



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:57 AM UTC

PDB ID : 2ZTM / pdb_00002ztm
Title : T190S mutant of D-3-hydroxybutyrate dehydrogenase
Authors : Nakashima, K.; Nakajima, Y.; Ito, K.; Yoshimoto, T.
Deposited on : 2008-10-07
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 ⓘ 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 ⓘ](#)) were used in the production of this report:

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

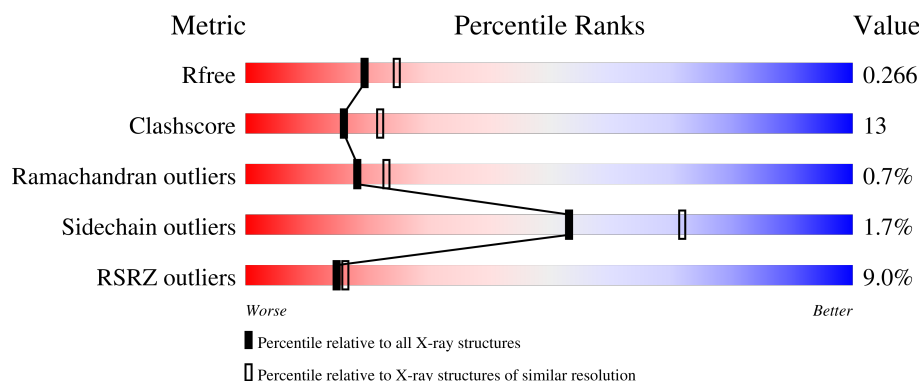
1 Overall quality at a glance

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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 $\leq 5\%$. 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	260	4% 79% 18% • •
1	B	260	20% 52% 32% • 14%
1	C	260	5% 78% 17% • •
1	D	260	5% 77% 17% • 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7551 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.

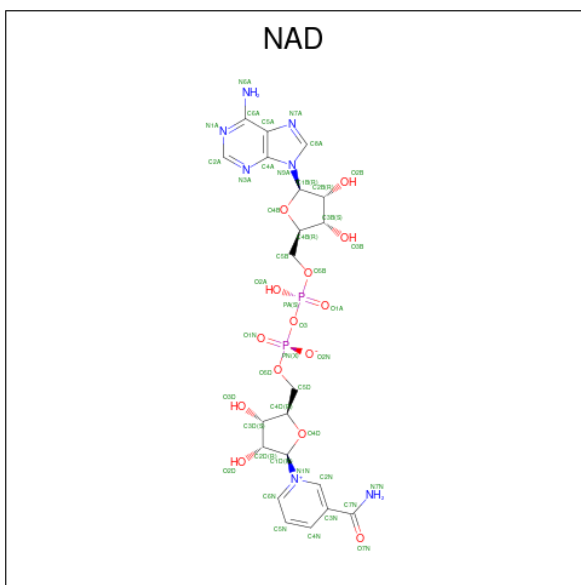
- Molecule 1 is a protein called D(-)-3-hydroxybutyrate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			1840	1155	327	355	3			
1	B	224	Total	C	N	O	S	0	0	0
			1598	1006	286	303	3			
1	C	250	Total	C	N	O	S	0	0	0
			1793	1131	317	341	4			
1	D	247	Total	C	N	O	S	0	0	0
			1771	1117	314	336	4			

There are 4 discrepancies between the modelled and reference sequences:

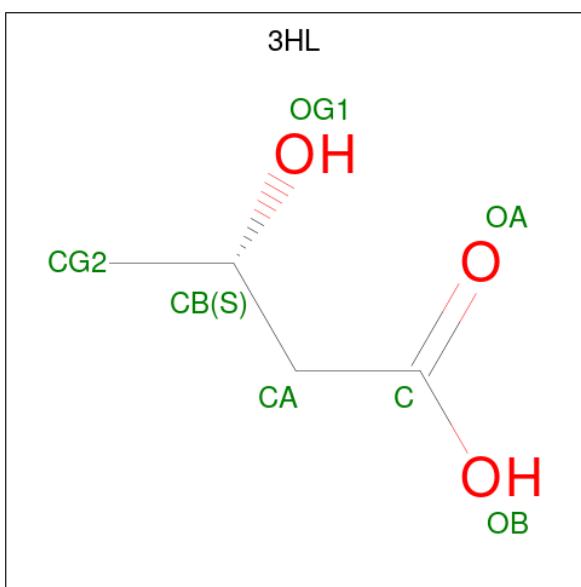
Chain	Residue	Modelled	Actual	Comment	Reference
A	190	SER	THR	engineered mutation	UNP Q5KST5
B	190	SER	THR	engineered mutation	UNP Q5KST5
C	190	SER	THR	engineered mutation	UNP Q5KST5
D	190	SER	THR	engineered mutation	UNP Q5KST5

- 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 44	C 21	N 7	O 14	P 2	0	0
2	C	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	D	1	Total 44	C 21	N 7	O 14	P 2	0	0

- Molecule 3 is (3S)-3-HYDROXYBUTANOIC ACID (CCD ID: 3HL) (formula: $C_4H_8O_3$).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mg 1	0	0
4	B	1	Total 1	Mg 1	0	0

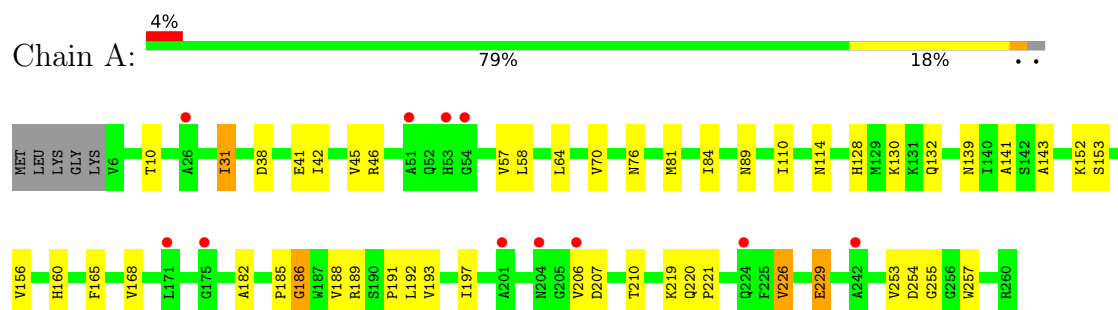
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	106	Total 106	O 106	0	0
5	B	54	Total 54	O 54	0	0
5	C	141	Total 141	O 141	0	0
5	D	107	Total 107	O 107	0	0

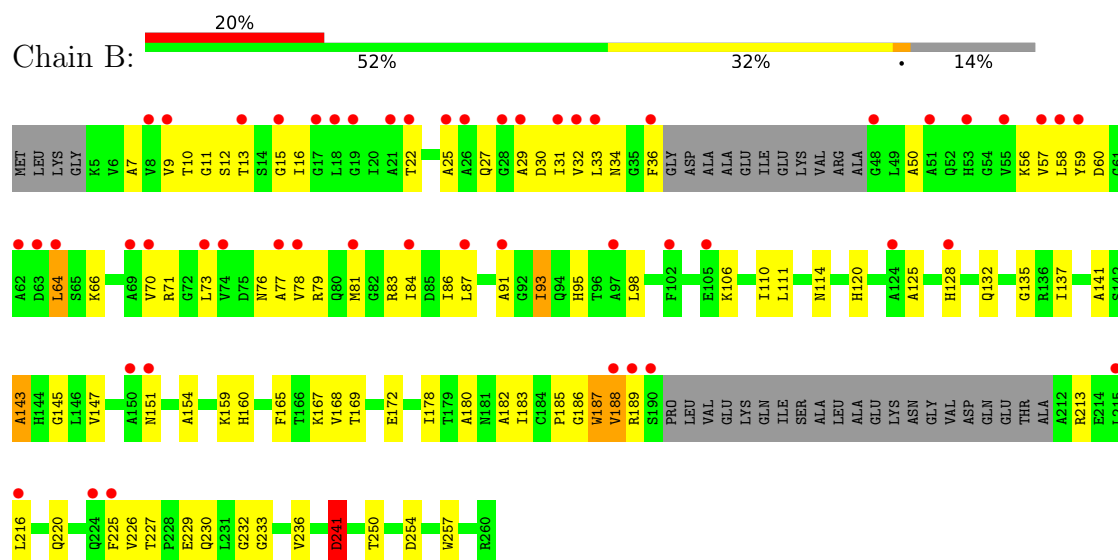
3 Residue-property plots [i](#)

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.

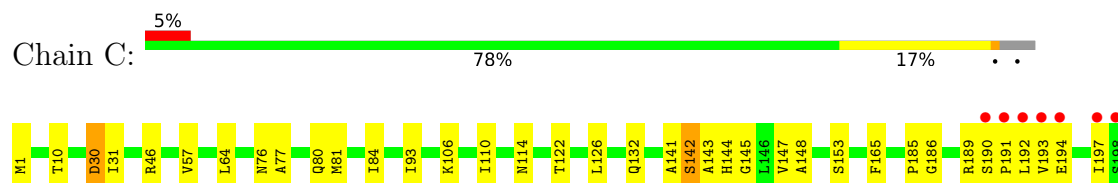
- Molecule 1: D(-)-3-hydroxybutyrate dehydrogenase

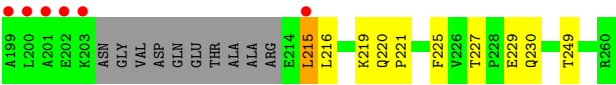


- Molecule 1: D(-)-3-hydroxybutyrate dehydrogenase

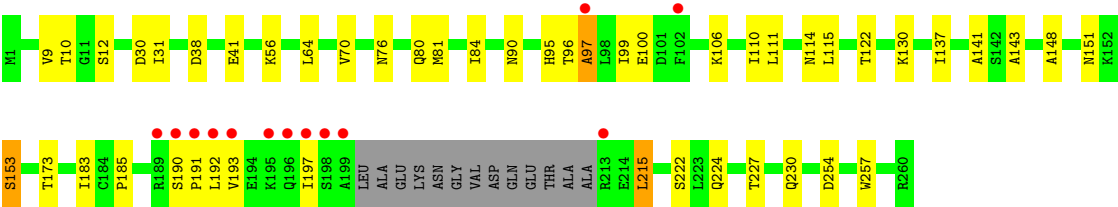
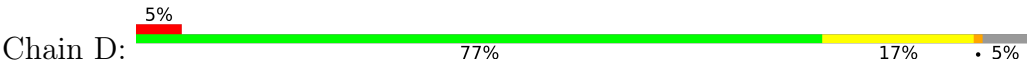


- Molecule 1: D(-)-3-hydroxybutyrate dehydrogenase





● Molecule 1: D(-)-3-hydroxybutyrate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.79Å 105.93Å 164.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.30) 99.7 (20.00-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.227 , 0.266 0.227 , 0.266	Depositor DCC
R_{free} test set	2950 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.1	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	7551	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% 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, 3HL, MG

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	0.39	0/1866	0.91	5/2534 (0.2%)
1	B	0.36	0/1621	0.92	5/2202 (0.2%)
1	C	0.45	0/1818	0.94	7/2467 (0.3%)
1	D	0.43	0/1796	0.92	3/2437 (0.1%)
All	All	0.41	0/7101	0.92	20/9640 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	GLY	N-CA-C	-7.58	101.76	112.60
1	B	143	ALA	N-CA-C	-7.31	103.46	112.38
1	C	186	GLY	N-CA-C	-7.23	102.26	112.60
1	A	143	ALA	N-CA-C	-7.22	103.41	111.71
1	C	31	ILE	N-CA-C	6.61	118.30	108.46
1	D	143	ALA	N-CA-C	-6.45	104.29	111.71
1	B	186	GLY	N-CA-C	-6.26	103.64	112.60
1	C	77	ALA	N-CA-C	-6.26	104.53	111.36
1	A	31	ILE	N-CA-C	6.11	117.28	108.48
1	A	226	VAL	N-CA-C	-5.69	103.85	110.05
1	C	145	GLY	N-CA-C	-5.69	105.71	115.61
1	A	160	HIS	N-CA-C	-5.44	105.43	111.36
1	B	160	HIS	N-CA-C	-5.38	105.42	111.28
1	D	215	LEU	N-CA-C	-5.35	105.49	112.23
1	D	31	ILE	N-CA-C	5.33	116.16	108.48
1	B	145	GLY	N-CA-C	-5.28	106.42	115.61
1	C	30	ASP	N-CA-C	-5.23	102.74	110.48
1	C	143	ALA	N-CA-C	-5.21	106.02	112.38
1	B	241	ASP	N-CA-C	-5.15	106.96	113.20
1	C	215	LEU	N-CA-C	-5.02	107.31	112.93

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	1840	0	1863	36	0
1	B	1598	0	1601	89	0
1	C	1793	0	1822	30	0
1	D	1771	0	1800	43	0
2	A	44	0	26	3	0
2	C	44	0	26	3	0
2	D	44	0	26	5	0
3	A	7	0	7	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	106	0	0	1	0
5	B	54	0	0	6	0
5	C	141	0	0	2	0
5	D	107	0	0	0	0
All	All	7551	0	7171	187	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 13.

All (187) 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:B:188:VAL:HG12	1:B:189:ARG:H	1.34	0.90
1:B:34:ASN:HD21	1:B:64:LEU:HD11	1.35	0.89
1:C:84:ILE:H	1:C:132:GLN:HE22	1.24	0.85
1:B:59:TYR:HD2	1:B:60:ASP:N	1.76	0.83
1:B:216:LEU:HD11	1:B:225:PHE:CE1	2.14	0.83
1:A:206:VAL:HG12	1:A:207:ASP:H	1.45	0.81
1:B:76:ASN:HA	1:B:79:ARG:HE	1.46	0.81
1:B:64:LEU:H	1:B:64:LEU:HD12	1.45	0.80
1:B:83:ARG:HA	5:B:1395:HOH:O	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:HIS:O	1:B:132:GLN:HG3	1.87	0.74
1:B:98:LEU:HD23	1:B:151:ASN:ND2	2.07	0.70
1:C:194:GLU:HA	1:C:197:ILE:HD12	1.74	0.69
1:C:10:THR:HB	1:C:64:LEU:HD11	1.75	0.69
1:B:29:ALA:O	1:B:31:ILE:HD12	1.94	0.68
1:B:78:VAL:HG21	1:B:128:HIS:CD2	2.28	0.68
1:B:232:GLY:O	1:B:236:VAL:HG23	1.93	0.68
1:B:56:LYS:HG2	1:B:58:LEU:HD11	1.74	0.68
1:B:188:VAL:HG12	1:B:189:ARG:N	2.08	0.67
1:B:9:VAL:HG12	1:B:12:SER:HB3	1.78	0.66
1:D:185:PRO:HG2	2:D:300:NAD:C5N	2.25	0.66
1:A:58:LEU:HD22	1:A:81:MET:HE3	1.77	0.65
1:D:56:LYS:HG2	1:D:81:MET:HE1	1.78	0.65
1:B:27:GLN:HG2	5:B:1360:HOH:O	1.96	0.65
1:A:152:LYS:O	1:A:156:VAL:HG23	1.98	0.64
1:B:93:ILE:HD13	1:B:93:ILE:H	1.62	0.64
1:A:206:VAL:HG12	1:A:207:ASP:N	2.13	0.64
1:D:197:ILE:HG12	1:D:215:LEU:HD13	1.80	0.63
1:D:64:LEU:HD22	1:D:70:VAL:HG22	1.81	0.63
1:A:206:VAL:HG11	1:A:210:THR:HG21	1.82	0.62
1:C:93:ILE:HG13	1:C:110:ILE:HD13	1.81	0.62
1:C:192:LEU:HD23	1:C:192:LEU:O	2.00	0.62
1:B:95:HIS:CD2	1:B:106:LYS:HG3	2.35	0.62
1:B:76:ASN:HA	1:B:79:ARG:NE	2.15	0.61
1:B:167:LYS:HE2	5:B:1405:HOH:O	1.99	0.61
1:C:193:VAL:O	1:C:197:ILE:HG13	2.01	0.61
1:B:16:ILE:HD11	1:B:188:VAL:HG21	1.82	0.60
1:B:110:ILE:O	1:B:114:ASN:HB2	2.02	0.60
1:A:165:PHE:CE1	1:C:153:SER:HB2	2.36	0.59
1:B:56:LYS:HG2	1:B:58:LEU:CD1	2.33	0.59
1:D:185:PRO:HG2	2:D:300:NAD:H5N	1.84	0.58
1:B:172:GLU:OE1	1:D:99:ILE:HG13	2.04	0.58
1:B:71:ARG:HE	1:B:120:HIS:HD2	1.51	0.58
1:A:130:LYS:HD2	5:C:1273:HOH:O	2.04	0.57
1:B:59:TYR:CD2	1:B:60:ASP:N	2.67	0.57
1:C:1:MET:HG2	5:C:1359:HOH:O	2.03	0.57
1:B:183:ILE:O	1:B:185:PRO:HD3	2.04	0.57
1:D:193:VAL:O	1:D:197:ILE:HG13	2.06	0.56
1:C:227:THR:OG1	1:C:230:GLN:HG3	2.05	0.55
1:A:76:ASN:HB3	5:A:1406:HOH:O	2.07	0.55
1:B:25:ALA:HA	1:B:31:ILE:HD11	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:THR:HB	1:C:64:LEU:CD1	2.37	0.54
1:D:10:THR:HB	1:D:64:LEU:CD1	2.38	0.54
1:B:77:ALA:O	1:B:81:MET:HB2	2.08	0.54
1:C:30:ASP:HB3	1:C:81:MET:HE3	1.90	0.53
1:B:9:VAL:CG1	1:B:12:SER:HB3	2.38	0.53
1:B:7:ALA:HA	1:B:86:ILE:O	2.09	0.53
1:B:71:ARG:HE	1:B:120:HIS:CD2	2.28	0.52
1:B:84:ILE:O	1:B:84:ILE:HG23	2.08	0.52
1:D:95:HIS:CD2	1:D:106:LYS:HG3	2.44	0.52
1:C:185:PRO:HB2	2:C:300:NAD:H5N	1.92	0.52
1:A:41:GLU:O	1:A:45:VAL:HG23	2.10	0.52
1:A:226:VAL:HG21	1:A:253:VAL:O	2.10	0.52
1:B:110:ILE:HD12	1:B:154:ALA:HB3	1.91	0.52
1:C:106:LYS:O	1:C:110:ILE:HG12	2.10	0.51
1:B:11:GLY:H	1:B:34:ASN:HD22	1.59	0.51
1:D:30:ASP:HB3	1:D:81:MET:HE3	1.93	0.51
1:B:233:GLY:O	1:B:236:VAL:HB	2.10	0.51
1:B:59:TYR:CD2	1:B:59:TYR:C	2.89	0.51
1:D:122:THR:HG23	1:D:137:ILE:HD13	1.93	0.50
1:A:64:LEU:HD22	1:A:70:VAL:HG22	1.92	0.50
1:B:213:ARG:HB2	1:B:213:ARG:NH1	2.26	0.50
1:B:167:LYS:NZ	5:B:1186:HOH:O	2.43	0.50
1:B:187:TRP:HA	1:B:187:TRP:CE3	2.46	0.50
1:B:241:ASP:HB2	5:B:1163:HOH:O	2.10	0.50
1:D:96:THR:O	1:D:97:ALA:HB2	2.12	0.50
1:B:64:LEU:HB3	1:B:70:VAL:HG22	1.94	0.50
1:D:215:LEU:HD23	1:D:215:LEU:O	2.12	0.49
1:D:222:SER:O	1:D:224:GLN:HG3	2.12	0.49
1:B:91:ALA:HB3	5:B:1397:HOH:O	2.12	0.49
1:B:168:VAL:HG22	1:D:148:ALA:HB1	1.95	0.49
1:C:215:LEU:HD12	1:C:215:LEU:O	2.12	0.49
1:A:46:ARG:HG2	1:A:46:ARG:HH11	1.78	0.48
1:A:110:ILE:O	1:A:114:ASN:HB2	2.12	0.48
1:C:76:ASN:ND2	1:C:80:GLN:OE1	2.46	0.48
1:A:38:ASP:O	1:A:42:ILE:HG13	2.13	0.48
1:B:64:LEU:C	1:B:66:LYS:H	2.21	0.48
1:D:76:ASN:O	1:D:80:GLN:HG2	2.12	0.48
1:B:73:LEU:O	1:B:73:LEU:HD12	2.13	0.48
1:D:84:ILE:HG23	1:D:84:ILE:O	2.14	0.48
1:C:84:ILE:O	1:C:84:ILE:HG23	2.13	0.48
1:C:189:ARG:HE	1:C:194:GLU:HG2	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:110:ILE:O	1:D:114:ASN:HB2	2.14	0.48
1:D:64:LEU:CD2	1:D:70:VAL:HG22	2.43	0.47
1:B:31:ILE:HD12	1:B:31:ILE:N	2.29	0.47
1:B:60:ASP:HB3	1:B:73:LEU:HD13	1.96	0.47
1:B:22:THR:O	1:B:25:ALA:HB3	2.13	0.47
1:B:56:LYS:O	1:B:81:MET:HE1	2.14	0.47
1:C:229:GLU:H	1:C:229:GLU:CD	2.22	0.47
1:B:59:TYR:HD2	1:B:59:TYR:C	2.23	0.47
1:B:93:ILE:HD13	1:B:93:ILE:N	2.30	0.47
1:B:254:ASP:HB2	1:B:257:TRP:HB2	1.96	0.46
1:A:84:ILE:HG23	1:A:84:ILE:O	2.15	0.46
1:A:188:VAL:HG22	1:A:226:VAL:HB	1.96	0.46
1:A:206:VAL:HG11	1:A:210:THR:CG2	2.44	0.46
1:A:191:PRO:HD2	2:A:300:NAD:O2A	2.16	0.46
1:B:188:VAL:O	1:B:225:PHE:HD2	1.99	0.46
1:B:30:ASP:HB3	1:B:81:MET:CE	2.46	0.46
1:B:58:LEU:HD12	1:B:58:LEU:N	2.31	0.46
1:B:227:THR:OG1	1:B:230:GLN:HG3	2.16	0.46
1:A:219:LYS:HB3	1:A:257:TRP:CE2	2.51	0.46
1:B:187:TRP:HA	1:B:187:TRP:HE3	1.81	0.46
1:B:165:PHE:HE1	1:D:153:SER:HB2	1.81	0.46
1:B:98:LEU:HD23	1:B:151:ASN:HD21	1.80	0.45
1:B:168:VAL:CG2	1:D:148:ALA:HB1	2.46	0.45
1:B:87:LEU:HD22	1:B:125:ALA:HB2	1.99	0.45
1:D:38:ASP:HB3	1:D:41:GLU:HB2	1.97	0.45
1:D:10:THR:HB	1:D:64:LEU:HD11	1.98	0.45
1:A:221:PRO:HD2	1:A:255:GLY:O	2.17	0.45
1:B:7:ALA:HB2	1:B:86:ILE:HB	1.98	0.45
1:C:46:ARG:HG3	1:C:57:VAL:HG12	1.98	0.45
1:D:190:SER:HB2	1:D:191:PRO:HD2	1.99	0.45
1:D:141:ALA:O	2:D:300:NAD:H6N	2.17	0.45
1:B:64:LEU:H	1:B:64:LEU:CD1	2.22	0.45
1:B:50:ALA:HB2	1:B:57:VAL:HB	1.99	0.45
1:C:122:THR:O	1:C:126:LEU:HG	2.17	0.45
1:A:31:ILE:O	1:A:57:VAL:HA	2.16	0.44
1:B:169:THR:HA	1:D:99:ILE:HD11	1.98	0.44
1:A:128:HIS:O	1:A:132:GLN:HG2	2.17	0.44
1:B:30:ASP:HB3	1:B:81:MET:HE3	1.98	0.44
1:B:172:GLU:CD	1:D:99:ILE:HG13	2.41	0.44
1:B:229:GLU:OE1	1:B:229:GLU:N	2.51	0.44
1:C:141:ALA:O	2:C:300:NAD:H6N	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:215:LEU:HD12	1:C:215:LEU:C	2.42	0.44
1:A:186:GLY:O	2:A:300:NAD:H4N	2.17	0.44
1:D:130:LYS:HE2	1:D:173:THR:HG22	1.99	0.44
1:D:111:LEU:HD23	1:D:111:LEU:O	2.18	0.44
1:B:59:TYR:HD2	1:B:60:ASP:H	1.61	0.44
1:B:183:ILE:C	1:B:185:PRO:HD3	2.43	0.43
1:B:135:GLY:HA3	1:B:178:ILE:HG23	2.00	0.43
1:D:215:LEU:HD23	1:D:215:LEU:C	2.44	0.43
1:B:25:ALA:CA	1:B:31:ILE:HD11	2.49	0.43
1:B:165:PHE:CE1	1:D:153:SER:HB2	2.53	0.43
1:C:216:LEU:HD21	1:C:225:PHE:CE2	2.53	0.43
1:B:137:ILE:O	1:B:180:ALA:HA	2.19	0.43
1:A:192:LEU:HD23	1:A:192:LEU:O	2.19	0.43
1:B:188:VAL:O	1:B:225:PHE:CD2	2.71	0.43
1:D:254:ASP:HB2	1:D:257:TRP:HB3	2.01	0.43
1:A:219:LYS:O	1:A:221:PRO:HD3	2.18	0.43
1:B:159:LYS:HD3	1:B:159:LYS:HA	1.83	0.43
1:D:9:VAL:HG12	1:D:12:SER:HB3	2.01	0.42
1:A:46:ARG:HG2	1:A:46:ARG:NH1	2.34	0.42
1:B:31:ILE:HD12	1:B:31:ILE:H	1.83	0.42
1:A:168:VAL:HG22	1:C:148:ALA:HB1	2.02	0.42
1:D:192:LEU:C	1:D:192:LEU:HD23	2.44	0.42
1:B:188:VAL:HG13	1:B:226:VAL:O	2.20	0.42
1:A:254:ASP:HB2	1:A:257:TRP:CB	2.49	0.42
1:B:143:ALA:O	1:B:147:VAL:HG22	2.19	0.42
1:C:142:SER:HB2	2:C:300:NAD:H6N	2.00	0.42
1:A:139:ASN:O	1:A:182:ALA:HA	2.20	0.42
1:B:34:ASN:ND2	1:B:64:LEU:HD11	2.18	0.42
1:B:111:LEU:HD23	1:D:111:LEU:HD21	2.02	0.42
1:D:96:THR:HA	1:D:151:ASN:O	2.20	0.42
1:A:229:GLU:CD	1:A:229:GLU:H	2.28	0.42
1:D:192:LEU:HD23	1:D:192:LEU:O	2.19	0.42
1:B:84:ILE:H	1:B:132:GLN:HE22	1.68	0.41
1:A:193:VAL:O	1:A:197:ILE:HG13	2.20	0.41
1:D:227:THR:OG1	1:D:230:GLN:HG3	2.20	0.41
1:C:219:LYS:O	1:C:221:PRO:HD3	2.21	0.41
1:D:111:LEU:HD21	1:D:115:LEU:HD23	2.02	0.41
1:D:183:ILE:HG22	1:D:185:PRO:HD3	2.03	0.41
1:A:89:ASN:HB2	1:A:139:ASN:HD22	1.86	0.41
1:B:182:ALA:HB3	1:B:250:THR:HG22	2.03	0.41
1:C:110:ILE:O	1:C:114:ASN:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:144:HIS:HA	1:C:147:VAL:O	2.21	0.41
1:B:7:ALA:CB	1:B:86:ILE:HB	2.51	0.41
1:B:13:THR:HG21	1:B:36:PHE:O	2.21	0.41
1:D:257:TRP:CD1	1:D:257:TRP:C	2.99	0.41
1:A:10:THR:HB	1:A:64:LEU:CD1	2.50	0.40
1:C:190:SER:HB2	1:C:191:PRO:CD	2.51	0.40
1:D:90:ASN:OD1	2:D:300:NAD:H4D	2.22	0.40
1:A:153:SER:HB2	1:C:165:PHE:CE1	2.56	0.40
1:B:32:VAL:HG22	1:B:81:MET:HE2	2.03	0.40
1:A:185:PRO:HB2	2:A:300:NAD:C5N	2.52	0.40
1:A:192:LEU:HD23	1:A:192:LEU:C	2.46	0.40
1:B:10:THR:HA	1:B:34:ASN:HB3	2.02	0.40
1:B:60:ASP:CG	1:B:73:LEU:HD13	2.46	0.40
1:D:185:PRO:O	2:D:300:NAD:H5N	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	253/260 (97%)	237 (94%)	14 (6%)	2 (1%)	16	20
1	B	218/260 (84%)	198 (91%)	17 (8%)	3 (1%)	9	9
1	C	246/260 (95%)	236 (96%)	9 (4%)	1 (0%)	30	38
1	D	243/260 (94%)	233 (96%)	9 (4%)	1 (0%)	30	38
All	All	960/1040 (92%)	904 (94%)	49 (5%)	7 (1%)	18	23

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	141	ALA

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Mol	Chain	Res	Type
1	D	97	ALA
1	A	141	ALA
1	B	15	GLY
1	C	142	SER
1	A	189	ARG
1	B	188	VAL

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	183/187 (98%)	181 (99%)	2 (1%)	65	81
1	B	156/187 (83%)	150 (96%)	6 (4%)	29	44
1	C	177/187 (95%)	175 (99%)	2 (1%)	65	81
1	D	175/187 (94%)	173 (99%)	2 (1%)	65	81
All	All	691/748 (92%)	679 (98%)	12 (2%)	53	72

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	220	GLN
1	A	229	GLU
1	B	33	LEU
1	B	64	LEU
1	B	93	ILE
1	B	187	TRP
1	B	220	GLN
1	B	241	ASP
1	C	220	GLN
1	C	249	THR
1	D	100	GLU
1	D	153	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	GLN
1	A	128	HIS
1	A	139	ASN
1	A	176	GLN
1	A	204	ASN
1	A	230	GLN
1	A	245	GLN
1	B	34	ASN
1	B	76	ASN
1	B	80	GLN
1	B	89	ASN
1	B	95	HIS
1	B	120	HIS
1	B	128	HIS
1	B	132	GLN
1	B	139	ASN
1	B	151	ASN
1	B	176	GLN
1	B	224	GLN
1	C	76	ASN
1	C	132	GLN
1	C	139	ASN
1	D	27	GLN
1	D	76	ASN
1	D	95	HIS
1	D	128	HIS
1	D	132	GLN
1	D	139	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAD	A	300	-	46,48,48	2.84	13 (28%)	64,73,73	1.81	10 (15%)
2	NAD	C	300	-	46,48,48	2.82	13 (28%)	64,73,73	1.93	10 (15%)
2	NAD	D	300	-	46,48,48	2.80	13 (28%)	64,73,73	1.90	9 (14%)
3	3HL	A	301	-	6,6,6	0.99	0	6,7,7	1.18	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	A	300	-	-	14/30/62/62	0/5/5/5
2	NAD	C	300	-	-	9/30/62/62	0/5/5/5
2	NAD	D	300	-	-	7/30/62/62	0/5/5/5
3	3HL	A	301	-	-	2/4/4/4	-

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	300	NAD	C2N-N1N	9.80	1.45	1.35
2	A	300	NAD	C2N-N1N	9.75	1.45	1.35
2	D	300	NAD	C2N-N1N	9.58	1.45	1.35
2	C	300	NAD	C4N-C3N	9.28	1.53	1.39
2	D	300	NAD	C4N-C3N	9.13	1.53	1.39
2	A	300	NAD	C4N-C3N	9.06	1.53	1.39
2	A	300	NAD	C5N-C4N	7.76	1.52	1.38
2	D	300	NAD	C5N-C4N	7.40	1.51	1.38
2	C	300	NAD	C5N-C4N	7.34	1.51	1.38
2	D	300	NAD	PN-O3	5.10	1.65	1.59
2	C	300	NAD	PN-O3	4.99	1.64	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300	NAD	PN-O3	4.90	1.64	1.59
2	A	300	NAD	C2A-N3A	3.65	1.40	1.33
2	D	300	NAD	C2A-N3A	3.63	1.40	1.33
2	D	300	NAD	C4A-N3A	3.57	1.41	1.34
2	C	300	NAD	C2A-N3A	3.53	1.40	1.33
2	A	300	NAD	C4A-N3A	3.49	1.40	1.34
2	A	300	NAD	C6N-N1N	3.41	1.43	1.35
2	C	300	NAD	C4A-N3A	3.40	1.40	1.34
2	D	300	NAD	C6N-N1N	3.40	1.43	1.35
2	C	300	NAD	C6N-N1N	3.34	1.42	1.35
2	A	300	NAD	C2N-C3N	-2.87	1.34	1.39
2	C	300	NAD	O4D-C1D	2.83	1.44	1.40
2	D	300	NAD	O4D-C1D	2.67	1.44	1.40
2	C	300	NAD	C2N-C3N	-2.62	1.34	1.39
2	C	300	NAD	C5A-N7A	-2.62	1.34	1.39
2	A	300	NAD	O4D-C1D	2.61	1.44	1.40
2	A	300	NAD	C5A-N7A	-2.53	1.34	1.39
2	D	300	NAD	C2N-C3N	-2.47	1.35	1.39
2	D	300	NAD	C5A-C6A	2.43	1.47	1.41
2	C	300	NAD	C3D-C4D	2.37	1.59	1.53
2	D	300	NAD	C5A-N7A	-2.37	1.34	1.39
2	A	300	NAD	C5A-C6A	2.30	1.47	1.41
2	A	300	NAD	O4D-C4D	2.28	1.50	1.45
2	D	300	NAD	O4D-C4D	2.19	1.49	1.45
2	C	300	NAD	C5A-C6A	2.17	1.47	1.41
2	C	300	NAD	O2B-C2B	-2.15	1.37	1.43
2	A	300	NAD	C3D-C4D	2.14	1.58	1.53
2	D	300	NAD	C3D-C4D	2.03	1.58	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300	NAD	C5N-C4N-C3N	-10.13	110.41	120.36
2	D	300	NAD	C5N-C4N-C3N	-9.79	110.74	120.36
2	A	300	NAD	C5N-C4N-C3N	-8.69	111.83	120.36
2	C	300	NAD	N3A-C2A-N1A	-4.41	121.90	128.58
2	A	300	NAD	N3A-C2A-N1A	-4.38	121.94	128.58
2	D	300	NAD	N3A-C2A-N1A	-4.35	122.00	128.58
2	C	300	NAD	C2A-N1A-C6A	4.22	125.67	118.73
2	D	300	NAD	C2A-N1A-C6A	4.22	125.66	118.73
2	A	300	NAD	C2A-N1A-C6A	4.15	125.54	118.73
2	C	300	NAD	C6N-N1N-C2N	-3.63	118.79	121.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	300	NAD	C6N-N1N-C2N	-3.21	119.14	121.88
2	A	300	NAD	C6N-N1N-C2N	-2.92	119.39	121.88
2	A	300	NAD	C1B-N9A-C8A	2.81	133.33	127.09
2	D	300	NAD	C1B-N9A-C8A	2.68	133.05	127.09
2	D	300	NAD	C4A-N9A-C8A	-2.65	102.95	105.74
2	A	300	NAD	C4A-N9A-C8A	-2.57	103.04	105.74
2	C	300	NAD	C4A-N9A-C8A	-2.53	103.08	105.74
2	C	300	NAD	C1B-N9A-C8A	2.53	132.71	127.09
2	C	300	NAD	C5A-N7A-C8A	2.43	107.27	103.45
2	D	300	NAD	C5A-N7A-C8A	2.39	107.20	103.45
2	C	300	NAD	C2N-C3N-C4N	2.35	120.99	118.26
2	A	300	NAD	C2N-C3N-C4N	2.34	120.98	118.26
2	D	300	NAD	C2N-C3N-C4N	2.34	120.98	118.26
2	A	300	NAD	C5A-N7A-C8A	2.34	107.12	103.45
2	A	300	NAD	O2B-C2B-C3B	2.31	119.21	111.82
2	A	300	NAD	C3N-C7N-N7N	2.20	120.45	117.74
2	C	300	NAD	N6A-C6A-N1A	2.09	123.03	118.38
2	C	300	NAD	O2B-C2B-C3B	2.07	118.47	111.82
2	D	300	NAD	O4B-C4B-C5B	-2.03	102.83	109.33

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	300	NAD	C5D-O5D-PN-O3
2	A	300	NAD	C5D-O5D-PN-O1N
2	A	300	NAD	C5D-O5D-PN-O2N
2	A	300	NAD	O4D-C1D-N1N-C2N
2	A	300	NAD	O4D-C1D-N1N-C6N
2	A	300	NAD	C2D-C1D-N1N-C6N
2	C	300	NAD	C5D-O5D-PN-O3
2	C	300	NAD	C5D-O5D-PN-O1N
2	C	300	NAD	C5D-O5D-PN-O2N
2	C	300	NAD	O4D-C1D-N1N-C6N
2	C	300	NAD	C2D-C1D-N1N-C6N
2	D	300	NAD	C5D-O5D-PN-O3
2	D	300	NAD	C5D-O5D-PN-O1N
2	D	300	NAD	C5D-O5D-PN-O2N
3	A	301	3HL	C-CA-CB-OG1
3	A	301	3HL	C-CA-CB-CG2
2	A	300	NAD	PN-O3-PA-O5B
2	A	300	NAD	PA-O3-PN-O1N

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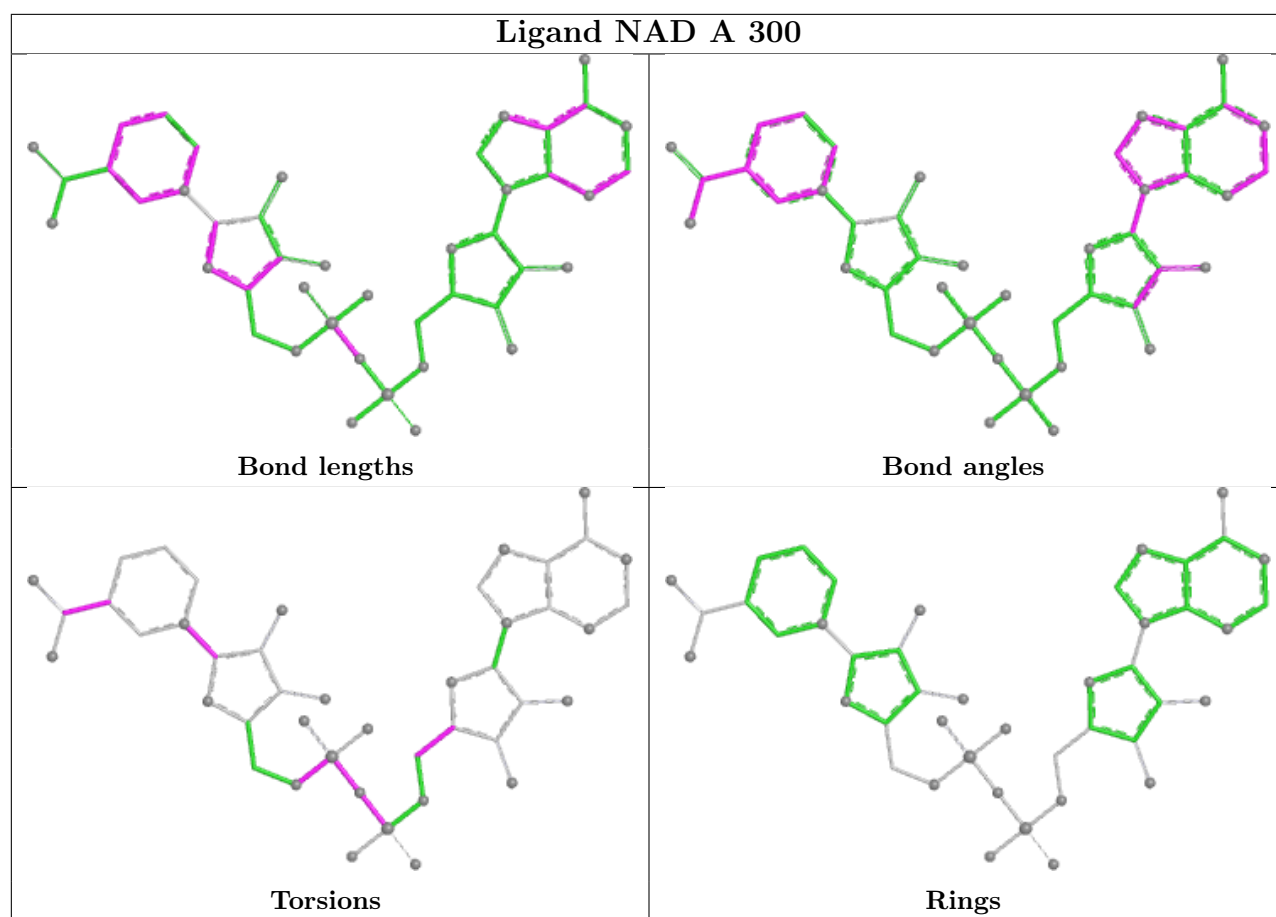
Mol	Chain	Res	Type	Atoms
2	C	300	NAD	PA-O3-PN-O2N
2	A	300	NAD	C4N-C3N-C7N-N7N
2	A	300	NAD	PA-O3-PN-O2N
2	D	300	NAD	PA-O3-PN-O2N
2	A	300	NAD	C4N-C3N-C7N-O7N
2	D	300	NAD	O4D-C1D-N1N-C6N
2	C	300	NAD	PN-O3-PA-O5B
2	A	300	NAD	C2N-C3N-C7N-N7N
2	C	300	NAD	PA-O3-PN-O1N
2	D	300	NAD	PA-O3-PN-O1N
2	C	300	NAD	O4B-C4B-C5B-O5B
2	A	300	NAD	C2N-C3N-C7N-O7N
2	D	300	NAD	O4B-C4B-C5B-O5B
2	A	300	NAD	O4B-C4B-C5B-O5B

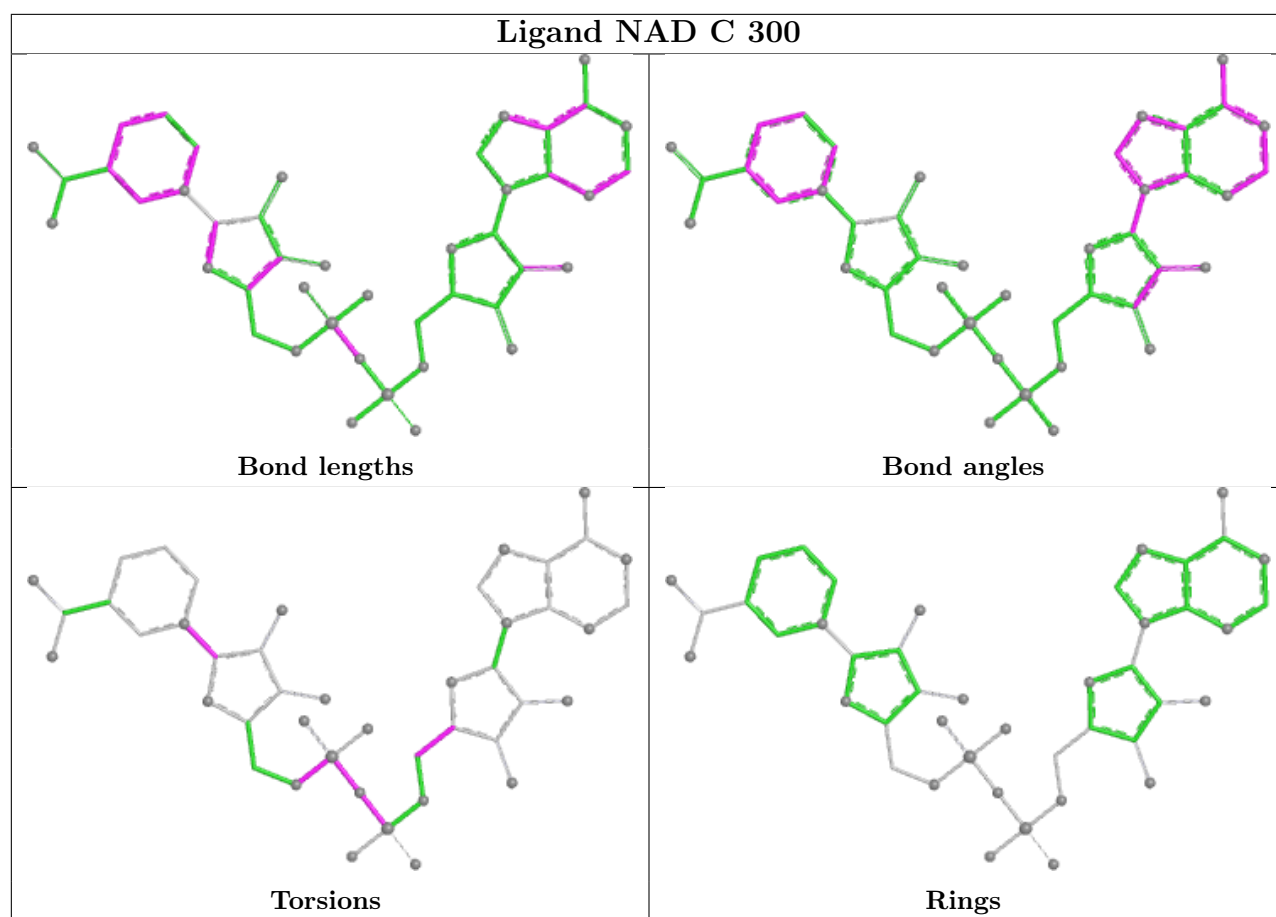
There are no ring outliers.

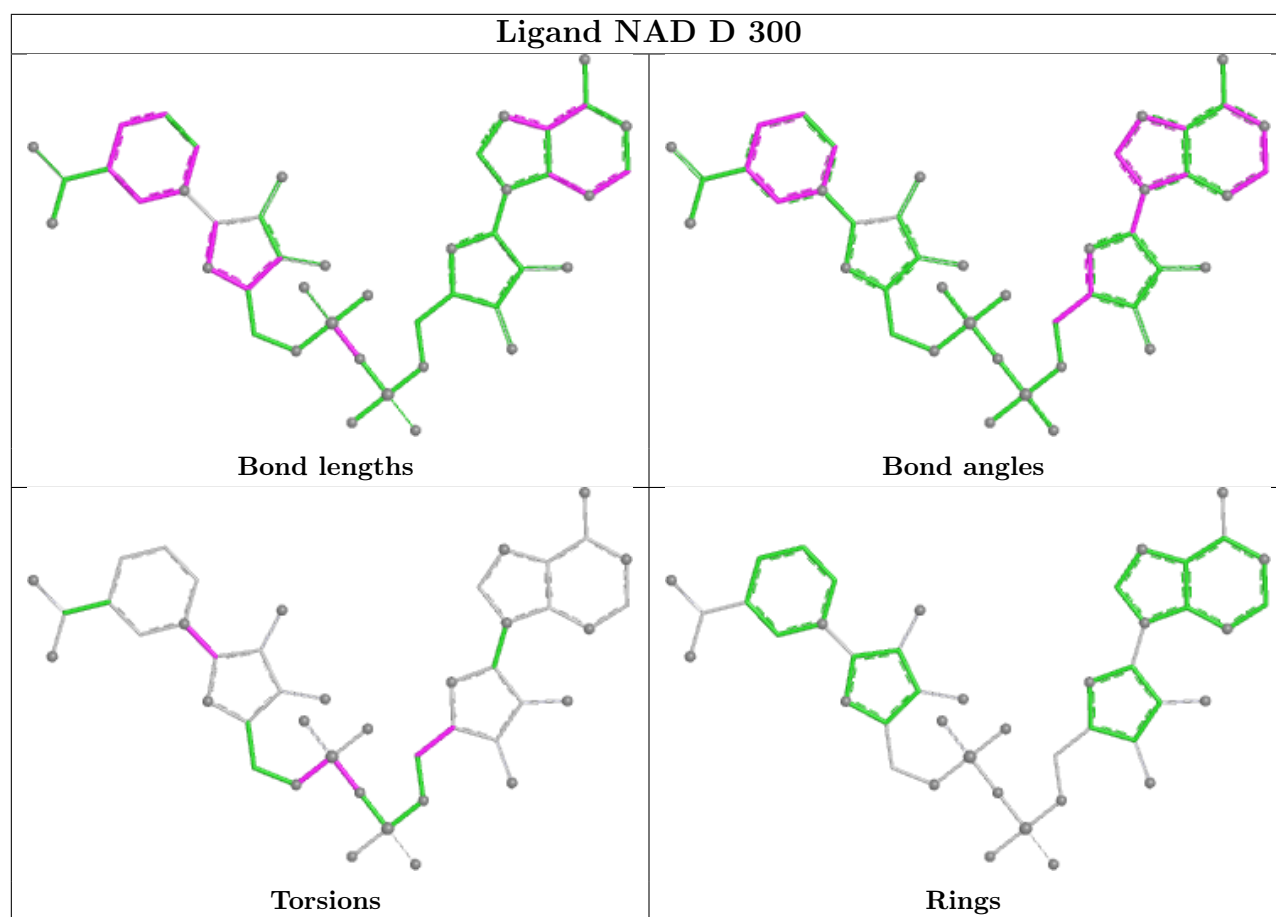
3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	NAD	3	0
2	C	300	NAD	3	0
2	D	300	NAD	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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	255/260 (98%)	0.38	11 (4%) 40 41	20, 40, 64, 78	0
1	B	224/260 (86%)	1.45	51 (22%) 2 2	32, 56, 79, 83	0
1	C	250/260 (96%)	-0.14	13 (5%) 33 34	19, 28, 64, 79	0
1	D	247/260 (95%)	0.14	13 (5%) 32 34	21, 34, 65, 85	1 (0%)
All	All	976/1040 (93%)	0.43	88 (9%) 15 16	19, 39, 74, 85	1 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	199	ALA	5.2
1	C	193	VAL	5.0
1	B	62	ALA	4.9
1	C	191	PRO	4.6
1	C	200	LEU	4.3
1	D	193	VAL	4.1
1	C	201	ALA	4.1
1	A	54	GLY	4.0
1	B	55	VAL	3.9
1	D	192	LEU	3.8
1	B	31	ILE	3.7
1	C	197	ILE	3.7
1	B	36	PHE	3.6
1	B	26	ALA	3.6
1	B	58	LEU	3.5
1	B	18	LEU	3.3
1	B	188	VAL	3.2
1	D	199	ALA	3.1
1	B	33	LEU	3.1
1	B	48	GLY	3.1
1	B	32	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	64	LEU	3.1
1	B	84	ILE	3.0
1	B	225	PHE	3.0
1	B	8	VAL	3.0
1	D	190	SER	3.0
1	D	191	PRO	3.0
1	B	53	HIS	2.9
1	C	202	GLU	2.9
1	B	70	VAL	2.9
1	C	192	LEU	2.8
1	D	189	ARG	2.7
1	B	57	VAL	2.7
1	B	13	THR	2.7
1	C	194	GLU	2.7
1	B	25	ALA	2.6
1	B	91	ALA	2.6
1	B	74	VAL	2.6
1	C	198	SER	2.6
1	D	195	LYS	2.6
1	B	21	ALA	2.6
1	B	9	VAL	2.6
1	B	189	ARG	2.6
1	B	63	ASP	2.6
1	A	224	GLN	2.5
1	B	73	LEU	2.4
1	D	196	GLN	2.4
1	B	97	ALA	2.4
1	C	190	SER	2.4
1	B	102	PHE	2.4
1	A	53	HIS	2.4
1	B	15	GLY	2.4
1	B	69	ALA	2.4
1	C	203	LYS	2.3
1	B	216	LEU	2.3
1	A	51	ALA	2.3
1	B	78	VAL	2.3
1	A	171	LEU	2.3
1	A	175	GLY	2.3
1	D	213	ARG	2.3
1	B	128	HIS	2.2
1	B	190	SER	2.2
1	D	102	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	224	GLN	2.2
1	D	197	ILE	2.2
1	A	206	VAL	2.2
1	A	204	ASN	2.2
1	B	22	THR	2.2
1	B	51	ALA	2.2
1	B	29	ALA	2.2
1	B	124	ALA	2.2
1	D	97	ALA	2.2
1	D	198	SER	2.2
1	B	215	LEU	2.1
1	C	215	LEU	2.1
1	B	151	ASN	2.1
1	B	28	GLY	2.1
1	B	87	LEU	2.1
1	B	59	TYR	2.1
1	A	201	ALA	2.1
1	B	17	GLY	2.1
1	A	26	ALA	2.0
1	A	242	ALA	2.0
1	B	77	ALA	2.0
1	B	81	MET	2.0
1	B	105	GLU	2.0
1	B	150	ALA	2.0
1	B	19	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

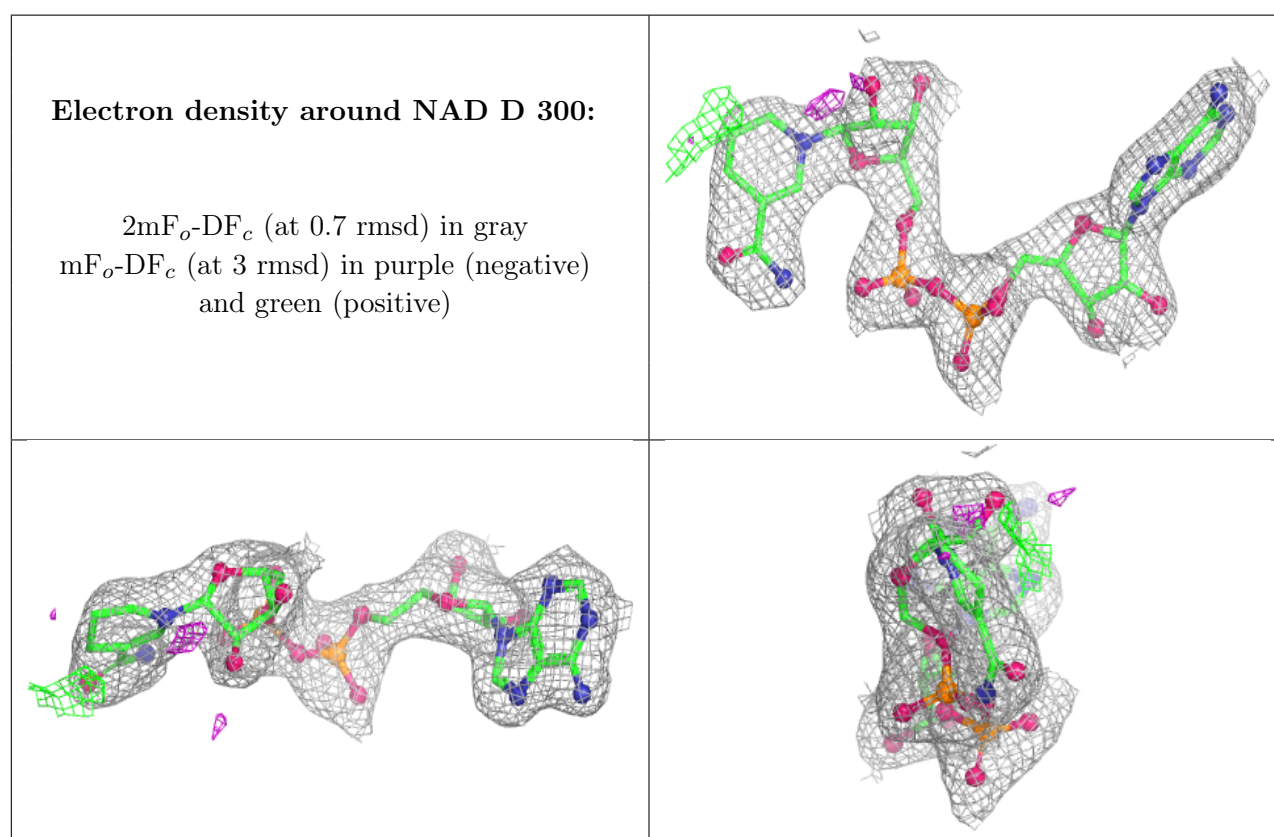
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

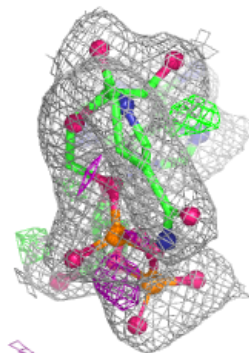
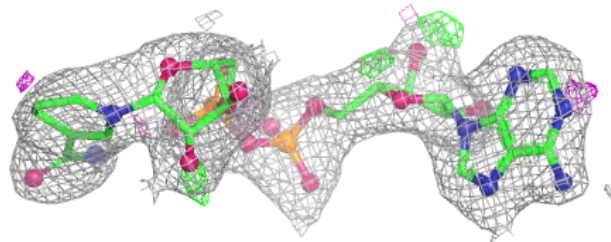
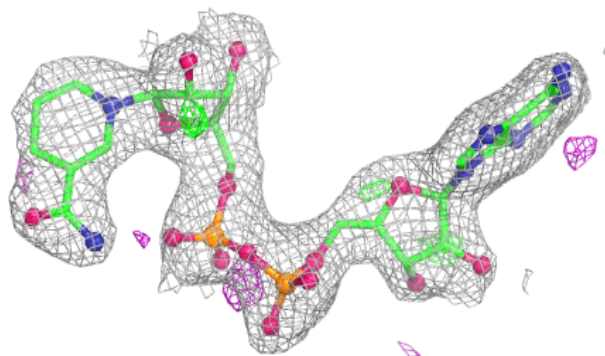
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	406	1/1	0.87	0.12	42,42,42,42	0
3	3HL	A	301	7/7	0.92	0.12	45,47,48,48	0
2	NAD	D	300	44/44	0.92	0.09	39,46,51,52	0
4	MG	B	401	1/1	0.92	0.12	42,42,42,42	0
2	NAD	A	300	44/44	0.93	0.09	34,40,43,44	0
2	NAD	C	300	44/44	0.95	0.07	31,38,44,47	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.

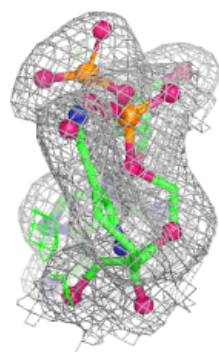
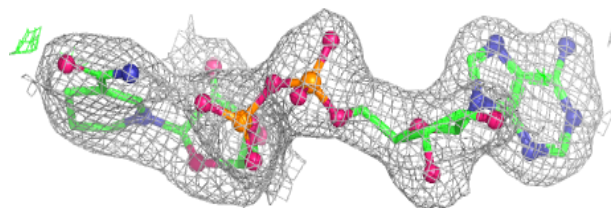
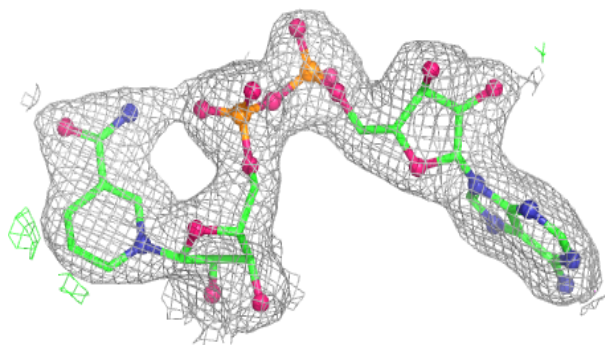


Electron density around NAD A 300:

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

**Electron density around NAD C 300:**

$2mF_o-DF_c$ (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.