O2N
2	A	351	NAD	O4D-C1D-N1N-C2N
2	A	351	NAD	O4D-C1D-N1N-C6N
2	A	351	NAD	C2D-C1D-N1N-C2N

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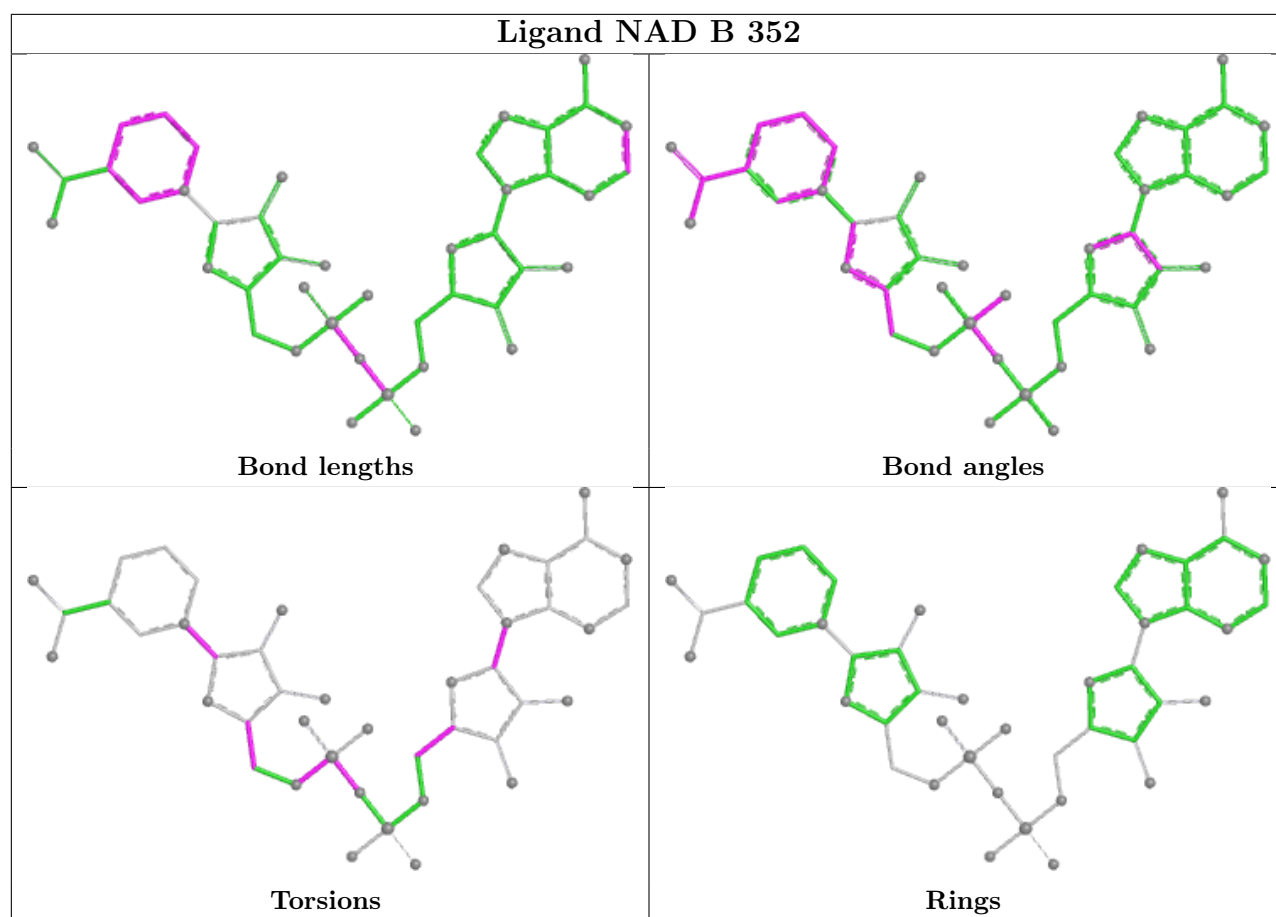
Mol	Chain	Res	Type	Atoms
2	A	351	NAD	C2D-C1D-N1N-C6N
2	B	352	NAD	C5D-O5D-PN-O3
2	B	352	NAD	C5D-O5D-PN-O1N
2	B	352	NAD	C5D-O5D-PN-O2N
2	B	352	NAD	O4D-C1D-N1N-C2N
2	B	352	NAD	O4D-C1D-N1N-C6N
2	B	352	NAD	C2D-C1D-N1N-C2N
2	B	352	NAD	C2D-C1D-N1N-C6N
2	B	352	NAD	C3D-C4D-C5D-O5D
2	B	352	NAD	O4D-C4D-C5D-O5D
2	A	351	NAD	O4D-C4D-C5D-O5D
2	A	351	NAD	C3D-C4D-C5D-O5D
2	A	351	NAD	C2B-C1B-N9A-C8A
2	B	352	NAD	C2B-C1B-N9A-C8A
2	A	351	NAD	C5D-O5D-PN-O1N
2	B	352	NAD	PA-O3-PN-O2N
2	A	351	NAD	O4B-C4B-C5B-O5B
2	B	352	NAD	O4B-C4B-C5B-O5B

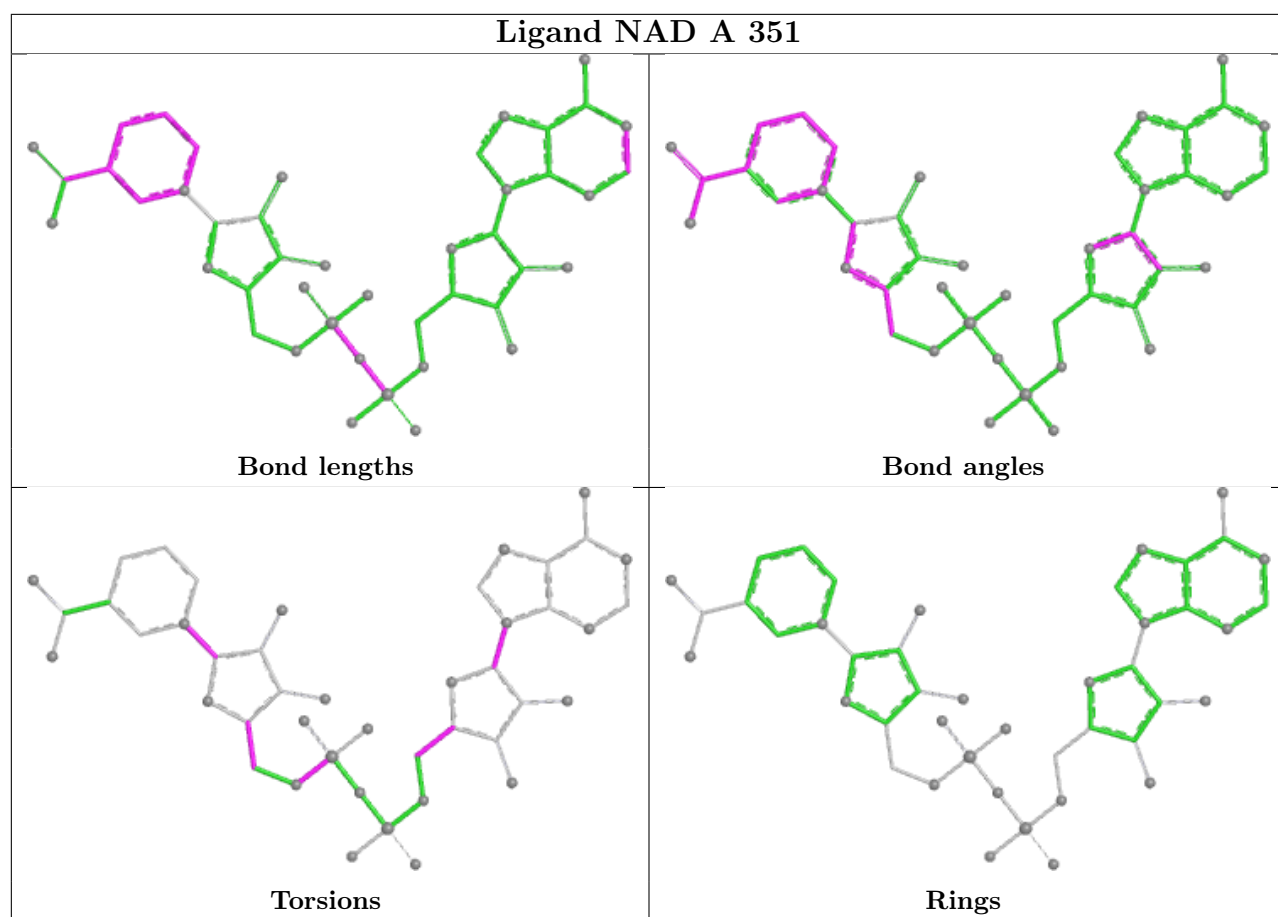
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	352	NAD	2	0
2	A	351	NAD	2	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	330/330 (100%)	-0.20	0 100 100	15, 26, 38, 52	0
1	B	330/330 (100%)	-0.16	1 (0%) 90 89	14, 26, 39, 60	0
All	All	660/660 (100%)	-0.18	1 (0%) 91 90	14, 26, 39, 60	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	237	VAL	2.4

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

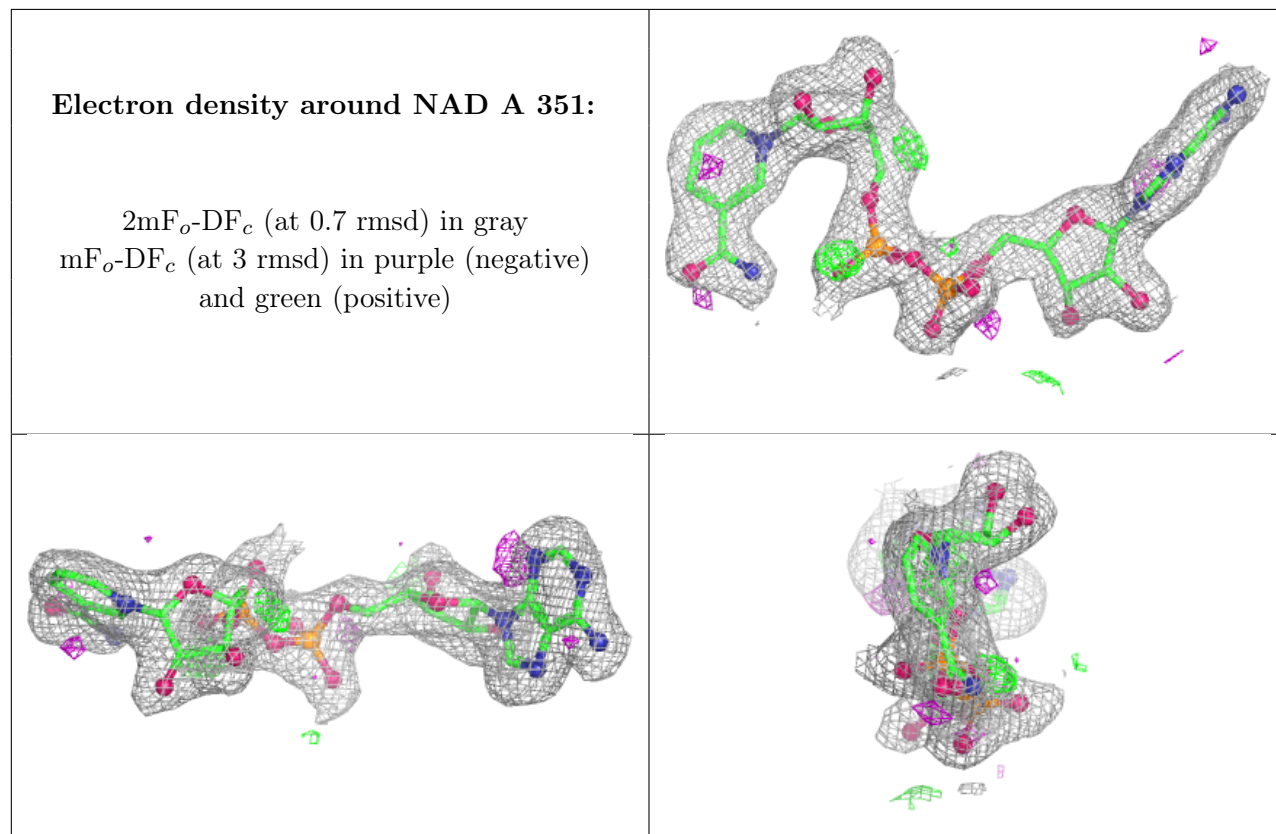
6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAD	A	351	44/44	0.94	0.08	12,29,34,37	0
2	NAD	B	352	44/44	0.95	0.07	14,25,31,33	0

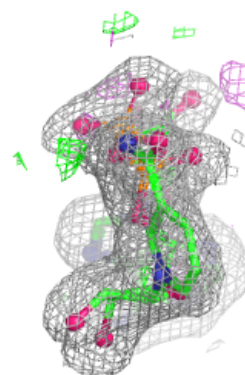
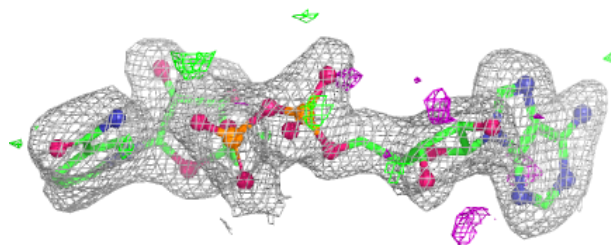
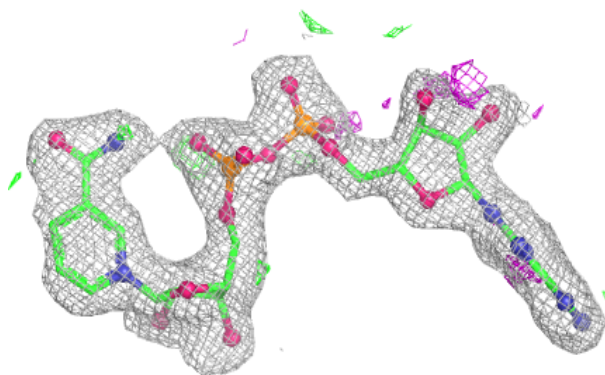
The following is a graphical depiction of the model fit to experimental electron density of all

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 B 352:

$2mF_o-DF_c$ (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.