



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:20 PM UTC

PDB ID : 1I3N / pdb_00001i3n
Title : MOLECULAR BASIS FOR SEVERE EPIMERASE-DEFICIENCY GALACTOSEMIA: X-RAY STRUCTURE OF THE HUMAN V94M-SUBSTITUTED UDP-GALACTOSE 4-EPIMERASE
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Deposited on : 2001-02-15
Resolution : 1.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

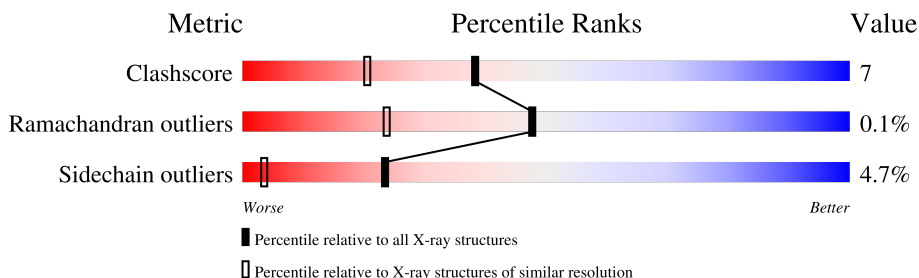
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	348	 85% 12% •
1	B	348	 82% 12% 5% •

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6500 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 UDP-GLUCOSE 4-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	2	0
			2697	1707	471	503	16			
1	B	345	Total	C	N	O	S	0	8	0
			2706	1714	467	509	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	MET	VAL	engineered mutation	UNP Q14376
B	94	MET	VAL	engineered mutation	UNP Q14376

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cl	0	0
			2	2		
2	B	2	Total	Cl	0	0
			2	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

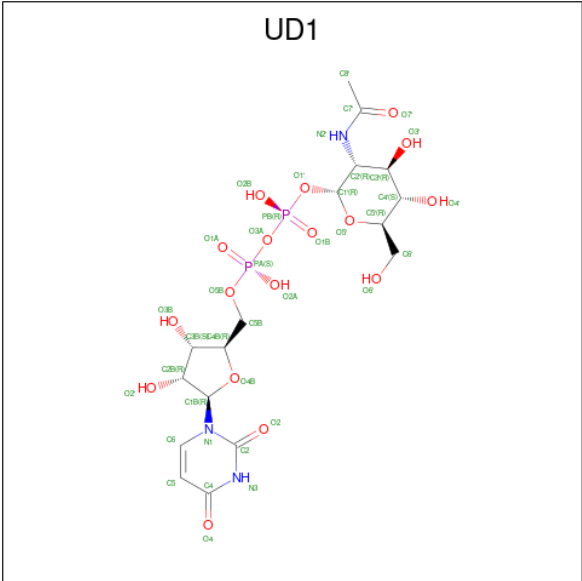
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

- Molecule 5 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (CCD ID: UD1) (formula: C₁₇H₂₇N₃O₁₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	1
			40	17	3	18	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

- Molecule 6 is water.

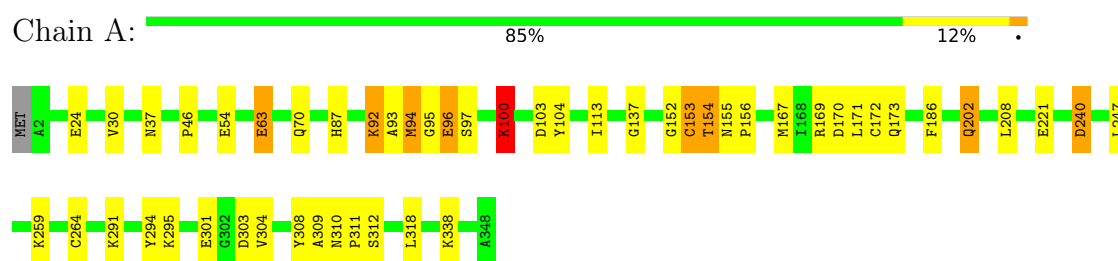
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	498	Total	O	0	0
			498	498		
6	B	427	Total	O	0	0
			427	427		

3 Residue-property plots

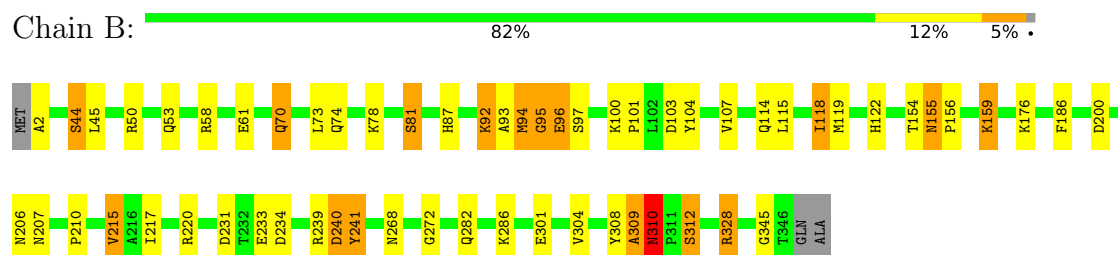
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: UDP-GLUCOSE 4-EPIMERASE



• Molecule 1: UDP-GLUCOSE 4-EPIMERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.10 Å 90.00 Å 96.80 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.50	Depositor
% Data completeness (in resolution range)	97.3 (50.00-1.50)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.180 , 0.215	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6500	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAD, UD1, CL, 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	0.95	1/2762 (0.0%)	1.39	24/3733 (0.6%)
1	B	0.98	0/2799	1.39	23/3787 (0.6%)
All	All	0.97	1/5561 (0.0%)	1.39	47/7520 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	54	GLU	CD-OE2	5.06	1.34	1.25

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	CYS	CA-C-N	10.31	141.24	121.54
1	A	153	CYS	C-N-CA	10.31	141.24	121.54
1	B	96	GLU	N-CA-C	10.11	122.38	111.36
1	A	100	LYS	CB-CA-C	-10.04	97.61	110.96
1	A	155	ASN	N-CA-C	9.59	121.82	110.31
1	A	153	CYS	N-CA-C	9.58	125.21	110.36
1	A	154	THR	N-CA-C	9.20	130.39	110.80
1	B	95	GLY	CA-C-N	8.84	132.84	120.29
1	B	95	GLY	C-N-CA	8.84	132.84	120.29
1	B	94	MET	N-CA-C	-8.59	101.04	112.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	92	LYS	CA-C-N	-7.96	110.18	121.72
1	B	92	LYS	C-N-CA	-7.96	110.18	121.72
1	A	96	GLU	N-CA-C	7.79	119.56	111.14
1	A	208	LEU	N-CA-C	7.62	119.27	110.97
1	A	94	MET	N-CA-C	-7.07	102.93	112.26
1	A	92	LYS	CA-C-N	-7.00	111.74	122.09
1	A	92	LYS	C-N-CA	-7.00	111.74	122.09
1	A	186	PHE	CA-CB-CG	-6.76	107.04	113.80
1	B	312	SER	N-CA-C	6.63	118.17	111.07
1	B	155	ASN	CA-CB-CG	6.39	118.99	112.60
1	A	240	ASP	CA-CB-CG	6.25	118.85	112.60
1	A	312	SER	N-CA-C	6.23	118.07	111.28
1	B	309	ALA	N-CA-C	6.16	120.11	110.32
1	B	240	ASP	CA-CB-CG	5.84	118.44	112.60
1	B	92	LYS	N-CA-CB	5.82	119.81	110.85
1	B	186	PHE	CA-CB-CG	-5.82	107.98	113.80
1	A	30	VAL	N-CA-C	-5.79	99.20	107.77
1	A	95	GLY	CA-C-N	5.76	128.28	120.44
1	A	95	GLY	C-N-CA	5.76	128.28	120.44
1	A	63	GLU	CA-CB-CG	-5.73	102.64	114.10
1	A	87	HIS	N-CA-C	5.52	118.28	107.57
1	B	328	ARG	NE-CZ-NH1	5.46	126.96	121.50
1	A	46	PRO	N-CA-C	-5.42	102.60	111.21
1	B	200	ASP	CA-CB-CG	5.41	118.00	112.60
1	B	241	TYR	CB-CA-C	5.34	118.55	109.84
1	B	310	ASN	N-CA-C	-5.28	99.18	108.69
1	B	94	MET	CB-CA-C	5.24	117.66	111.22
1	B	215	VAL	N-CA-C	-5.23	105.44	110.72
1	A	309	ALA	N-CA-C	5.22	118.24	110.52
1	B	107	VAL	N-CA-C	5.22	115.43	110.53
1	A	310	ASN	N-CA-C	-5.18	98.28	109.01
1	B	45	LEU	N-CA-C	5.16	117.01	109.84
1	A	221	GLU	N-CA-C	5.11	116.66	111.14
1	A	152	GLY	N-CA-C	5.08	118.64	112.49
1	B	234	ASP	CA-CB-CG	5.06	117.66	112.60
1	B	268	ASN	CB-CA-C	-5.05	100.99	109.53
1	B	87	HIS	N-CA-C	5.04	117.34	107.57

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	154	THR	CA

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	2697	0	2661	25	0
1	B	2706	0	2655	50	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
4	A	44	0	26	0	0
4	B	44	0	26	0	0
5	A	40	0	6	0	0
5	B	39	0	25	3	0
6	A	498	0	0	5	0
6	B	427	0	0	12	0
All	All	6500	0	5399	76	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 7.

All (76) 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:73:LEU:HD13	1:B:118[A]:ILE:HD13	1.40	1.03
5:B:901:UD1:H5'	6:B:1985:HOH:O	1.84	0.78
1:B:70:GLN:HG3	6:B:1270:HOH:O	1.83	0.77
1:B:155:ASN:HB3	1:B:156:PRO:HD2	1.66	0.76
1:A:63:GLU:HG3	6:A:1961:HOH:O	1.84	0.76
1:B:73:LEU:HD13	1:B:118[A]:ILE:CD1	2.20	0.68
1:B:73:LEU:HD22	1:B:118[A]:ILE:CD1	2.24	0.67
1:B:73:LEU:HB3	1:B:118[A]:ILE:HD12	1.75	0.67
1:A:97:SER:HA	1:A:104:TYR:CE1	2.34	0.63
1:A:97:SER:O	1:A:156:PRO:HG2	1.98	0.62
1:B:154:THR:HB	6:B:1939:HOH:O	1.99	0.62
1:B:310:ASN:HD22	1:B:310:ASN:C	2.09	0.61
1:B:94:MET:HB2	1:B:96:GLU:HG3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:SER:HA	1:B:104:TYR:CE1	2.36	0.59
1:B:345:GLY:HA2	6:B:1338:HOH:O	2.02	0.58
1:B:215:VAL:HG22	1:B:220:ARG:HB2	1.84	0.57
1:B:286:LYS:HE2	6:B:1761:HOH:O	2.05	0.57
1:B:93:ALA:C	1:B:95:GLY:H	2.10	0.57
1:A:94:MET:HB2	1:A:96:GLU:HG3	1.85	0.57
1:A:37:ASN:OD1	1:A:202:GLN:NE2	2.37	0.56
1:B:53:GLN:HG3	1:B:58:ARG:O	2.05	0.56
1:A:113:ILE:HG12	1:A:167:MET:HE1	1.88	0.56
1:B:44:SER:HB2	6:B:1613:HOH:O	2.06	0.55
1:B:73:LEU:HD22	1:B:118[A]:ILE:HD11	1.86	0.55
1:A:100:LYS:HD3	1:A:103:ASP:HB2	1.87	0.55
1:B:282[B]:GLN:HG3	6:B:1760:HOH:O	2.07	0.55
1:B:310:ASN:HD22	1:B:312:SER:H	1.55	0.55
1:B:70:GLN:NE2	1:B:114[A]:GLN:OE1	2.43	0.52
1:A:94:MET:HE2	6:A:1112:HOH:O	2.10	0.52
5:B:901:UD1:O1'	5:B:901:UD1:O7'	2.28	0.52
1:B:310:ASN:ND2	1:B:312:SER:H	2.08	0.51
1:B:73:LEU:HB3	1:B:118[A]:ILE:CD1	2.41	0.51
1:B:94:MET:HB2	1:B:96:GLU:OE2	2.09	0.51
1:A:338:LYS:HE3	6:A:1573:HOH:O	2.11	0.51
1:B:74:GLN:O	1:B:78:LYS:HG3	2.12	0.50
1:B:73:LEU:CD1	1:B:114[A]:GLN:HG3	2.42	0.50
1:A:70:GLN:OE1	6:A:1996:HOH:O	2.19	0.50
1:B:70:GLN:HG2	1:B:114[A]:GLN:OE1	2.12	0.50
1:A:100:LYS:HD2	1:A:104:TYR:HE1	1.77	0.49
1:A:240:ASP:HB2	1:A:308:TYR:HA	1.94	0.49
1:B:94:MET:HB2	1:B:96:GLU:CG	2.43	0.49
1:A:100:LYS:HD2	1:A:104:TYR:CE1	2.49	0.48
1:B:239:ARG:HB2	1:B:241:TYR:CE1	2.49	0.48
1:B:61:GLU:HG2	6:B:1598:HOH:O	2.13	0.47
1:B:207[B]:ASN:CG	1:B:210:PRO:HG2	2.39	0.47
1:B:240:ASP:HB2	1:B:308:TYR:HA	1.96	0.47
5:B:901:UD1:H6'1	6:B:1940:HOH:O	2.14	0.46
1:A:311:PRO:HD2	6:A:1545:HOH:O	2.14	0.46
1:B:73:LEU:HD12	1:B:114[A]:GLN:HG3	1.98	0.46
1:A:171:LEU:HA	1:B:101:PRO:HG2	1.98	0.45
1:B:115:LEU:O	1:B:118[A]:ILE:HG12	2.16	0.45
1:B:73:LEU:HD22	1:B:118[A]:ILE:HD12	1.98	0.45
1:A:94:MET:HB2	1:A:96:GLU:CG	2.47	0.44
1:B:328:ARG:NH2	6:B:1989:HOH:O	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:ALA:N	1:B:81:SER:HB3	2.32	0.44
1:A:318:LEU:HD12	1:A:318:LEU:HA	1.80	0.44
1:B:115:LEU:O	1:B:118[B]:ILE:HG22	2.18	0.44
1:B:122:HIS:ND1	6:B:1632:HOH:O	2.36	0.43
1:B:272:GLY:HA2	1:B:309:ALA:O	2.18	0.43
1:A:167:MET:HE3	1:A:167:MET:HB3	1.83	0.43
1:A:100:LYS:HE2	1:A:103:ASP:OD1	2.19	0.43
1:A:93:ALA:HA	1:A:104:TYR:OH	2.18	0.43
1:A:170:ASP:OD2	1:B:159:LYS:NZ	2.51	0.43
1:B:100:LYS:HE3	1:B:103:ASP:HB2	2.00	0.43
1:B:73:LEU:CD1	1:B:118[A]:ILE:HD13	2.29	0.42
1:B:70:GLN:HB3	6:B:1854:HOH:O	2.19	0.42
1:B:155:ASN:HB3	1:B:156:PRO:CD	2.36	0.42
1:B:118[B]:ILE:CG2	1:B:119:MET:N	2.82	0.42
1:A:172:CYS:SG	1:A:264:CYS:HB2	2.60	0.41
1:A:294:TYR:CD2	1:A:294:TYR:C	2.98	0.41
1:A:137:GLY:HA2	1:A:153:CYS:HB2	2.02	0.41
1:B:94:MET:HA	1:B:206:ASN:ND2	2.36	0.41
1:B:93:ALA:C	1:B:95:GLY:N	2.77	0.41
1:B:215:VAL:CG2	1:B:220:ARG:HB2	2.50	0.41
1:A:173:GLN:NE2	1:A:173:GLN:HA	2.37	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	347/348 (100%)	340 (98%)	6 (2%)	1 (0%)	36	17
1	B	352/348 (101%)	341 (97%)	11 (3%)	0	100	100
All	All	699/696 (100%)	681 (97%)	17 (2%)	1 (0%)	48	24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	THR

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	282/282 (100%)	270 (96%)	12 (4%)	26	4
1	B	287/282 (102%)	271 (94%)	16 (6%)	19	2
All	All	569/564 (101%)	541 (95%)	28 (5%)	23	3

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLU
1	A	92	LYS
1	A	100	LYS
1	A	169	ARG
1	A	202	GLN
1	A	247	LEU
1	A	259	LYS
1	A	291	LYS
1	A	295	LYS
1	A	301	GLU
1	A	303	ASP
1	A	304	VAL
1	B	44	SER
1	B	50	ARG
1	B	70	GLN
1	B	81	SER
1	B	92	LYS
1	B	118[A]	ILE
1	B	118[B]	ILE
1	B	159	LYS
1	B	176	LYS
1	B	217	ILE

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Mol	Chain	Res	Type
1	B	231[A]	ASP
1	B	231[B]	ASP
1	B	233	GLU
1	B	301	GLU
1	B	304	VAL
1	B	310	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	138	ASN
1	A	173	GLN
1	A	202	GLN
1	A	224	ASN
1	A	339	GLN
1	B	74	GLN
1	B	138	ASN
1	B	202	GLN
1	B	206	ASN
1	B	310	ASN
1	B	339	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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	$\# Z > 2$
5	UD1	A	401[B]	-	40,41,41	1.25	2 (5%)	59,62,62	1.08	4 (6%)
5	UD1	A	401[A]	-	40,41,41	1.25	2 (5%)	59,62,62	1.08	4 (6%)
5	UD1	B	901	-	40,41,41	1.00	3 (7%)	59,62,62	1.09	5 (8%)
4	NAD	A	400	-	46,48,48	1.30	6 (13%)	64,73,73	1.54	6 (9%)
4	NAD	B	900	-	46,48,48	1.11	5 (10%)	64,73,73	1.59	6 (9%)

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
5	UD1	A	401[B]	-	-	2/26/63/63	0/3/3/3
5	UD1	A	401[A]	-	-	2/26/63/63	0/3/3/3
5	UD1	B	901	-	-	4/26/63/63	0/3/3/3
4	NAD	A	400	-	-	6/30/62/62	0/5/5/5
4	NAD	B	900	-	-	7/30/62/62	0/5/5/5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	401[A]	UD1	PA-O3A	5.33	1.65	1.59
5	A	401[B]	UD1	PA-O3A	5.33	1.65	1.59
4	A	400	NAD	C2N-N1N	4.04	1.39	1.35
4	A	400	NAD	PN-O3	3.25	1.63	1.59
4	B	900	NAD	C4N-C3N	3.10	1.44	1.39
5	A	401[A]	UD1	PB-O3A	-2.98	1.56	1.59
5	A	401[B]	UD1	PB-O3A	-2.98	1.56	1.59
5	B	901	UD1	C2'-N2'	2.74	1.50	1.45
4	A	400	NAD	PA-O3	-2.51	1.56	1.59
5	B	901	UD1	PA-O3A	2.51	1.62	1.59
4	B	900	NAD	PN-O3	2.45	1.62	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	900	NAD	PA-O2A	-2.45	1.44	1.55
4	A	400	NAD	C2A-N1A	2.43	1.38	1.33
4	A	400	NAD	C4N-C3N	2.41	1.43	1.39
4	A	400	NAD	C3N-C7N	2.28	1.54	1.50
5	B	901	UD1	PB-O3A	-2.17	1.57	1.59
4	B	900	NAD	C6N-N1N	2.13	1.40	1.35
4	B	900	NAD	C8A-N7A	2.06	1.35	1.31

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	400	NAD	C5N-C4N-C3N	-6.52	113.96	120.36
4	B	900	NAD	C5N-C4N-C3N	-6.10	114.36	120.36
4	B	900	NAD	C5N-C6N-N1N	-5.95	112.26	120.38
4	A	400	NAD	C6N-C5N-C4N	5.93	127.99	119.45
4	B	900	NAD	C6N-N1N-C2N	4.96	126.09	121.88
4	A	400	NAD	C5N-C6N-N1N	-4.75	113.90	120.38
4	B	900	NAD	C6N-C5N-C4N	4.46	125.87	119.45
5	A	401[A]	UD1	C2'-N2'-C7'	-3.13	115.77	123.11
5	A	401[B]	UD1	C2'-N2'-C7'	-3.13	115.77	123.11
4	B	900	NAD	C3N-C7N-N7N	3.13	121.60	117.74
5	A	401[A]	UD1	O5'-C1'-O1'	-2.97	107.48	111.36
5	A	401[B]	UD1	O5'-C1'-O1'	-2.97	107.48	111.36
5	B	901	UD1	O5'-C5'-C4'	2.71	114.59	109.70
5	A	401[A]	UD1	C8'-C7'-N2'	2.61	120.45	116.12
5	A	401[B]	UD1	C8'-C7'-N2'	2.61	120.45	116.12
4	A	400	NAD	C3N-C7N-N7N	2.58	120.92	117.74
4	A	400	NAD	C2N-C3N-C4N	2.53	121.20	118.26
5	B	901	UD1	C4'-C3'-C2'	2.42	113.93	110.40
5	B	901	UD1	O3A-PB-O1B	-2.37	103.59	110.70
5	B	901	UD1	O5'-C5'-C6'	2.09	111.61	106.44
4	B	900	NAD	C2N-C3N-C4N	2.08	120.68	118.26
5	A	401[A]	UD1	O3A-PB-O1B	-2.06	104.50	110.70
5	A	401[B]	UD1	O3A-PB-O1B	-2.06	104.50	110.70
4	A	400	NAD	C6N-N1N-C2N	2.06	123.63	121.88
5	B	901	UD1	O3'-C3'-C4'	-2.00	105.65	110.38

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	400	NAD	O4D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
4	B	900	NAD	C5D-O5D-PN-O2N
4	B	900	NAD	O4D-C1D-N1N-C2N
4	B	900	NAD	O4D-C1D-N1N-C6N
5	A	401[A]	UD1	C1'-O1'-PB-O3A
5	A	401[B]	UD1	C1'-O1'-PB-O3A
5	B	901	UD1	C1'-O1'-PB-O3A
5	A	401[A]	UD1	C1'-O1'-PB-O2B
5	A	401[B]	UD1	C1'-O1'-PB-O2B
4	A	400	NAD	C5B-O5B-PA-O2A
4	B	900	NAD	C5D-O5D-PN-O3
4	A	400	NAD	PA-O3-PN-O2N
4	B	900	NAD	PA-O3-PN-O2N
5	B	901	UD1	PA-O3A-PB-O1B
5	B	901	UD1	C3'-C2'-N2'-C7'
4	B	900	NAD	C2D-C1D-N1N-C6N
4	A	400	NAD	O4D-C1D-N1N-C6N
4	B	900	NAD	PA-O3-PN-O1N
5	B	901	UD1	C1'-C2'-N2'-C7'
4	A	400	NAD	O4B-C4B-C5B-O5B
4	A	400	NAD	PA-O3-PN-O1N

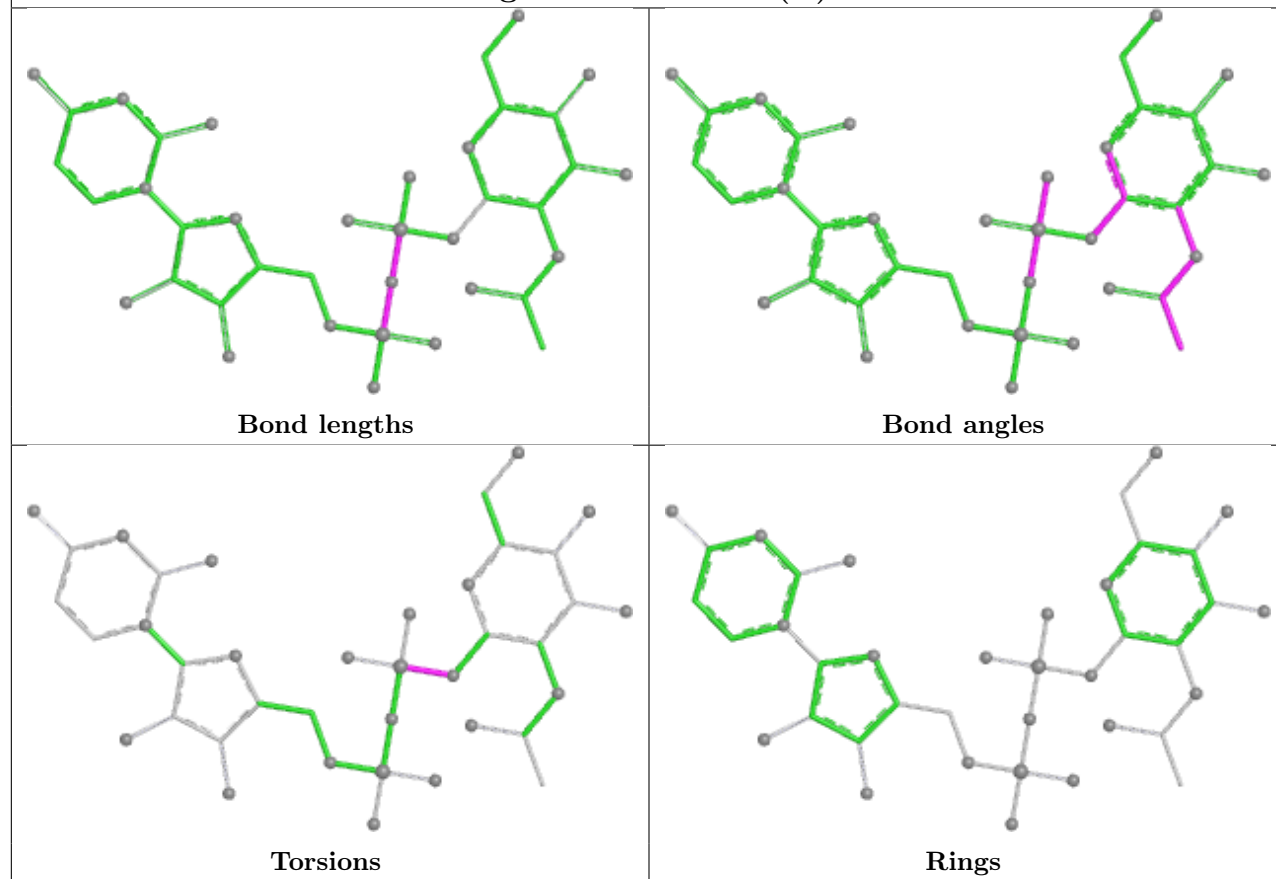
There are no ring outliers.

1 monomer is involved in 3 short contacts:

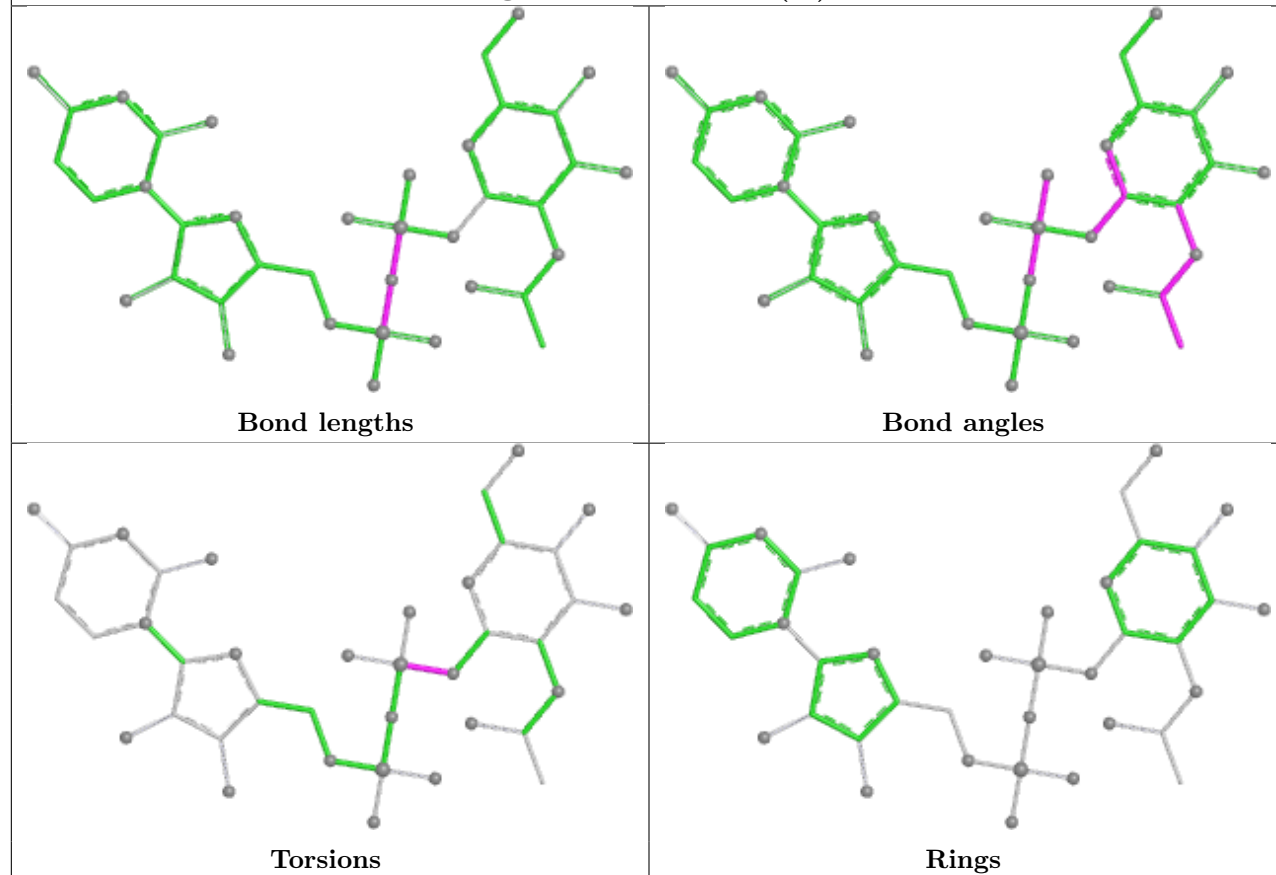
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	901	UD1	3	0

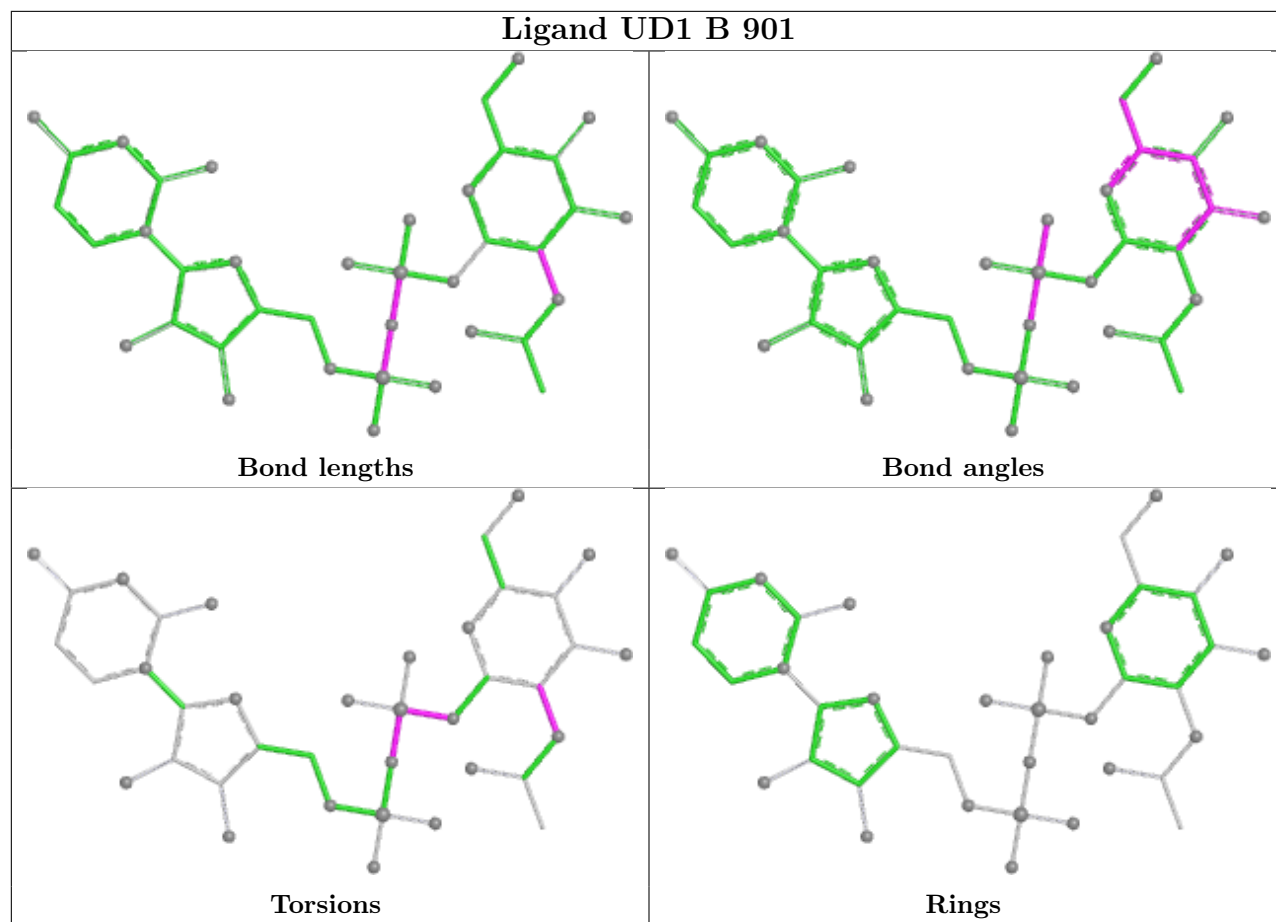
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

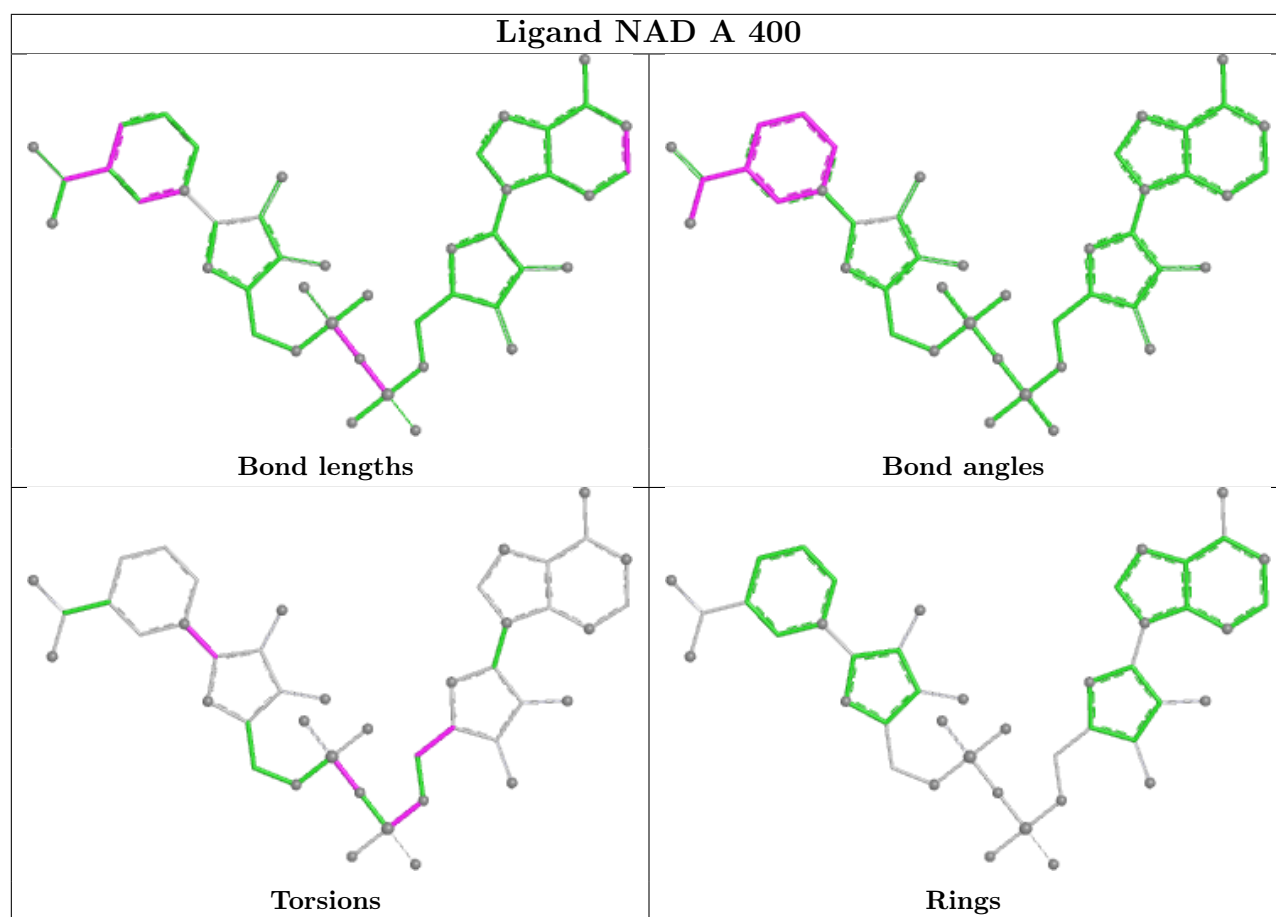
Ligand UD1 A 401 (B)

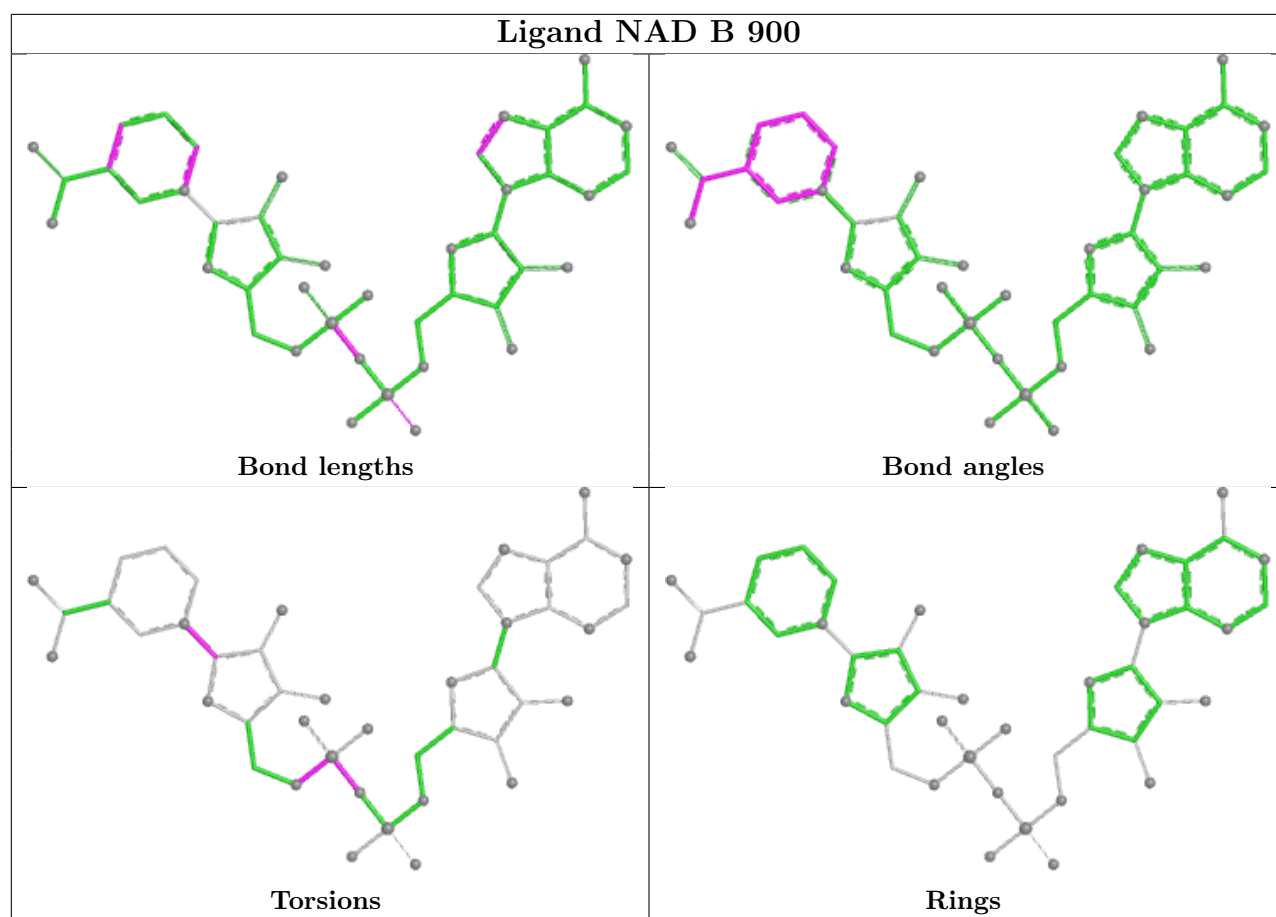


Ligand UD1 A 401 (A)









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.