



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 01:45 pm BST

PDB ID : 1ZXB
Title : Synthesis, Biological Activity, and X-Ray Crystal Structural Analysis of Diaryl Ether Inhibitors of Malarial Enoyl ACP Reductase. Part 1:4'-Substituted Triclosan Derivatives
Authors : Freundlich, J.S.; Anderson, J.W.; Sarantakis, D.; Shieh, H.M.; Yu, M.; Lucumi, E.; Kuo, M.; Schiehser, G.A.; Jacobus, D.P.; Jacobs Jr., W.R.; Fidock, D.A.; Sacchettini, J.C.
Deposited on : 2005-06-07
Resolution : 2.68 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 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)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

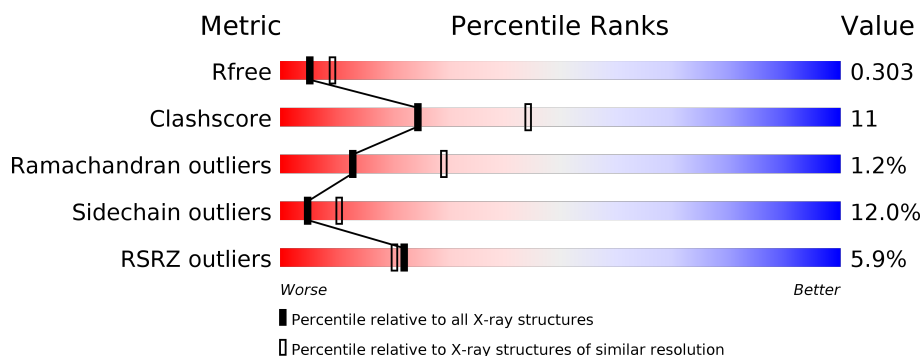
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	3863 (2.70-2.66)
Clashscore	141614	4210 (2.70-2.66)
Ramachandran outliers	138981	4141 (2.70-2.66)
Sidechain outliers	138945	4141 (2.70-2.66)
RSRZ outliers	127900	3780 (2.70-2.66)

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 ≥ 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	336	
1	B	336	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4702 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 enoyl-acyl carrier reductase.

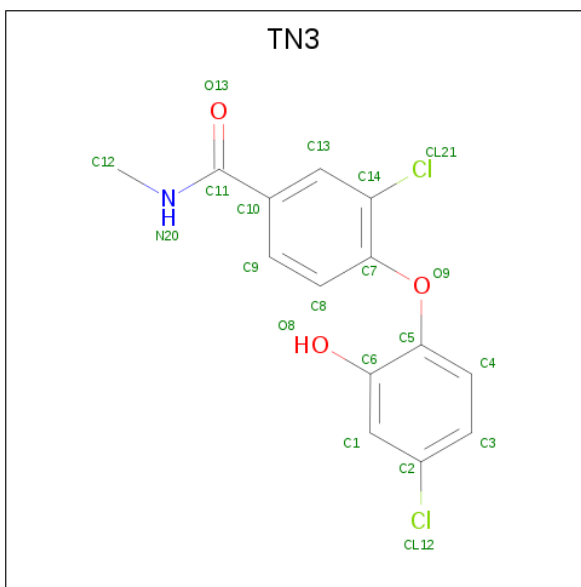
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2287	1458	384	434	11			
1	B	289	Total	C	N	O	S	0	0	0
			2287	1458	384	434	11			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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 3-CHLORO-4-(4-CHLORO-2-HYDROXYPHENOXY)-N-METHYLBENZAMIDE (three-letter code: TN3) (formula: C₁₄H₁₁Cl₂NO₃).

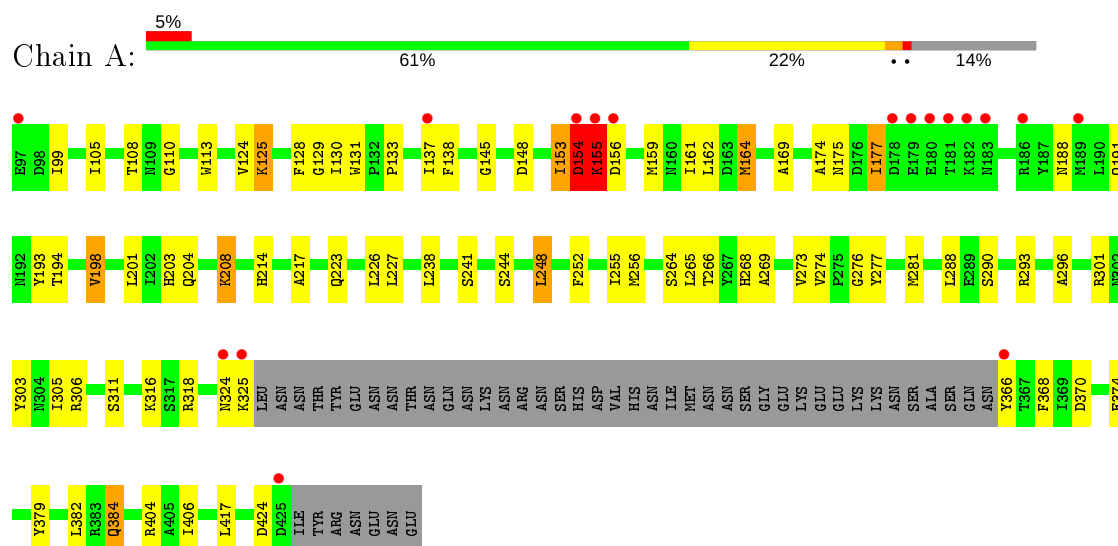


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			20	14	2	1	3		
3	B	1	Total	C	Cl	N	O	0	0
			20	14	2	1	3		

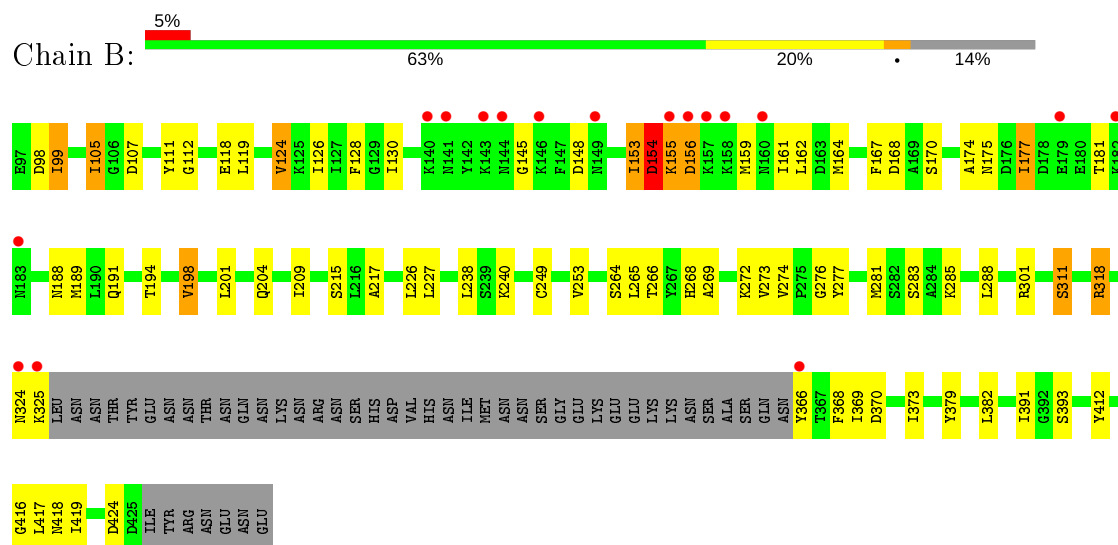
3 Residue-property plots [i](#)

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 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 ($\text{RSRZ} > 2$). 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: enoyl-acyl carrier reductase



- Molecule 1: enoyl-acyl carrier reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	131.43 Å 131.43 Å 82.95 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.68 29.75 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-2.68) 99.9 (29.75-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 2.68 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.233 , 0.301 0.234 , 0.303	Depositor DCC
R_{free} test set	1068 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4702	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 4.52% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAD, TN3

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	0.54	0/2329	0.65	0/3141
1	B	0.55	0/2329	0.67	0/3141
All	All	0.55	0/4658	0.66	0/6282

There are no bond length outliers.

There are no bond angle outliers.

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	2287	0	2297	53	0
1	B	2287	0	2297	52	0
2	A	44	0	26	2	0
2	B	44	0	26	0	0
3	A	20	0	10	1	0
3	B	20	0	11	1	0
All	All	4702	0	4667	102	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 11.

All (102) 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:170:SER:HB3	1:B:240:LYS:HE2	1.46	0.97
1:B:281:MET:HE1	1:B:285:LYS:HE2	1.63	0.80
1:A:223:GLN:NE2	1:A:324:ASN:HB3	1.99	0.78
1:A:203:HIS:HE1	1:A:208:LYS:NZ	1.93	0.67
1:B:119:LEU:O	1:B:124:VAL:HG13	1.95	0.67
1:A:276:GLY:HA2	1:A:325:LYS:HE3	1.77	0.67
1:B:170:SER:CB	1:B:240:LYS:HE2	2.23	0.66
1:A:203:HIS:HB2	1:A:255:ILE:HG21	1.78	0.66
1:B:274:VAL:O	1:B:274:VAL:HG23	1.96	0.66
1:A:203:HIS:HE1	1:A:208:LYS:HZ1	1.42	0.65
1:B:111:TYR:CE1	1:B:391:ILE:HD13	2.35	0.61
1:A:277:TYR:CE2	1:A:281:MET:HB3	2.36	0.61
1:B:168:ASP:OD1	1:B:170:SER:OG	2.18	0.60
1:A:161:ILE:HD13	1:A:164:MET:HE1	1.83	0.60
1:B:105:ILE:HG22	1:B:112:GLY:HA3	1.84	0.60
1:A:223:GLN:HE21	1:A:324:ASN:HB3	1.67	0.59
1:B:276:GLY:HA2	1:B:325:LYS:HE3	1.83	0.59
1:B:118:GLU:HG3	1:B:393:SER:N	2.18	0.59
1:B:154:ASP:OD1	1:B:154:ASP:N	2.34	0.59
1:A:301:ARG:NH2	1:B:379:TYR:HA	2.18	0.58
1:B:128:PHE:HB3	1:B:130:ILE:HD11	1.85	0.58
1:A:269:ALA:HB2	1:A:274:VAL:CG2	2.35	0.56
1:A:203:HIS:CE1	1:A:208:LYS:NZ	2.73	0.56
1:A:269:ALA:CB	1:A:274:VAL:HG22	2.35	0.56
1:A:194:THR:O	1:A:198:VAL:HG13	2.06	0.56
1:B:194:THR:O	1:B:198:VAL:HG13	2.06	0.56
1:B:274:VAL:HG23	1:B:277:TYR:HB2	1.89	0.55
1:B:153:ILE:HD11	1:B:159:MET:HA	1.89	0.54
1:B:366:TYR:O	1:B:370:ASP:HB2	2.07	0.54
1:A:274:VAL:HG23	1:A:277:TYR:CB	2.38	0.54
1:A:325:LYS:HB3	1:A:368:PHE:HB2	1.90	0.53
1:A:125:LYS:H	1:A:125:LYS:HD2	1.74	0.52
1:A:161:ILE:HD13	1:A:164:MET:CE	2.40	0.52
1:A:269:ALA:HB2	1:A:274:VAL:HG21	1.92	0.51
1:B:111:TYR:HE1	1:B:391:ILE:HD13	1.75	0.51
1:B:416:GLY:O	1:B:419:ILE:HG12	2.11	0.51
1:A:290:SER:HA	1:A:293:ARG:NH2	2.26	0.51
1:A:379:TYR:HA	1:B:301:ARG:NH2	2.25	0.50
1:B:281:MET:CE	1:B:285:LYS:HE2	2.38	0.50
1:A:269:ALA:CB	1:A:274:VAL:CG2	2.90	0.50
1:A:153:ILE:O	1:A:154:ASP:C	2.49	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:ASP:O	1:B:156:ASP:N	2.38	0.50
1:B:177:ILE:HG22	1:B:177:ILE:O	2.12	0.49
1:A:301:ARG:HH21	1:B:379:TYR:C	2.15	0.49
1:A:290:SER:HA	1:A:293:ARG:HH21	1.78	0.49
1:A:293:ARG:O	1:A:296:ALA:HB3	2.12	0.49
1:B:266:THR:HA	1:B:285:LYS:HD2	1.94	0.49
1:B:119:LEU:HB2	1:B:126:ILE:HD11	1.95	0.48
1:A:217:ALA:HB3	3:A:500:TN3:CL21	2.50	0.48
1:B:145:GLY:CA	1:B:148:ASP:OD2	2.61	0.48
1:A:145:GLY:HA2	1:A:148:ASP:OD2	2.14	0.48
1:B:269:ALA:CB	1:B:274:VAL:HG22	2.44	0.48
1:A:274:VAL:HG23	1:A:277:TYR:HB3	1.95	0.48
2:A:450:NAD:H2N	2:A:450:NAD:O2N	2.14	0.47
1:B:269:ALA:HA	1:B:272:LYS:O	2.13	0.47
1:B:217:ALA:HB3	3:B:501:TN3:CL21	2.51	0.47
1:B:177:ILE:CG2	1:B:177:ILE:O	2.61	0.47
1:A:129:GLY:C	1:A:130:ILE:HD13	2.34	0.47
1:B:369:ILE:O	1:B:373:ILE:HG13	2.14	0.47
1:A:105:ILE:HG13	1:A:130:ILE:HD12	1.97	0.47
1:A:252:PHE:O	1:A:256:MET:HG3	2.15	0.46
1:B:174:ALA:O	1:B:177:ILE:HB	2.16	0.46
1:B:269:ALA:HB1	1:B:274:VAL:HG22	1.96	0.46
1:B:161:ILE:HD13	1:B:164:MET:CE	2.46	0.45
1:B:161:ILE:HD13	1:B:164:MET:HE3	1.99	0.45
1:A:131:TRP:CZ2	1:A:133:PRO:HG2	2.53	0.44
1:B:288:LEU:O	1:B:288:LEU:HD23	2.17	0.44
1:A:177:ILE:HD11	1:A:193:TYR:CE2	2.52	0.44
1:B:215:SER:HA	1:B:265:LEU:HD23	2.00	0.44
1:B:264:SER:OG	1:B:288:LEU:HD21	2.18	0.44
1:A:384:GLN:HB3	1:A:384:GLN:HE21	1.56	0.44
1:B:99:ILE:HB	1:B:209:ILE:HG22	1.99	0.44
1:A:203:HIS:CE1	1:A:208:LYS:HZ2	2.36	0.44
1:B:418:ASN:H	1:B:418:ASN:HD22	1.66	0.43
1:A:266:THR:O	1:A:311:SER:HA	2.18	0.43
1:A:108:THR:HG22	1:A:113:TRP:CE2	2.53	0.43
1:A:153:ILE:HD11	1:A:159:MET:HA	2.00	0.43
1:A:174:ALA:O	1:A:177:ILE:HB	2.19	0.42
1:A:306:ARG:HD2	1:A:406:ILE:O	2.20	0.42
1:B:274:VAL:O	1:B:274:VAL:CG2	2.65	0.42
1:A:155:LYS:HD3	1:A:155:LYS:O	2.20	0.42
1:A:248:LEU:HD22	1:A:252:PHE:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:GLY:HA3	2:A:450:NAD:O2A	2.18	0.42
1:B:318:ARG:HB3	1:B:318:ARG:HH11	1.85	0.42
1:B:325:LYS:HB3	1:B:368:PHE:HB2	2.02	0.42
1:B:98:ASP:HB3	1:B:124:VAL:HB	2.01	0.42
1:A:128:PHE:HZ	1:A:159:MET:HE1	1.85	0.41
1:A:379:TYR:C	1:B:301:ARG:HH21	2.17	0.41
1:A:169:ALA:HB1	1:A:244:SER:HB2	2.02	0.41
1:A:223:GLN:HE22	1:A:324:ASN:HB3	1.80	0.41
1:A:138:PHE:HE2	1:A:164:MET:HE2	1.86	0.41
1:A:303:TYR:HB3	1:A:305:ILE:HD12	2.02	0.41
1:B:311:SER:HB3	1:B:412:TYR:CD1	2.56	0.41
1:A:214:HIS:O	1:A:264:SER:HA	2.21	0.41
1:A:274:VAL:HG23	1:A:277:TYR:HB2	2.02	0.41
1:B:268:HIS:HB2	1:B:412:TYR:HE1	1.85	0.41
1:A:154:ASP:O	1:A:156:ASP:N	2.53	0.41
1:B:130:ILE:O	1:B:167:PHE:N	2.54	0.41
1:A:316:LYS:NZ	1:A:370:ASP:OD1	2.53	0.41
1:B:153:ILE:HG13	1:B:153:ILE:H	1.77	0.40
1:B:249:CYS:O	1:B:253:VAL:HB	2.20	0.40
1:B:145:GLY:HA3	1:B:148:ASP:OD2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	285/336 (85%)	259 (91%)	22 (8%)	4 (1%)	11	25
1	B	285/336 (85%)	266 (93%)	16 (6%)	3 (1%)	14	31
All	All	570/672 (85%)	525 (92%)	38 (7%)	7 (1%)	13	29

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	ASP
1	A	155	LYS
1	B	154	ASP
1	B	155	LYS
1	A	404	ARG
1	A	268	HIS
1	B	156	ASP

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	250/295 (85%)	218 (87%)	32 (13%)	4	9
1	B	250/295 (85%)	222 (89%)	28 (11%)	6	12
All	All	500/590 (85%)	440 (88%)	60 (12%)	5	10

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	ILE
1	A	124	VAL
1	A	125	LYS
1	A	137	ILE
1	A	153	ILE
1	A	154	ASP
1	A	155	LYS
1	A	162	LEU
1	A	164	MET
1	A	175	ASN
1	A	177	ILE
1	A	188	ASN
1	A	191	GLN
1	A	198	VAL
1	A	201	LEU
1	A	204	GLN
1	A	208	LYS

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Mol	Chain	Res	Type
1	A	226	LEU
1	A	227	LEU
1	A	238	LEU
1	A	241	SER
1	A	248	LEU
1	A	265	LEU
1	A	273	VAL
1	A	288	LEU
1	A	318	ARG
1	A	366	TYR
1	A	374	GLU
1	A	382	LEU
1	A	384	GLN
1	A	417	LEU
1	A	424	ASP
1	B	99	ILE
1	B	105	ILE
1	B	107	ASP
1	B	124	VAL
1	B	153	ILE
1	B	154	ASP
1	B	155	LYS
1	B	162	LEU
1	B	175	ASN
1	B	177	ILE
1	B	181	THR
1	B	188	ASN
1	B	189	MET
1	B	191	GLN
1	B	198	VAL
1	B	201	LEU
1	B	204	GLN
1	B	226	LEU
1	B	227	LEU
1	B	238	LEU
1	B	273	VAL
1	B	283	SER
1	B	311	SER
1	B	318	ARG
1	B	324	ASN
1	B	382	LEU
1	B	417	LEU

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Mol	Chain	Res	Type
1	B	424	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	175	ASN
1	A	184	ASN
1	A	200	ASN
1	A	203	HIS
1	A	204	GLN
1	A	223	GLN
1	A	302	ASN
1	A	384	GLN
1	B	141	ASN
1	B	175	ASN
1	B	184	ASN
1	B	188	ASN
1	B	200	ASN
1	B	223	GLN
1	B	254	ASN
1	B	302	ASN
1	B	384	GLN
1	B	418	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

4 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
3	TN3	A	500	-	21,21,21	0.93	2 (9%)	29,29,29	1.57	4 (13%)
2	NAD	A	450	-	42,48,48	1.33	5 (11%)	50,73,73	1.44	8 (16%)
2	NAD	B	451	-	42,48,48	1.25	4 (9%)	50,73,73	1.22	5 (10%)
3	TN3	B	501	-	21,21,21	1.07	3 (14%)	29,29,29	1.45	3 (10%)

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	TN3	A	500	-	-	0/10/10/10	0/2/2/2
2	NAD	A	450	-	-	4/26/62/62	0/5/5/5
2	NAD	B	451	-	-	8/26/62/62	0/5/5/5
3	TN3	B	501	-	-	1/10/10/10	0/2/2/2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	450	NAD	O7N-C7N	4.99	1.33	1.24
2	B	451	NAD	O7N-C7N	4.72	1.33	1.24
2	A	450	NAD	C7N-N7N	-3.63	1.26	1.33
2	B	451	NAD	C7N-N7N	-3.62	1.26	1.33
2	A	450	NAD	C5A-C4A	2.75	1.48	1.40
2	B	451	NAD	C5A-C4A	2.55	1.47	1.40
3	B	501	TN3	C14-CL21	2.48	1.79	1.73
3	B	501	TN3	C2-CL12	2.46	1.79	1.74
2	B	451	NAD	O4D-C1D	2.32	1.44	1.41
2	A	450	NAD	C2A-N3A	2.23	1.35	1.32
3	A	500	TN3	C14-CL21	2.19	1.78	1.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	TN3	C2-CL12	2.12	1.79	1.74
2	A	450	NAD	O4D-C1D	2.07	1.44	1.41
3	B	501	TN3	C11-N20	-2.07	1.30	1.33

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	TN3	C10-C11-N20	5.93	124.77	116.89
2	A	450	NAD	N3A-C2A-N1A	-4.41	121.78	128.68
3	B	501	TN3	C10-C11-N20	4.30	122.60	116.89
3	B	501	TN3	C5-O9-C7	3.60	126.61	118.04
2	B	451	NAD	N3A-C2A-N1A	-3.53	123.16	128.68
2	B	451	NAD	C4A-C5A-N7A	-3.48	105.77	109.40
2	A	450	NAD	O4D-C1D-C2D	-3.15	102.32	106.93
3	A	500	TN3	O13-C11-N20	-2.95	118.00	122.47
3	B	501	TN3	O13-C11-N20	-2.84	118.17	122.47
2	A	450	NAD	C1B-N9A-C4A	-2.72	121.87	126.64
2	B	451	NAD	C3N-C7N-N7N	2.49	120.73	117.75
2	A	450	NAD	C6N-N1N-C2N	-2.48	119.72	121.97
2	A	450	NAD	O7N-C7N-N7N	-2.45	119.10	122.58
3	A	500	TN3	C8-C9-C10	-2.40	117.99	120.78
2	A	450	NAD	O4B-C1B-C2B	-2.32	103.53	106.93
2	B	451	NAD	O7N-C7N-N7N	-2.27	119.36	122.58
2	A	450	NAD	C3N-C7N-N7N	2.20	120.39	117.75
2	B	451	NAD	C1B-N9A-C4A	-2.15	122.87	126.64
2	A	450	NAD	C4A-C5A-N7A	-2.08	107.23	109.40
3	A	500	TN3	O13-C11-C10	-2.08	117.23	120.94

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	450	NAD	C5D-O5D-PN-O3
2	A	450	NAD	C5D-O5D-PN-O2N
2	B	451	NAD	C4B-C5B-O5B-PA
2	B	451	NAD	PA-O3-PN-O5D
2	B	451	NAD	C5D-O5D-PN-O1N
2	B	451	NAD	O4B-C4B-C5B-O5B
2	B	451	NAD	C3B-C4B-C5B-O5B
2	A	450	NAD	O4B-C4B-C5B-O5B
2	B	451	NAD	C5D-O5D-PN-O2N
2	B	451	NAD	C5B-O5B-PA-O3

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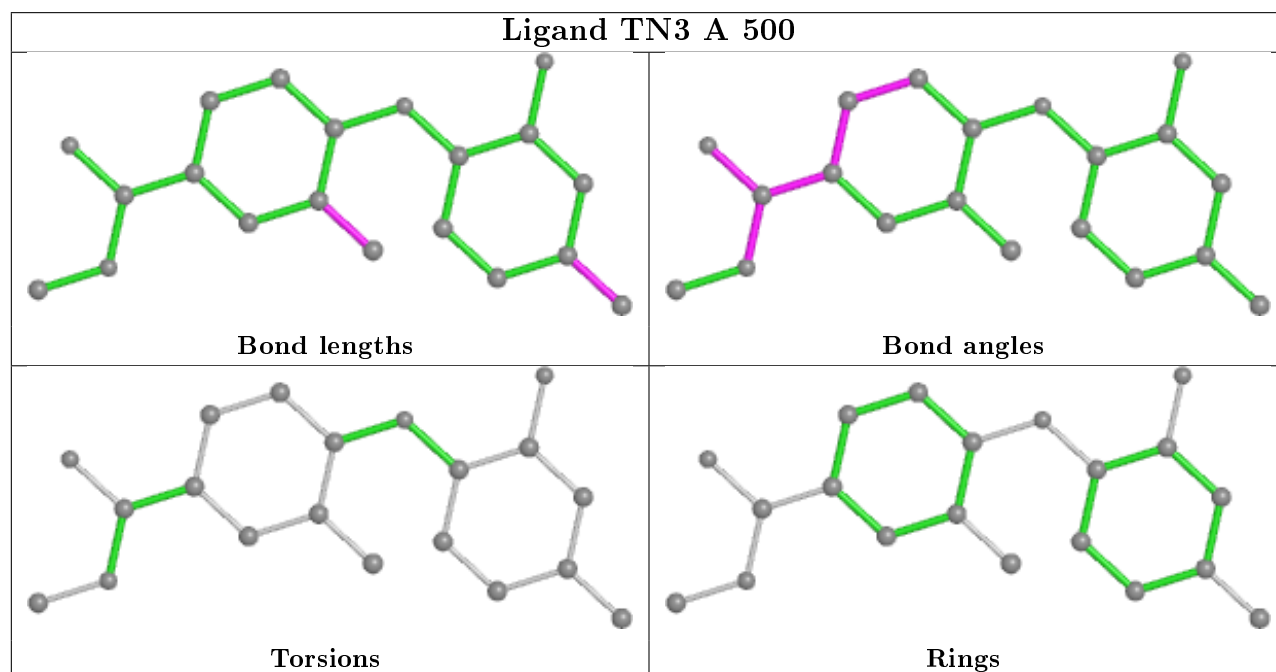
Mol	Chain	Res	Type	Atoms
2	B	451	NAD	C5D-O5D-PN-O3
2	A	450	NAD	C3B-C4B-C5B-O5B
3	B	501	TN3	C6-C5-O9-C7

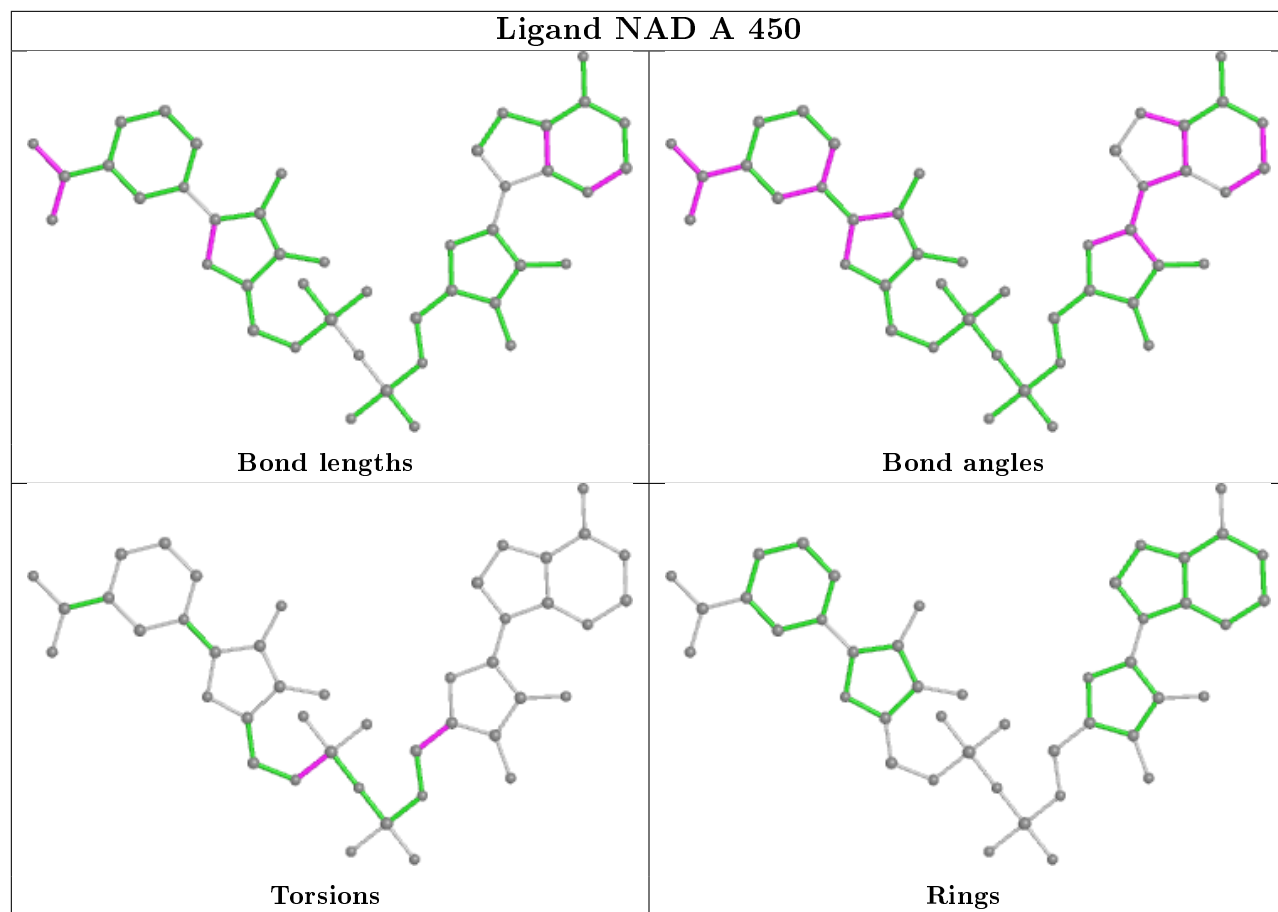
There are no ring outliers.

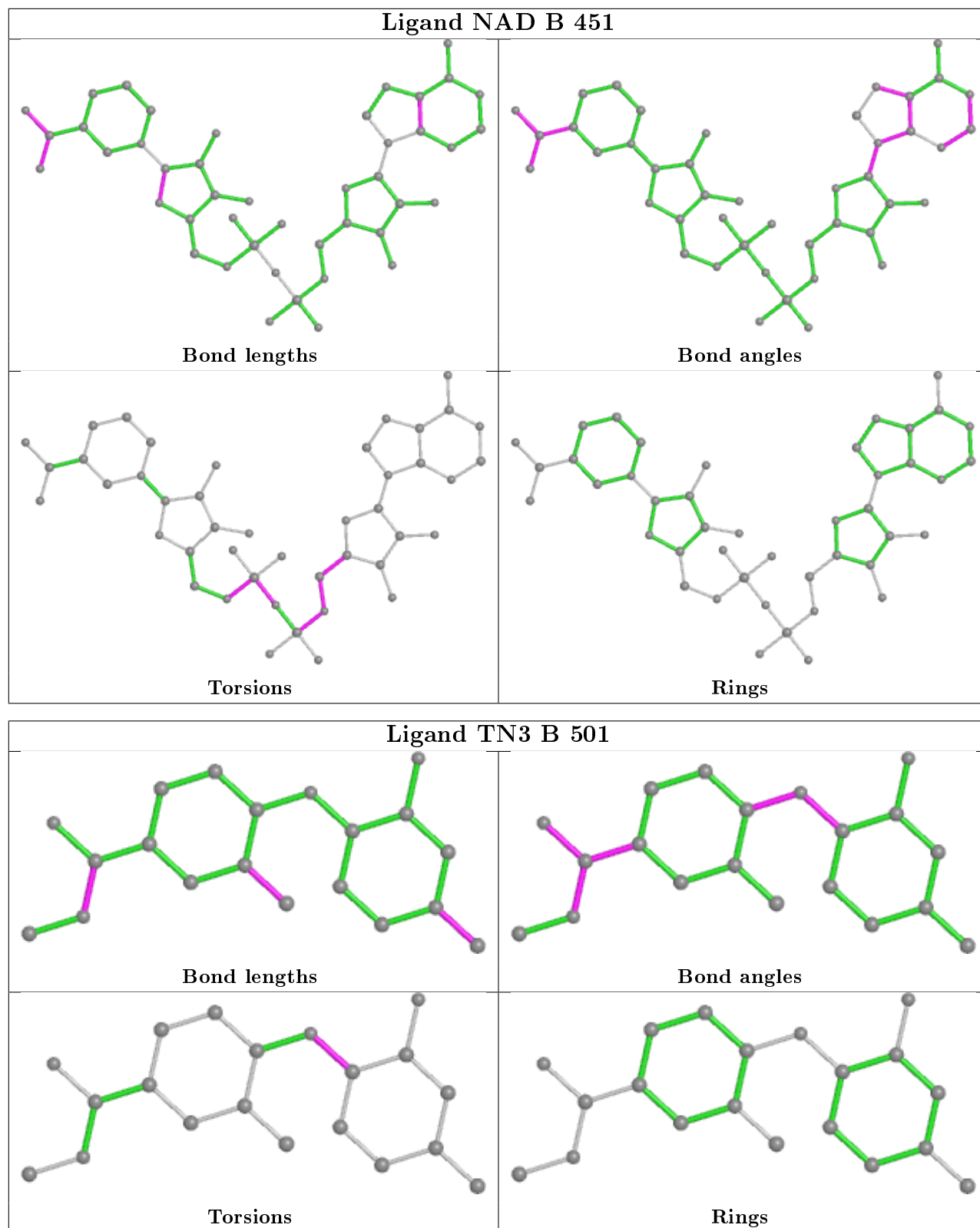
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	TN3	1	0
2	A	450	NAD	2	0
3	B	501	TN3	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 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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	289/336 (86%)	0.00	17 (5%)	22 20	14, 27, 67, 84	0
1	B	289/336 (86%)	0.13	17 (5%)	22 20	15, 28, 66, 83	0
All	All	578/672 (86%)	0.07	34 (5%)	22 20	14, 27, 67, 84	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	155	LYS	5.8
1	A	154	ASP	5.6
1	B	182	LYS	5.3
1	A	366	TYR	4.9
1	B	144	ASN	4.7
1	B	324	ASN	4.4
1	B	155	LYS	4.3
1	B	157	LYS	4.3
1	A	180	GLU	4.1
1	B	156	ASP	4.1
1	B	325	LYS	4.1
1	B	366	TYR	3.9
1	A	425	ASP	3.8
1	A	182	LYS	3.7
1	A	181	THR	3.6
1	A	186	ARG	3.5
1	A	183	ASN	3.5
1	B	183	ASN	3.5
1	B	141	ASN	3.1
1	A	179	GLU	3.0
1	A	178	ASP	2.9
1	A	325	LYS	2.9
1	B	146	LYS	2.8
1	B	158	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	97	GLU	2.6
1	B	149	ASN	2.5
1	B	179	GLU	2.4
1	A	189	MET	2.4
1	B	143	LYS	2.3
1	A	156	ASP	2.2
1	A	324	ASN	2.2
1	B	160	ASN	2.1
1	A	137	ILE	2.0
1	B	140	LYS	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 carbohydrates in this entry.

6.4 Ligands ⓘ

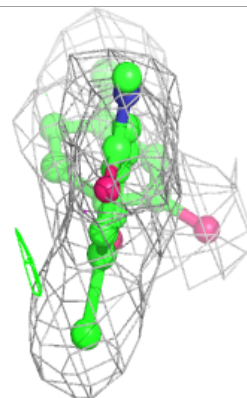
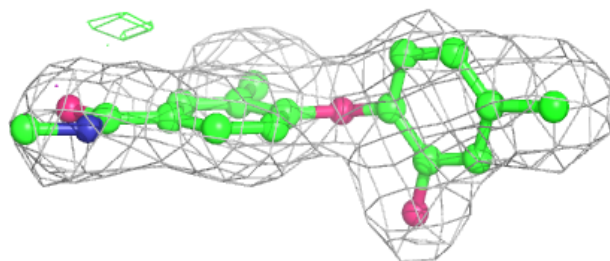
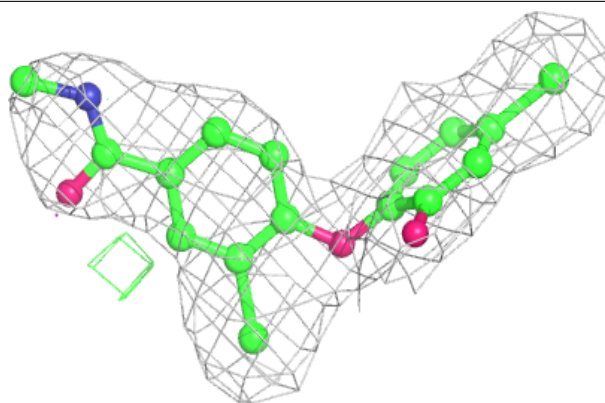
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	TN3	A	500	20/20	0.94	0.15	22,29,33,39	0
3	TN3	B	501	20/20	0.94	0.15	23,28,33,39	0
2	NAD	A	450	44/44	0.95	0.15	20,32,45,48	0
2	NAD	B	451	44/44	0.96	0.17	18,31,43,46	0

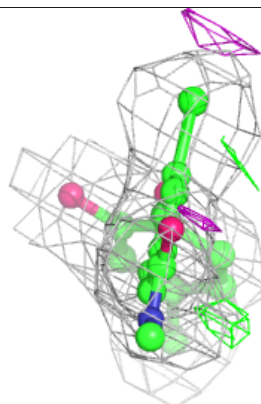
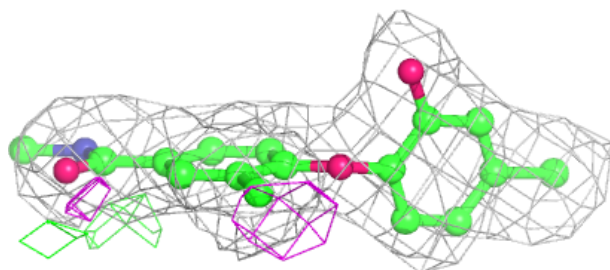
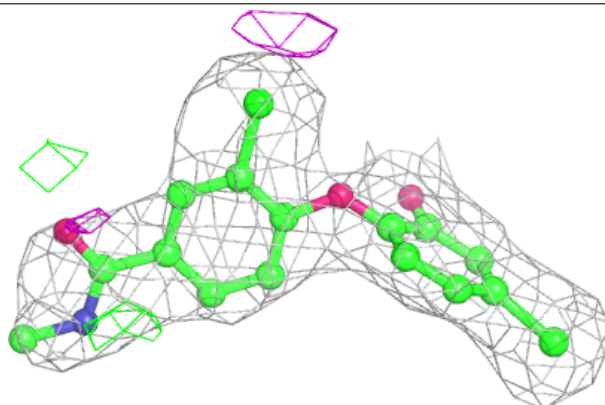
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TN3 A 500:

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

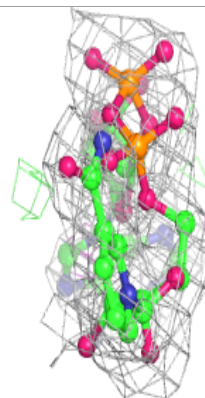
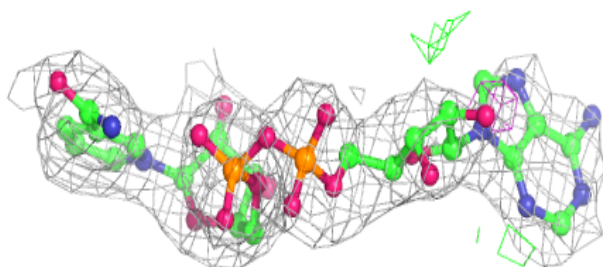
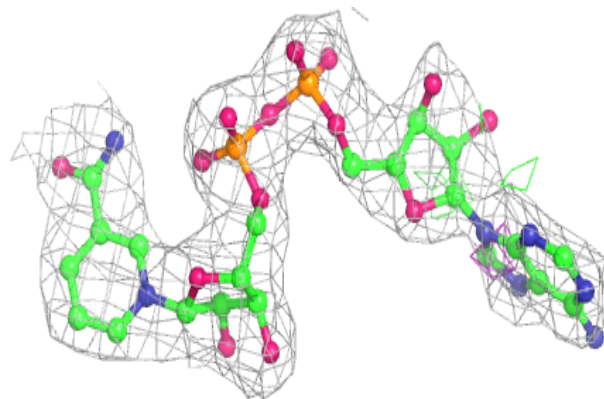
**Electron density around TN3 B 501:**

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

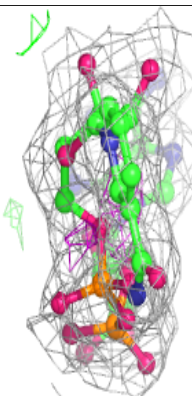
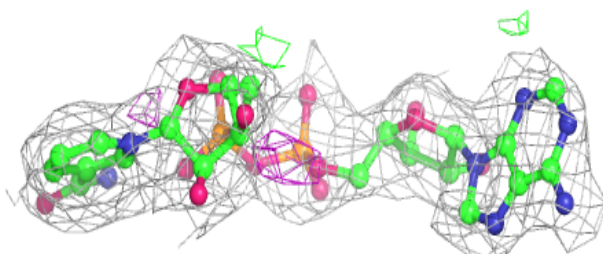
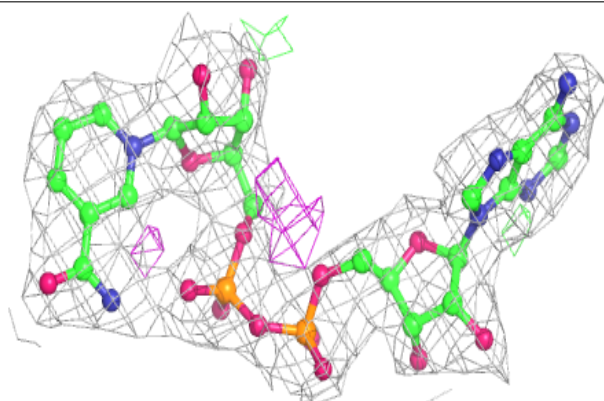


Electron density around NAD A 450:

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

**Electron density around NAD B 451:**

$2mF_o-DF_c$ (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.