



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:08 AM UTC

PDB ID : 7BTJ / pdb_00007btj
Title : Crystal structure of Pennisetum glaucum monodehydroascorbate reductase in complex with FADH2
Authors : Sonkar, K.S.; Achary, M.M.; Reddy, M.K.; Arulandu, A.
Deposited on : 2020-04-01
Resolution : 2.37 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

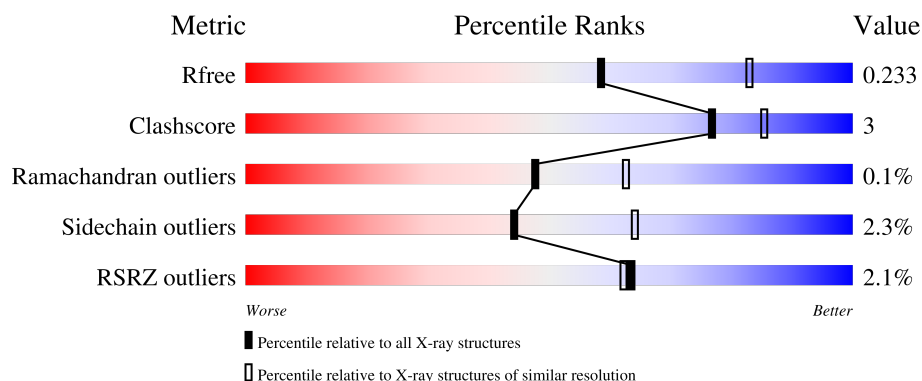
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	435	
1	B	435	
1	C	435	
1	D	435	

2 Entry composition [i](#)

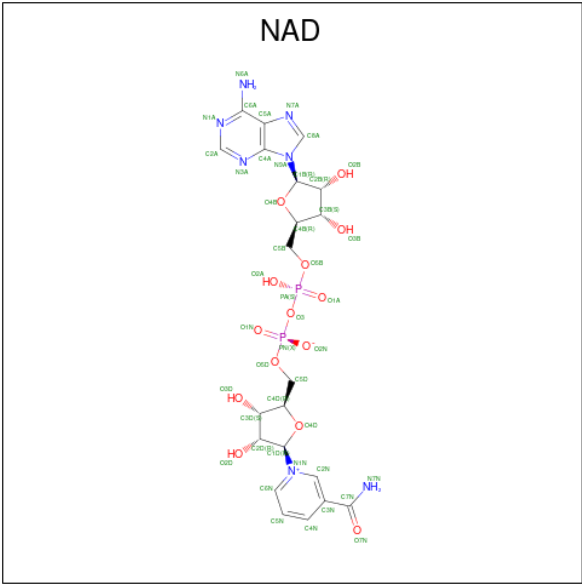
There are 4 unique types of molecules in this entry. The entry contains 13788 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pennisetum glaucum monodehydroascorbate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	432	Total	C	N	O	S	0	0	0
			3255	2098	529	623	5			
1	B	432	Total	C	N	O	S	0	0	0
			3259	2102	532	620	5			
1	C	435	Total	C	N	O	S	0	0	0
			3235	2083	525	621	6			
1	D	429	Total	C	N	O	S	0	0	0
			3098	1992	498	603	5			

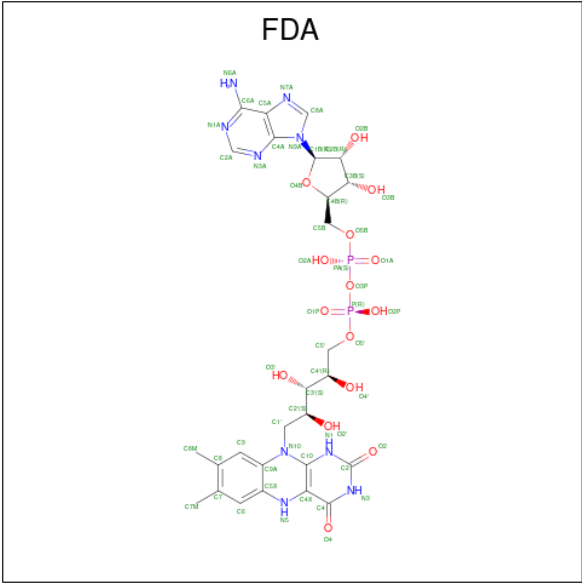
- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

- Molecule 3 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (CCD ID: FDA) (formula: C₂₇H₃₅N₉O₁₅P₂) (labeled as "Ligand of Interest" by depositor).



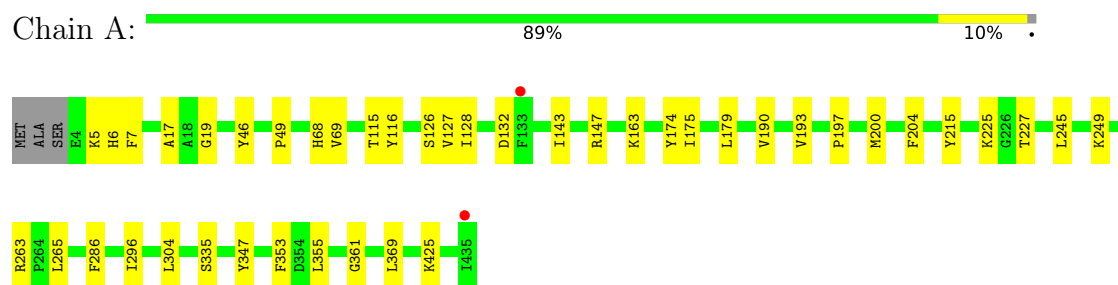
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	47	Total	O	0	0
			47	47		

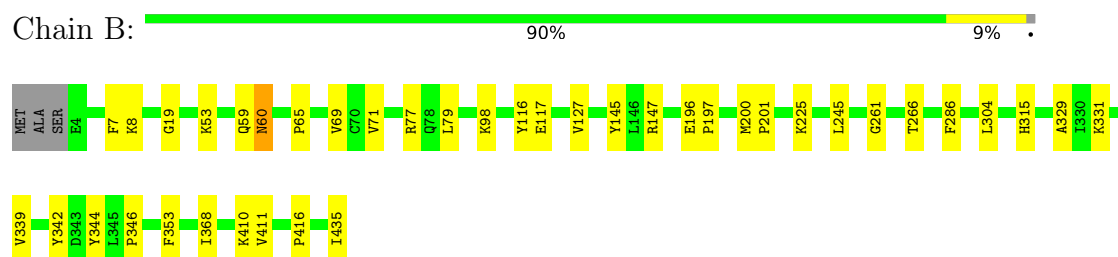
3 Residue-property plots

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.

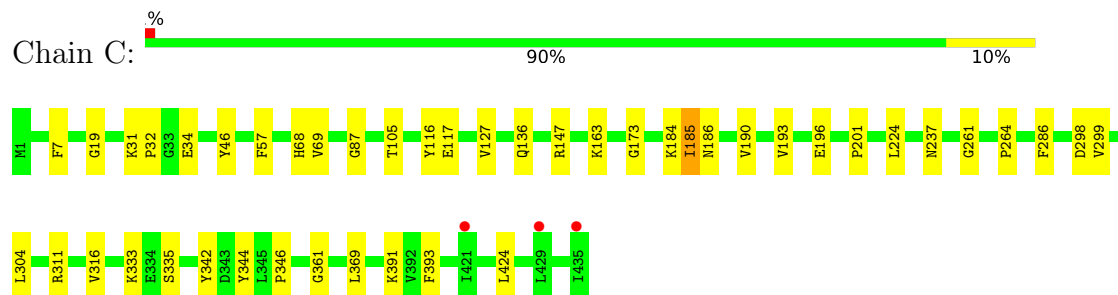
- Molecule 1: *Pennisetum glaucum* monodehydroascorbate reductase



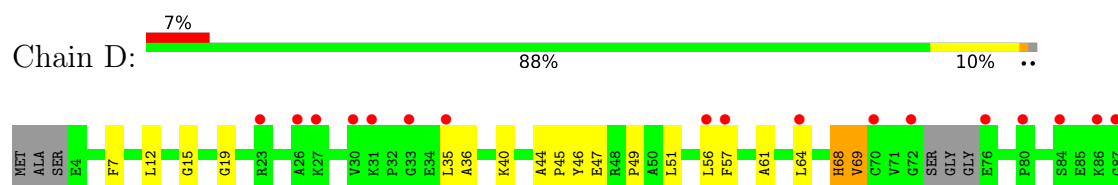
- Molecule 1: *Pennisetum glaucum* monodehydroascorbate reductase

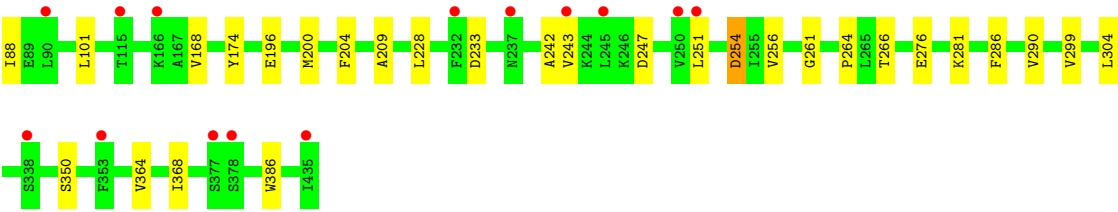


- Molecule 1: *Pennisetum glaucum* monodehydroascorbate reductase



- Molecule 1: *Pennisetum glaucum* monodehydroascorbate reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	65.73Å 79.12Å 90.69Å 96.39° 97.27° 111.92°	Depositor
Resolution (Å)	61.45 – 2.37 61.45 – 2.37	Depositor EDS
% Data completeness (in resolution range)	96.6 (61.45-2.37) 96.6 (61.45-2.37)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.37Å)	Xtriage
Refinement program	BUSTER 2.10.4 (11-DEC-2020)	Depositor
R, R_{free}	0.186 , 0.243 0.180 , 0.233	Depositor DCC
R_{free} test set	3231 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13788	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% 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: FDA, 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	0.79	1/3327 (0.0%)	1.03	7/4509 (0.2%)
1	B	0.78	2/3331 (0.1%)	1.06	10/4514 (0.2%)
1	C	0.75	0/3307	1.08	8/4489 (0.2%)
1	D	0.73	0/3168	1.08	12/4320 (0.3%)
All	All	0.76	3/13133 (0.0%)	1.06	37/17832 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	53	LYS	CA-C	8.04	1.56	1.52
1	B	353	PHE	CA-C	6.01	1.58	1.53
1	A	353	PHE	CA-C	5.20	1.57	1.53

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	286	PHE	CA-CB-CG	-8.21	105.59	113.80
1	B	286	PHE	CA-CB-CG	-7.25	106.55	113.80
1	C	186	ASN	CA-C-N	7.07	132.16	122.19
1	C	186	ASN	C-N-CA	7.07	132.16	122.19
1	C	361	GLY	N-CA-C	7.04	120.11	111.45
1	B	60	ASN	CA-CB-CG	6.87	119.47	112.60
1	D	261	GLY	N-CA-C	6.53	119.61	112.29
1	C	298	ASP	CA-CB-CG	6.09	118.69	112.60
1	C	190	VAL	N-CA-C	6.07	116.67	108.17
1	A	190	VAL	N-CA-C	6.06	116.66	108.17
1	C	261	GLY	N-CA-C	6.05	118.21	112.04
1	D	286	PHE	CA-CB-CG	-5.98	107.82	113.80
1	A	361	GLY	N-CA-C	5.96	118.79	111.45
1	B	315	HIS	CA-C-N	5.96	128.63	120.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	315	HIS	C-N-CA	5.96	128.63	120.46
1	B	416	PRO	N-CA-C	5.90	116.13	110.47
1	D	69	VAL	CA-C-N	5.85	132.24	121.70
1	D	69	VAL	C-N-CA	5.85	132.24	121.70
1	C	286	PHE	CA-CB-CG	-5.78	108.02	113.80
1	A	204	PHE	CA-CB-CG	5.75	119.55	113.80
1	C	117	GLU	N-CA-C	-5.73	105.96	113.12
1	D	254	ASP	CA-CB-CG	5.62	118.22	112.60
1	B	117	GLU	N-CA-C	-5.50	106.62	113.55
1	B	261	GLY	N-CA-C	5.28	117.43	112.04
1	B	353	PHE	CA-CB-CG	5.28	119.08	113.80
1	A	425	LYS	CA-C-N	5.24	127.72	120.29
1	A	425	LYS	C-N-CA	5.24	127.72	120.29
1	D	204	PHE	CA-CB-CG	5.23	119.03	113.80
1	D	304	LEU	CA-C-N	5.19	128.61	120.82
1	D	304	LEU	C-N-CA	5.19	128.61	120.82
1	D	47	GLU	CA-C-N	5.17	127.41	120.58
1	D	47	GLU	C-N-CA	5.17	127.41	120.58
1	B	266	THR	CA-C-N	5.11	127.64	120.28
1	B	266	THR	C-N-CA	5.11	127.64	120.28
1	D	209	ALA	CA-C-N	5.03	127.27	120.38
1	D	209	ALA	C-N-CA	5.03	127.27	120.38
1	A	143	ILE	N-CA-C	-5.02	99.72	106.85

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	3255	0	3216	18	0
1	B	3259	0	3229	18	0
1	C	3235	0	3155	16	0
1	D	3098	0	2868	23	0
2	A	44	0	26	1	0
2	B	44	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	44	0	26	1	0
2	D	44	0	26	0	0
3	A	53	0	33	1	0
3	B	53	0	33	0	0
3	C	53	0	33	0	0
3	D	53	0	33	3	0
4	A	191	0	0	1	0
4	B	180	0	0	0	0
4	C	135	0	0	1	0
4	D	47	0	0	0	0
All	All	13788	0	12704	75	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 3.

All (75) 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:D:44:ALA:HB3	1:D:68:HIS:HE2	1.49	0.76
1:D:64:LEU:HD22	1:D:68:HIS:ND1	2.02	0.74
1:B:196:GLU:HG2	1:B:201:PRO:HG3	1.68	0.74
1:D:35:LEU:HD23	1:D:88:ILE:HD12	1.76	0.67
1:B:127:VAL:CG2	1:B:147:ARG:HB2	2.25	0.65
1:D:44:ALA:HB3	1:D:68:HIS:NE2	2.10	0.65
1:C:46:TYR:HA	1:C:68:HIS:HA	1.79	0.65
1:C:196:GLU:HG2	1:C:201:PRO:HG3	1.79	0.64
1:A:126:SER:HB3	1:A:265:LEU:HD13	1.80	0.61
1:B:411:VAL:HG22	1:B:435:ILE:HD12	1.83	0.60
1:C:31:LYS:O	1:C:34:GLU:HG2	2.02	0.60
1:D:276:GLU:HG3	1:D:281:LYS:HD3	1.85	0.58
1:B:410:LYS:HE2	1:B:435:ILE:HG22	1.86	0.58
1:C:32:PRO:HB3	1:C:87:GLY:HA3	1.88	0.55
1:A:6:HIS:ND1	1:A:115:THR:OG1	2.39	0.54
1:D:15:GLY:O	1:D:45:PRO:HB3	2.09	0.53
1:C:19:GLY:HA3	1:C:69:VAL:HG21	1.91	0.52
1:C:193:VAL:HG12	1:C:224:LEU:HB2	1.92	0.52
1:B:127:VAL:HG21	1:B:147:ARG:HB2	1.90	0.52
1:D:168:VAL:HG11	1:D:243:VAL:HG21	1.92	0.50
1:A:193:VAL:HG13	1:A:227:THR:HG22	1.93	0.50
1:D:266:THR:HG22	1:D:299:VAL:HG13	1.94	0.50
1:C:344:TYR:CE2	1:C:346:PRO:HA	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:174:TYR:CD2	1:D:200:MET:HE1	2.48	0.49
1:C:7:PHE:O	1:C:116:TYR:HA	2.13	0.48
1:C:173:GLY:HA3	2:C:601:NAD:H52A	1.95	0.48
1:A:127:VAL:CG2	1:A:147:ARG:HB2	2.43	0.48
1:B:65:PRO:HB2	1:B:79:LEU:CD2	2.43	0.48
1:B:19:GLY:HA3	1:B:69:VAL:HG21	1.97	0.47
1:A:175:ILE:O	1:A:179:LEU:HB2	2.15	0.47
1:B:7:PHE:O	1:B:116:TYR:HA	2.14	0.47
1:A:174:TYR:CZ	1:A:200:MET:HE1	2.50	0.47
1:A:163:LYS:HD2	4:A:811:HOH:O	2.14	0.46
1:C:342:TYR:CE2	1:C:344:TYR:HB2	2.49	0.46
1:D:49:PRO:HA	3:D:602:FDA:N5	2.30	0.46
1:D:101:LEU:HD23	1:D:290:VAL:HG21	1.96	0.46
1:D:233:ASP:HB2	1:D:242:ALA:HB3	1.97	0.46
1:D:168:VAL:HG13	1:D:256:VAL:HG13	1.96	0.46
1:D:40:LYS:HG3	3:D:602:FDA:C4A	2.46	0.46
1:A:128:ILE:HD11	1:A:263:ARG:HG3	1.98	0.45
1:D:57:PHE:HB2	1:D:61:ALA:HA	1.99	0.45
1:C:264:PRO:HB3	1:C:299:VAL:HA	1.99	0.45
1:B:344:TYR:CE2	1:B:346:PRO:HA	2.52	0.45
1:A:17:ALA:HB1	1:A:296:ILE:HD12	1.99	0.45
1:B:304:LEU:HD11	1:B:346:PRO:HG3	1.99	0.44
1:D:19:GLY:HA3	1:D:69:VAL:HG21	1.98	0.44
1:A:197:PRO:O	1:A:225:LYS:HD2	2.17	0.44
1:A:215:TYR:CD1	1:A:355:LEU:HD11	2.53	0.44
1:B:329:ALA:HA	1:B:339:VAL:HG21	2.00	0.44
1:C:136:GLN:HB2	4:C:716:HOH:O	2.18	0.44
1:D:40:LYS:HG3	3:D:602:FDA:N3A	2.33	0.43
1:D:168:VAL:HG11	1:D:243:VAL:CG2	2.48	0.43
1:B:127:VAL:HG11	1:B:145:TYR:HB3	1.99	0.43
1:D:46:TYR:HE2	1:D:51:LEU:HD11	1.83	0.43
1:D:264:PRO:HB3	1:D:299:VAL:HA	1.99	0.43
1:A:19:GLY:HA3	1:A:69:VAL:HG21	1.99	0.43
1:D:7:PHE:CD1	1:D:36:ALA:HB2	2.54	0.43
1:A:7:PHE:O	1:A:116:TYR:HA	2.19	0.42
1:A:49:PRO:HA	3:A:602:FDA:N5	2.34	0.42
1:C:57:PHE:CD1	1:C:185:ILE:HD11	2.54	0.42
1:B:197:PRO:O	1:B:225:LYS:HD2	2.19	0.42
1:B:342:TYR:CE2	1:B:344:TYR:HB2	2.54	0.42
1:C:304:LEU:CD2	1:C:311:ARG:HG3	2.49	0.42
1:A:46:TYR:HA	1:A:68:HIS:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:391:LYS:HG3	1:C:393:PHE:HE2	1.85	0.41
1:A:174:TYR:HB3	2:A:601:NAD:C4N	2.50	0.41
1:B:331:LYS:HA	1:B:331:LYS:HD2	1.89	0.41
1:A:245:LEU:HD12	1:A:249:LYS:HB3	2.02	0.41
1:D:364:VAL:O	1:D:386:TRP:HB3	2.20	0.41
1:B:59:GLN:O	1:B:60:ASN:HB3	2.19	0.41
1:B:65:PRO:O	1:B:77:ARG:HD3	2.20	0.41
1:D:44:ALA:HB3	1:D:68:HIS:CD2	2.55	0.41
1:C:127:VAL:CG2	1:C:147:ARG:HB2	2.51	0.41
1:A:347:TYR:CD1	1:A:347:TYR:C	3.00	0.40
1:B:200:MET:HE3	1:B:200:MET:HB3	1.78	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	430/435 (99%)	413 (96%)	17 (4%)	0		
1	B	430/435 (99%)	416 (97%)	14 (3%)	0		
1	C	433/435 (100%)	417 (96%)	15 (4%)	1 (0%)		
1	D	425/435 (98%)	402 (95%)	23 (5%)	0		
All	All	1718/1740 (99%)	1648 (96%)	69 (4%)	1 (0%)		

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	163	LYS

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	331/341 (97%)	326 (98%)	5 (2%)	57	75
1	B	331/341 (97%)	326 (98%)	5 (2%)	57	75
1	C	323/341 (95%)	314 (97%)	9 (3%)	38	58
1	D	288/341 (84%)	278 (96%)	10 (4%)	32	50
All	All	1273/1364 (93%)	1244 (98%)	29 (2%)	44	64

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	132	ASP
1	A	304	LEU
1	A	335	SER
1	A	369	LEU
1	B	8	LYS
1	B	71	VAL
1	B	98	LYS
1	B	245	LEU
1	B	368	ILE
1	C	105	THR
1	C	184	LYS
1	C	185	ILE
1	C	237	ASN
1	C	316	VAL
1	C	333	LYS
1	C	335	SER
1	C	369	LEU
1	C	424	LEU
1	D	12	LEU
1	D	56	LEU
1	D	68	HIS
1	D	196	GLU
1	D	228	LEU

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Mol	Chain	Res	Type
1	D	247	ASP
1	D	251	LEU
1	D	254	ASP
1	D	350	SER
1	D	368	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	GLN
1	B	160	GLN
1	B	186	ASN
1	B	415	GLN
1	C	68	HIS
1	C	142	ASN
1	C	415	GLN
1	D	235	ASN
1	D	318	HIS

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](#)

8 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	FDA	C	602	-	57,58,58	0.33	0	78,89,89	0.54	1 (1%)
2	NAD	C	601	-	46,48,48	0.34	0	64,73,73	0.72	2 (3%)
3	FDA	B	602	-	57,58,58	0.37	0	78,89,89	0.48	0
3	FDA	A	602	-	57,58,58	0.33	0	78,89,89	0.45	0
2	NAD	A	601	-	46,48,48	0.39	0	64,73,73	0.67	2 (3%)
3	FDA	D	602	-	57,58,58	0.28	0	78,89,89	0.45	0
2	NAD	D	601	-	46,48,48	0.41	0	64,73,73	0.59	0
2	NAD	B	601	-	46,48,48	0.40	0	64,73,73	0.78	2 (3%)

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	FDA	C	602	-	-	0/34/50/50	0/6/6/6
2	NAD	C	601	-	-	3/30/62/62	0/5/5/5
3	FDA	B	602	-	-	2/34/50/50	0/6/6/6
3	FDA	A	602	-	-	1/34/50/50	0/6/6/6
2	NAD	A	601	-	-	8/30/62/62	0/5/5/5
3	FDA	D	602	-	-	2/34/50/50	0/6/6/6
2	NAD	D	601	-	-	9/30/62/62	0/5/5/5
2	NAD	B	601	-	-	8/30/62/62	0/5/5/5

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	NAD	C2N-N1N-C1D	-2.86	112.83	119.13
2	C	601	NAD	C2N-N1N-C1D	-2.73	113.11	119.13
2	C	601	NAD	C6N-N1N-C1D	2.69	125.00	119.73
2	B	601	NAD	C6N-N1N-C1D	2.60	124.82	119.73
2	A	601	NAD	C2N-N1N-C1D	-2.08	114.53	119.13
2	A	601	NAD	O5B-PA-O1A	2.05	117.08	108.94
3	C	602	FDA	O2A-PA-O1A	2.05	121.97	112.44

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	NAD	C5B-O5B-PA-O1A
2	B	601	NAD	C2B-C1B-N9A-C8A
2	B	601	NAD	O4D-C1D-N1N-C2N
2	D	601	NAD	C5B-O5B-PA-O1A
2	D	601	NAD	O4D-C1D-N1N-C2N
2	A	601	NAD	C2B-C1B-N9A-C8A
2	D	601	NAD	C2B-C1B-N9A-C8A
3	B	602	FDA	O4B-C4B-C5B-O5B
3	B	602	FDA	C3B-C4B-C5B-O5B
2	A	601	NAD	C2B-C1B-N9A-C4A
2	B	601	NAD	C2B-C1B-N9A-C4A
2	A	601	NAD	PN-O3-PA-O5B
2	B	601	NAD	PN-O3-PA-O5B
2	D	601	NAD	PN-O3-PA-O5B
2	A	601	NAD	PA-O3-PN-O1N
2	B	601	NAD	PA-O3-PN-O1N
2	C	601	NAD	PA-O3-PN-O1N
2	D	601	NAD	PA-O3-PN-O1N
2	A	601	NAD	C5B-O5B-PA-O1A
3	D	602	FDA	O4B-C4B-C5B-O5B
2	C	601	NAD	PN-O3-PA-O1A
2	D	601	NAD	C2B-C1B-N9A-C4A
2	D	601	NAD	O4D-C1D-N1N-C6N
2	C	601	NAD	PA-O3-PN-O2N
2	D	601	NAD	PA-O3-PN-O2N
2	D	601	NAD	O4B-C1B-N9A-C8A
3	A	602	FDA	O4B-C4B-C5B-O5B
3	D	602	FDA	C3B-C4B-C5B-O5B
2	A	601	NAD	O4B-C1B-N9A-C8A
2	A	601	NAD	PA-O3-PN-O2N
2	B	601	NAD	PA-O3-PN-O2N
2	A	601	NAD	O4B-C4B-C5B-O5B
2	B	601	NAD	O4B-C4B-C5B-O5B

There are no ring outliers.

4 monomers are involved in 6 short contacts:

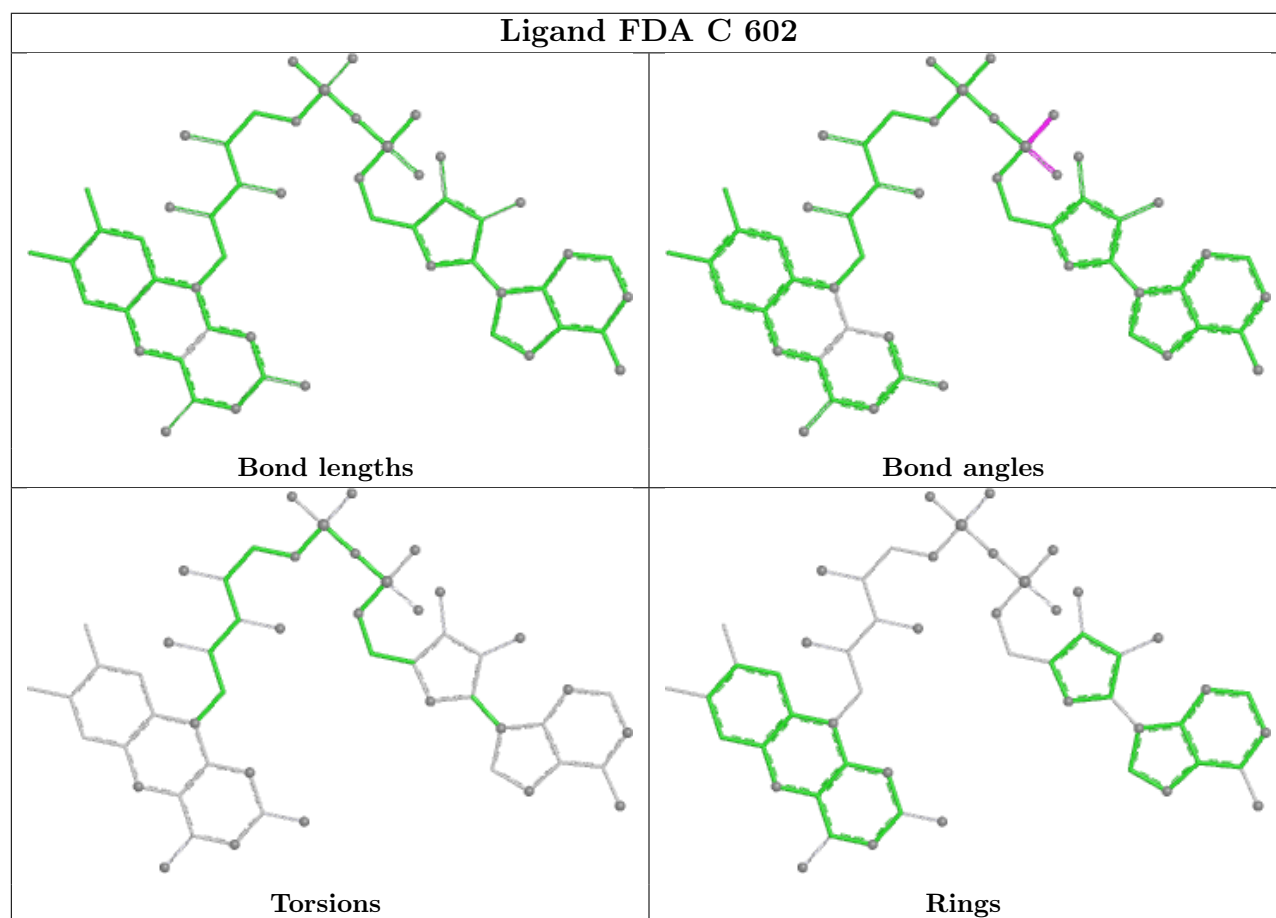
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	NAD	1	0
3	A	602	FDA	1	0

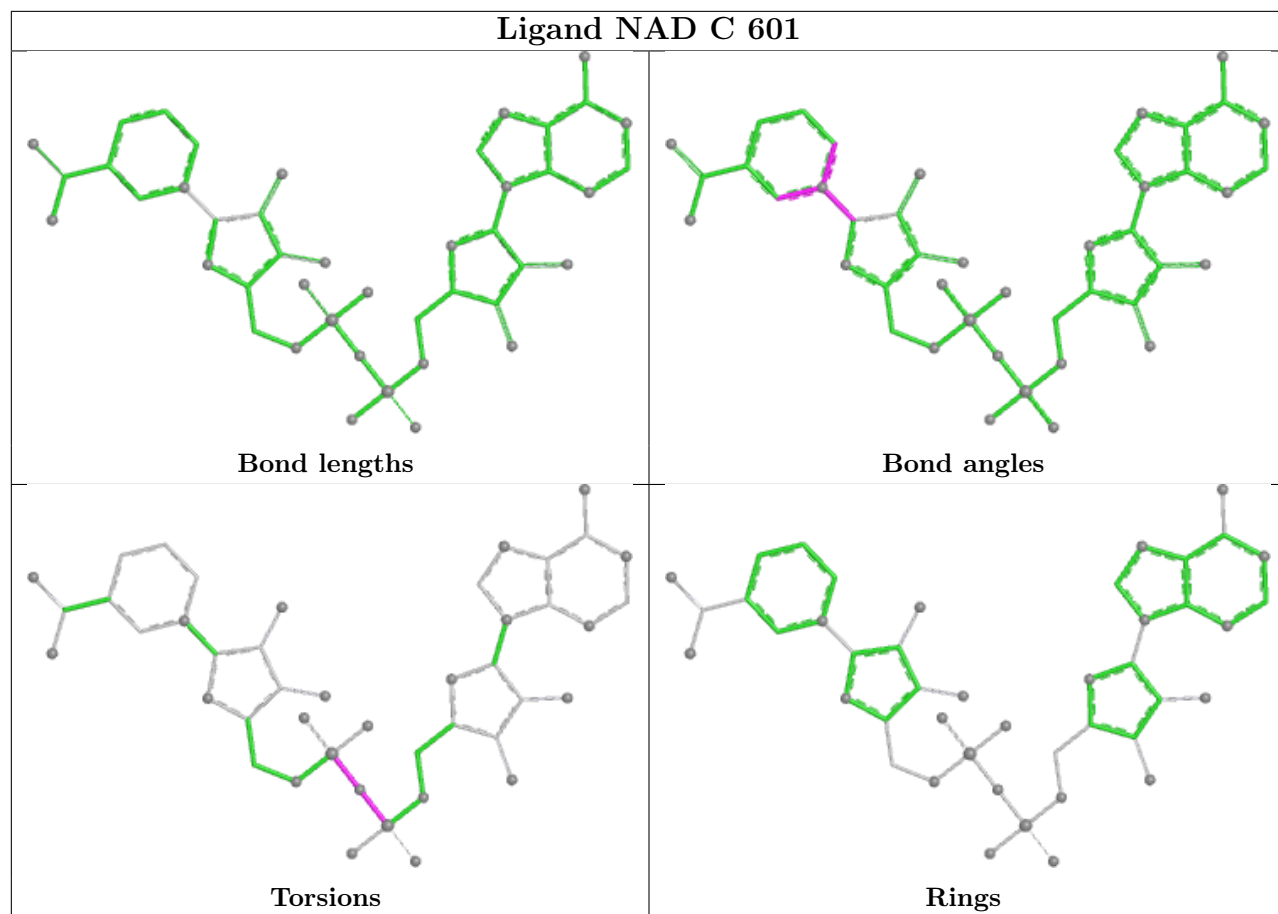
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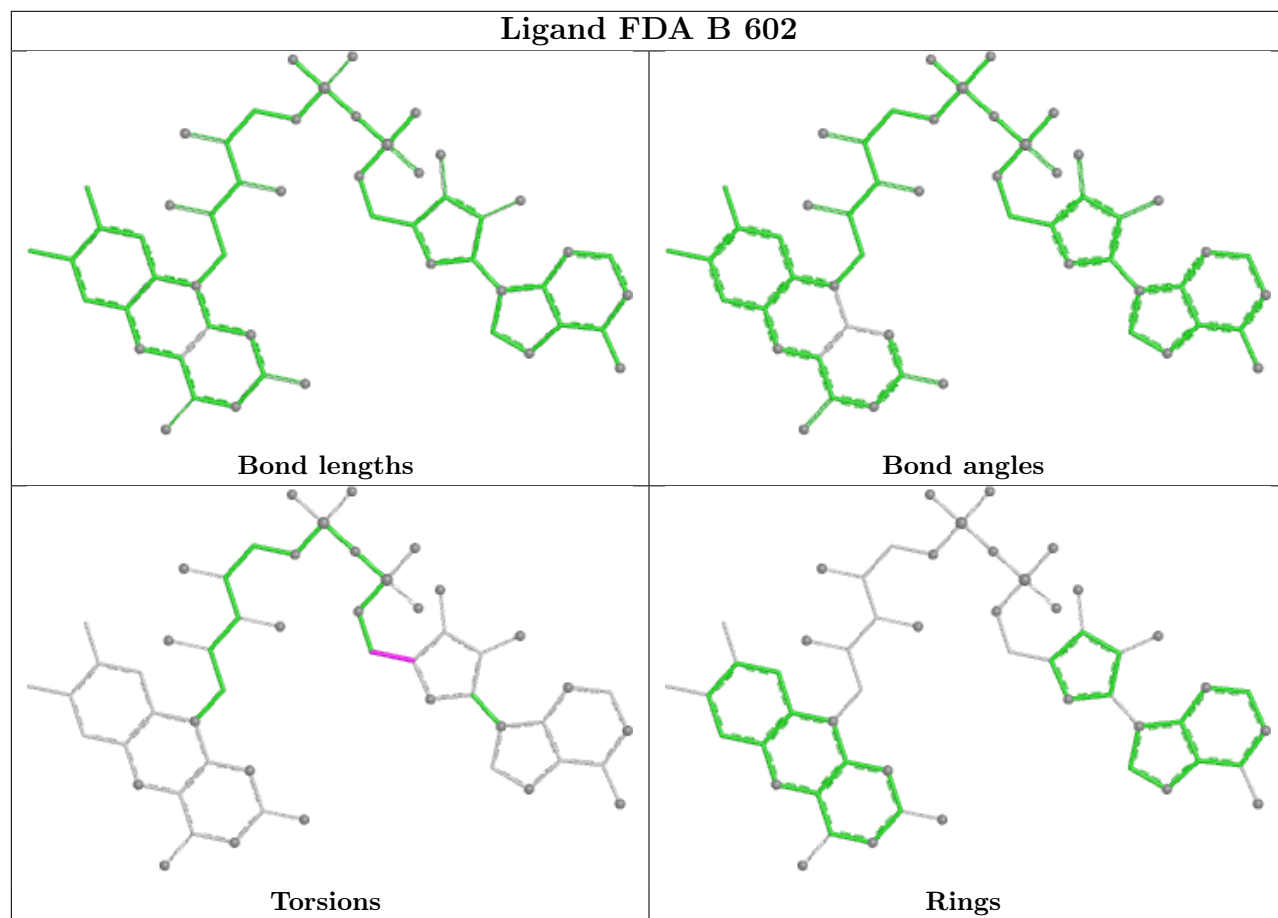
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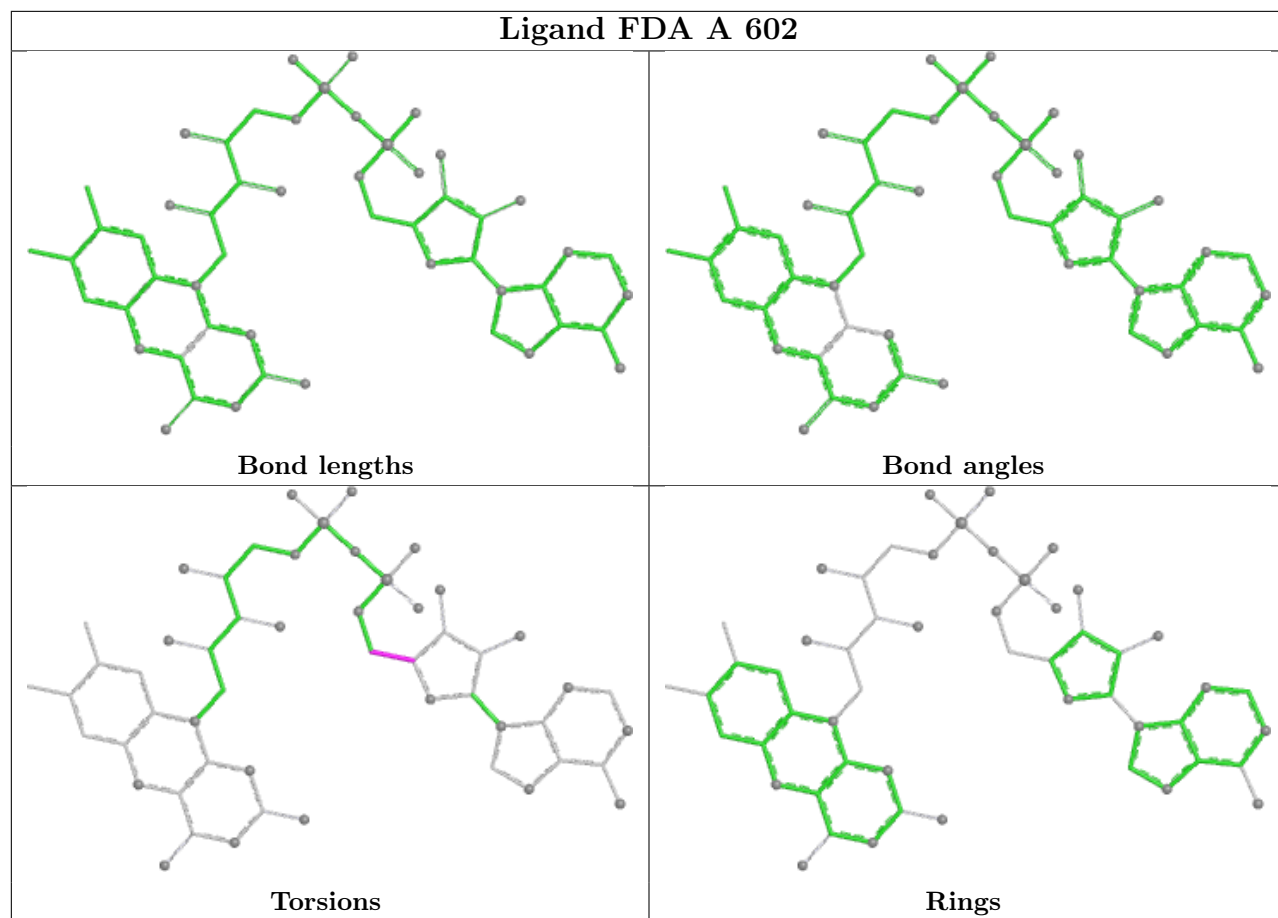
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	NAD	1	0
3	D	602	FDA	3	0

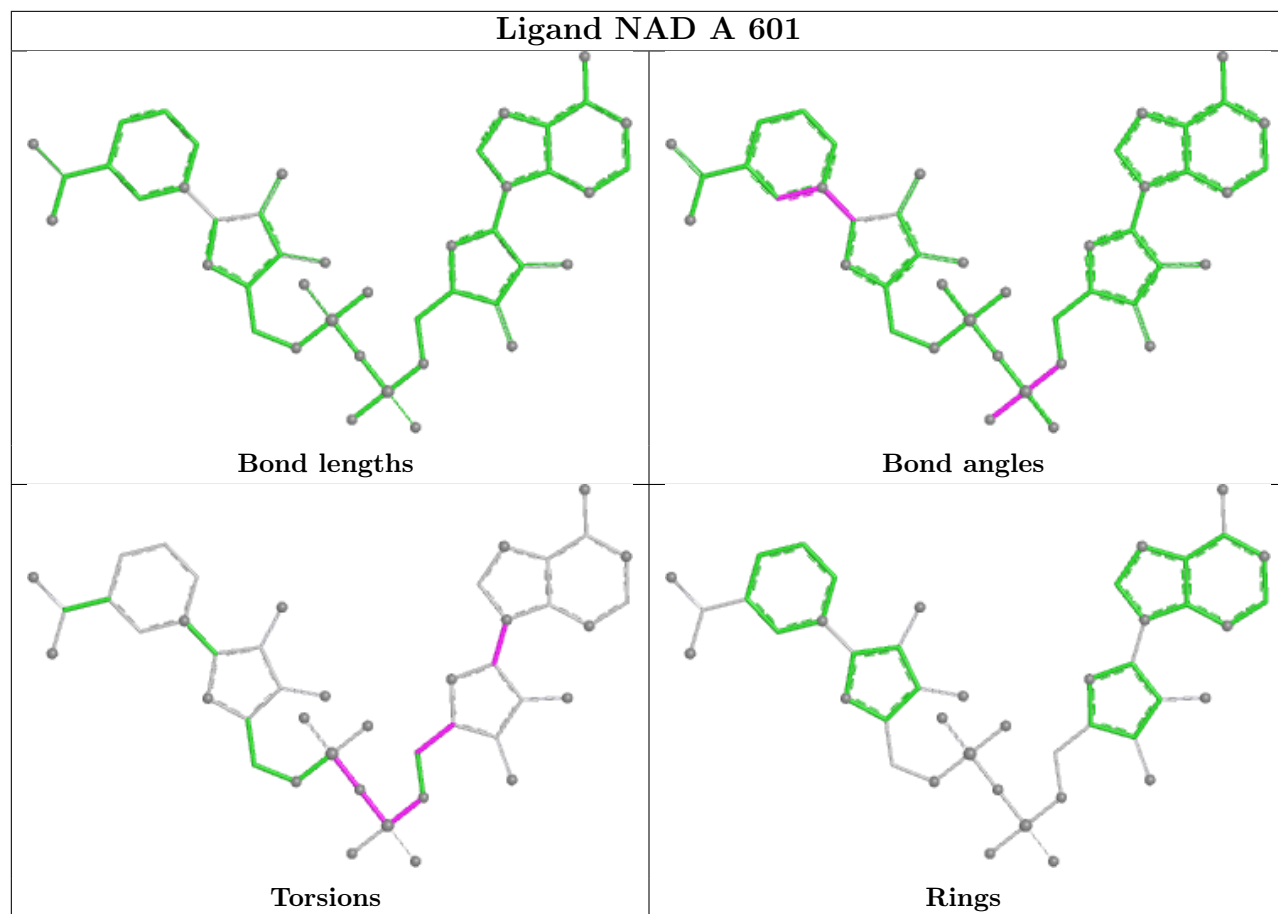
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

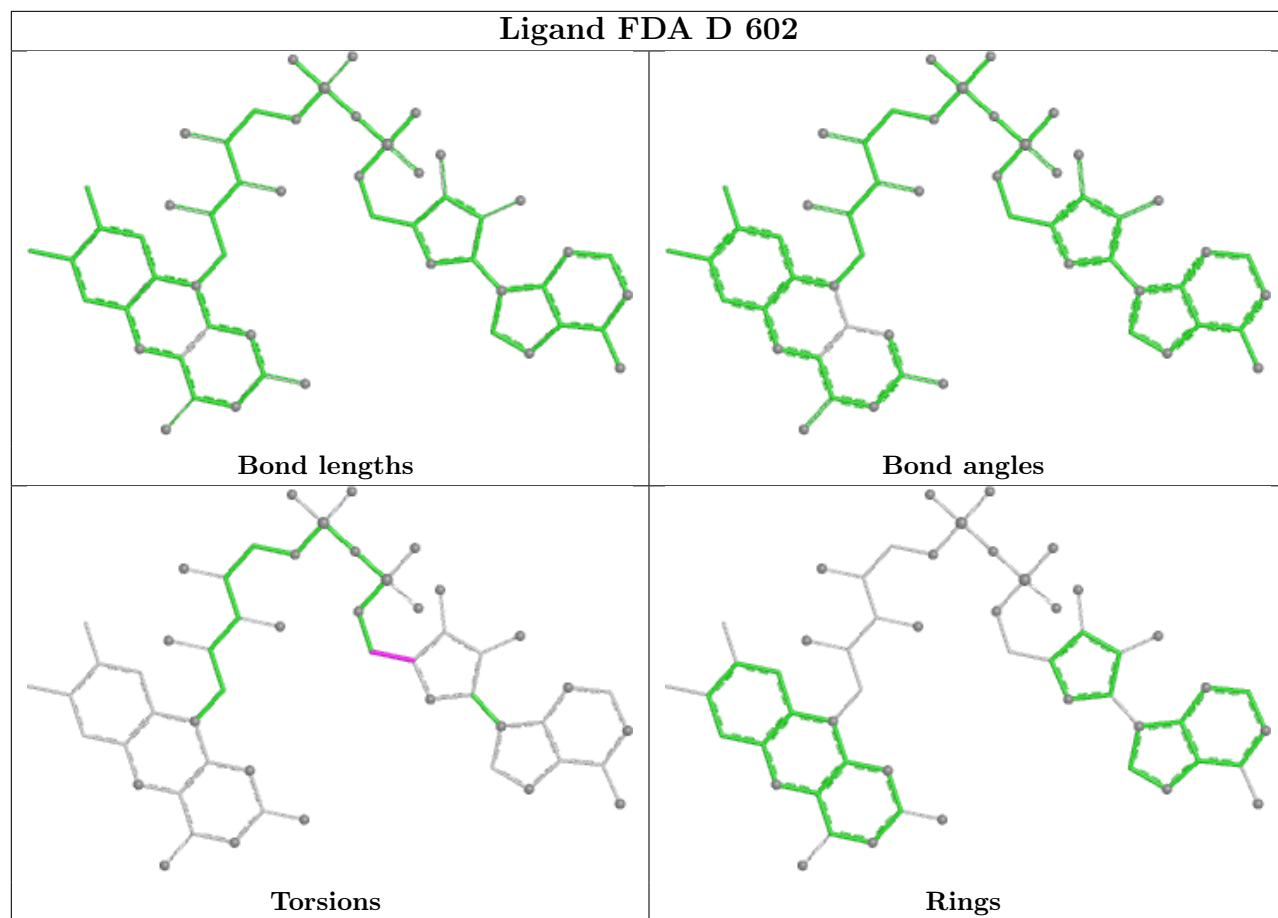


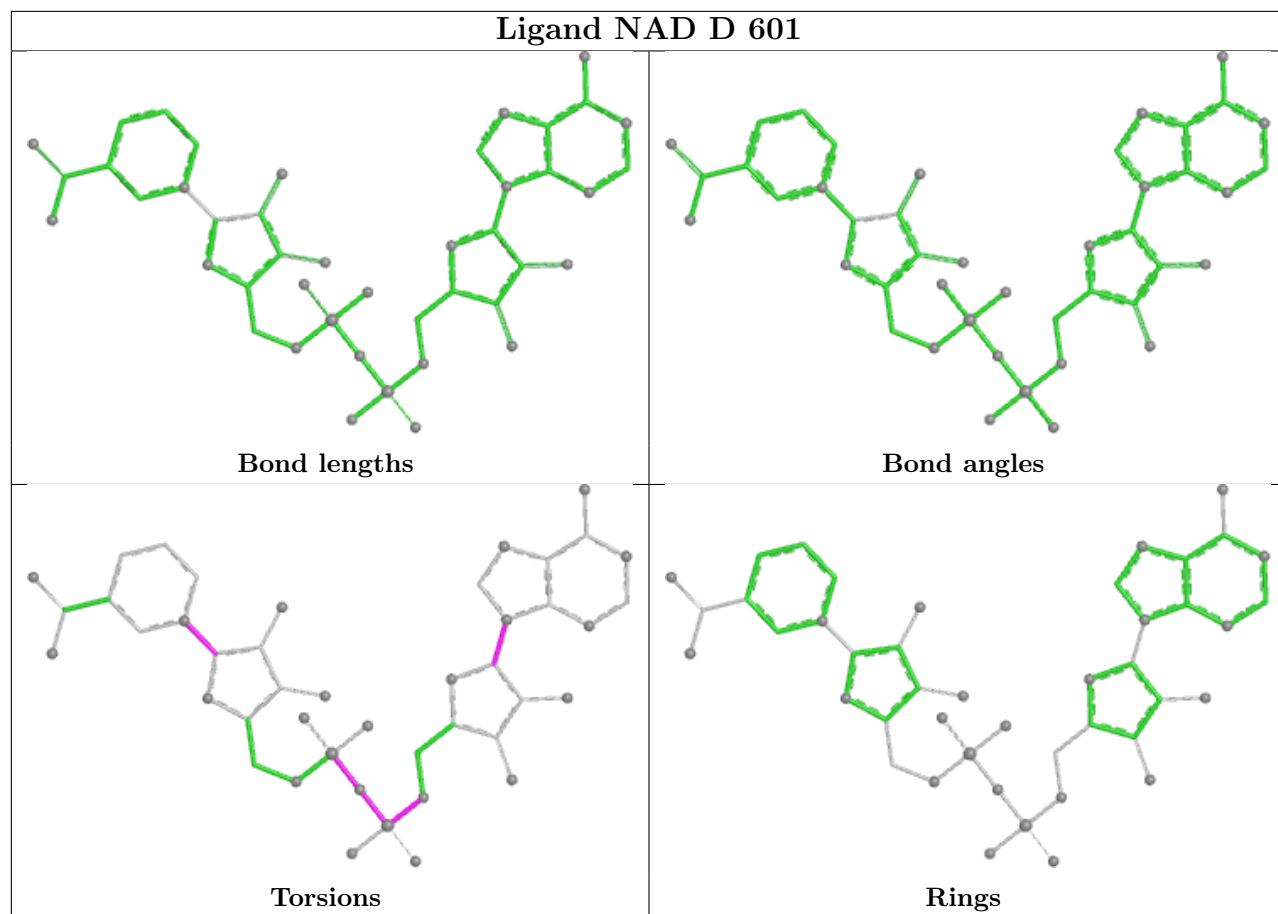


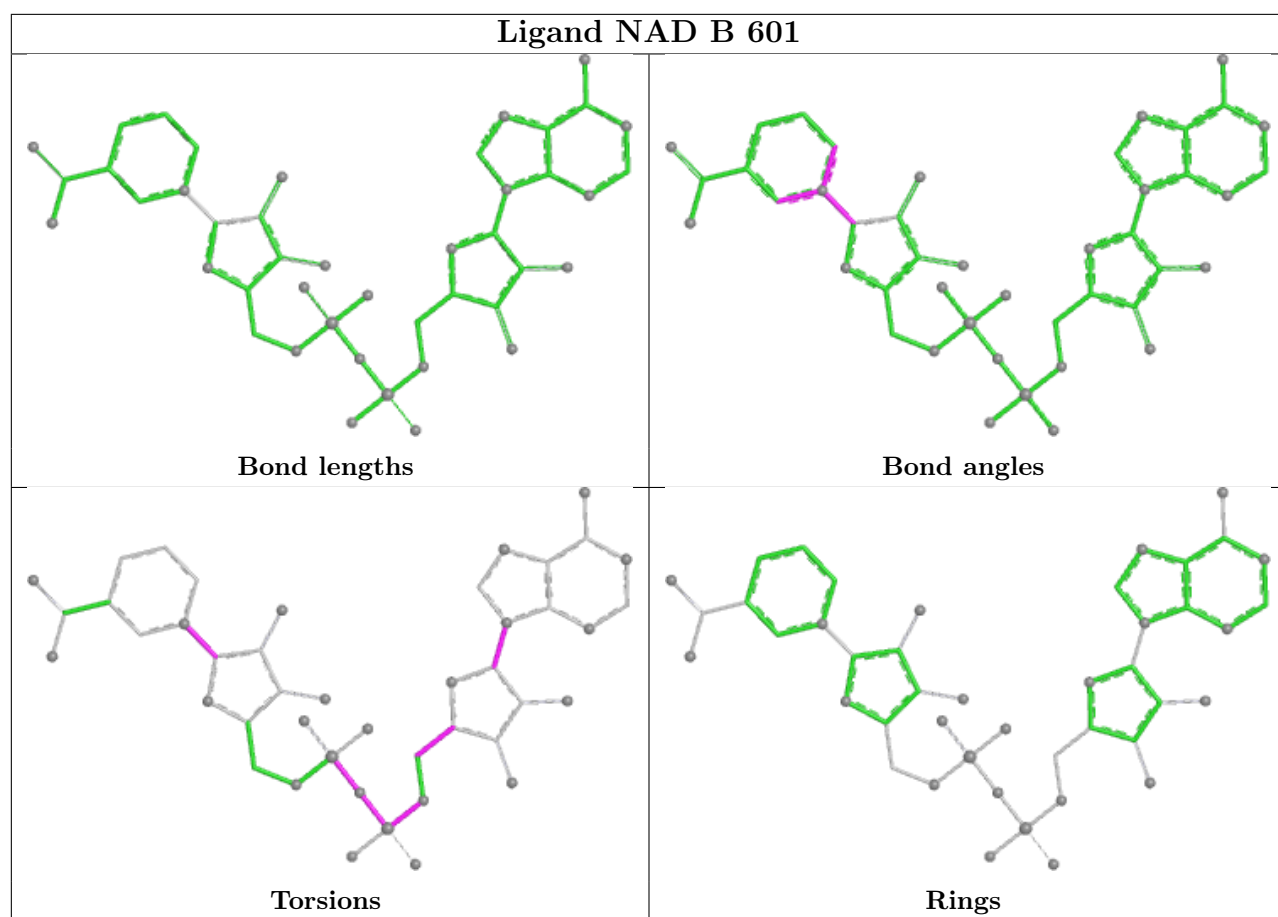












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	432/435 (99%)	-0.29	2 (0%) 87 86	23, 32, 50, 70	0
1	B	432/435 (99%)	-0.24	0 100 100	22, 33, 51, 59	0
1	C	435/435 (100%)	-0.03	3 (0%) 84 83	22, 38, 57, 77	0
1	D	429/435 (98%)	0.77	31 (7%) 21 20	38, 58, 76, 83	0
All	All	1728/1740 (99%)	0.05	36 (2%) 63 62	22, 38, 70, 83	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	435	ILE	4.0
1	A	435	ILE	3.3
1	D	90	LEU	2.9
1	D	435	ILE	2.8
1	D	84	SER	2.7
1	D	87	GLY	2.7
1	D	23	ARG	2.6
1	D	377	SER	2.6
1	D	245	LEU	2.5
1	D	80	PRO	2.5
1	D	27	LYS	2.4
1	D	76	GLU	2.4
1	C	421	ILE	2.4
1	D	250	VAL	2.3
1	D	232	PHE	2.3
1	D	56	LEU	2.3
1	D	64	LEU	2.3
1	D	30	VAL	2.3
1	D	31	LYS	2.3
1	D	243	VAL	2.3
1	D	72	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	133	PHE	2.1
1	D	57	PHE	2.1
1	D	70	CYS	2.1
1	D	26	ALA	2.1
1	D	115	THR	2.1
1	D	338	SER	2.1
1	D	378	SER	2.1
1	D	35	LEU	2.1
1	D	237	ASN	2.1
1	D	353	PHE	2.1
1	D	33	GLY	2.0
1	C	429	LEU	2.0
1	D	251	LEU	2.0
1	D	86	LYS	2.0
1	D	166	LYS	2.0

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](#)

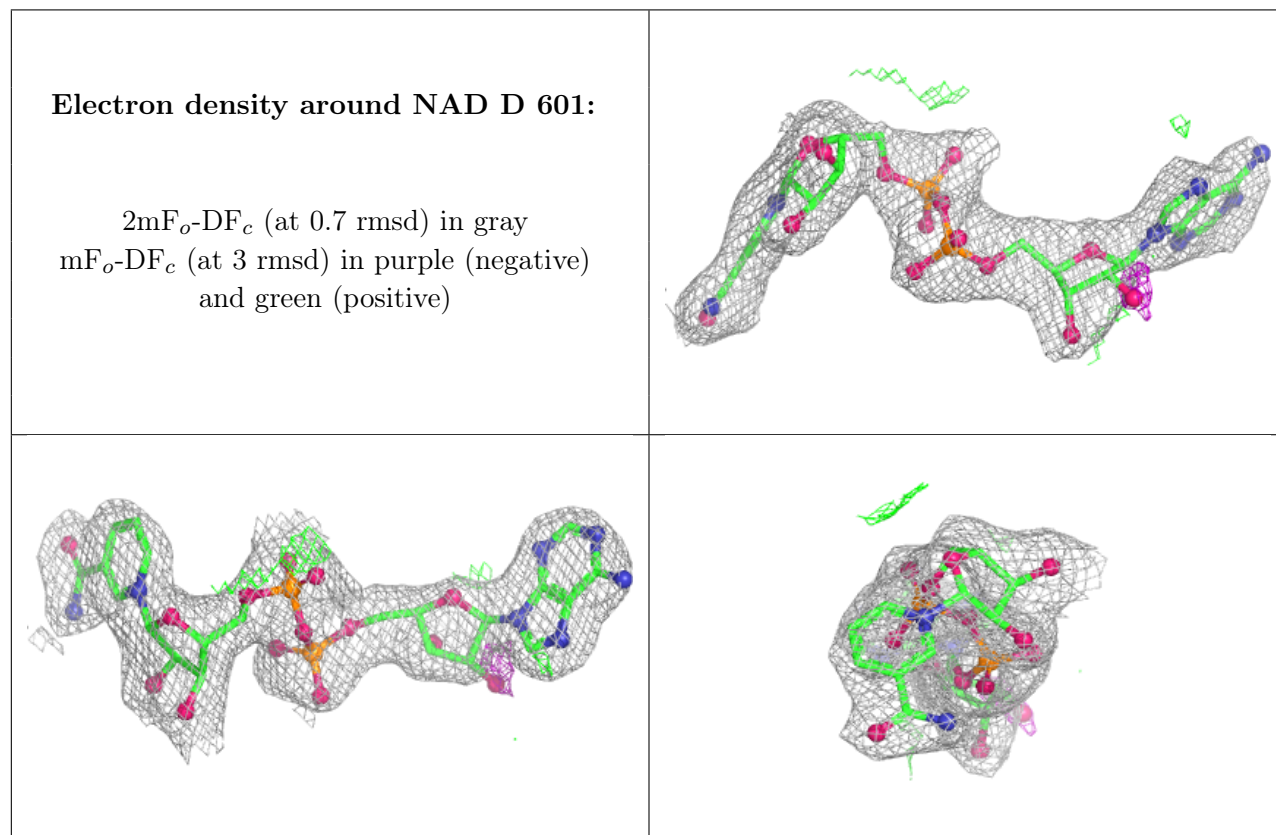
There are no oligosaccharides in this entry.

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.

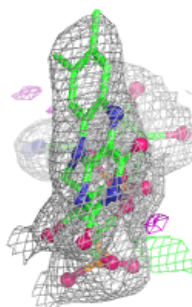
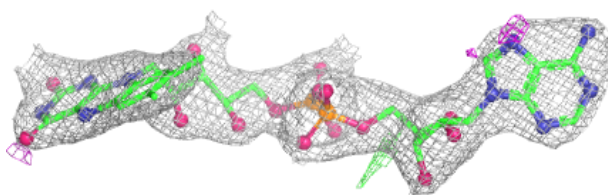
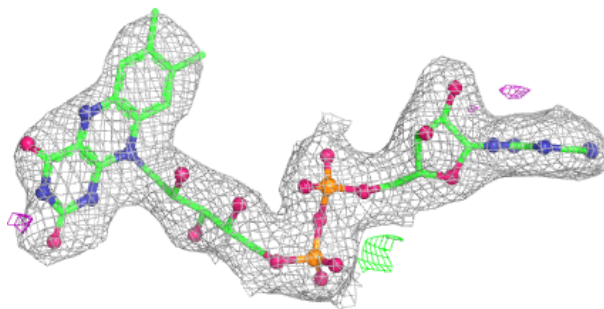
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAD	D	601	44/44	0.93	0.09	36,43,54,54	0
3	FDA	D	602	53/53	0.94	0.08	43,46,55,55	0
2	NAD	C	601	44/44	0.95	0.08	28,32,54,55	0
2	NAD	B	601	44/44	0.97	0.06	20,32,39,40	0
3	FDA	C	602	53/53	0.97	0.05	25,27,31,31	0
2	NAD	A	601	44/44	0.97	0.06	23,28,40,41	0
3	FDA	A	602	53/53	0.98	0.05	23,25,27,28	0
3	FDA	B	602	53/53	0.98	0.04	22,25,31,32	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.

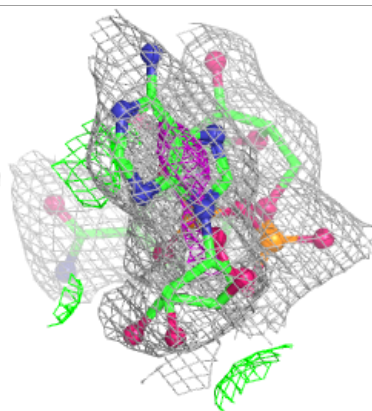
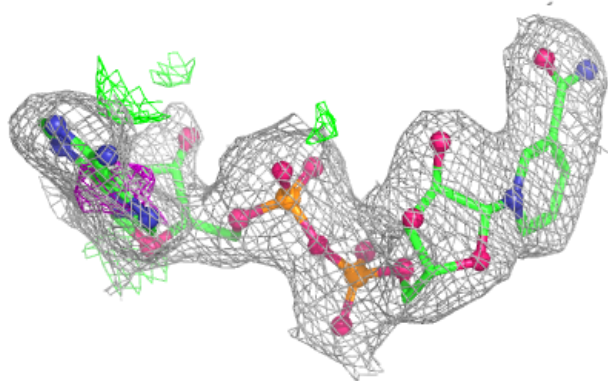
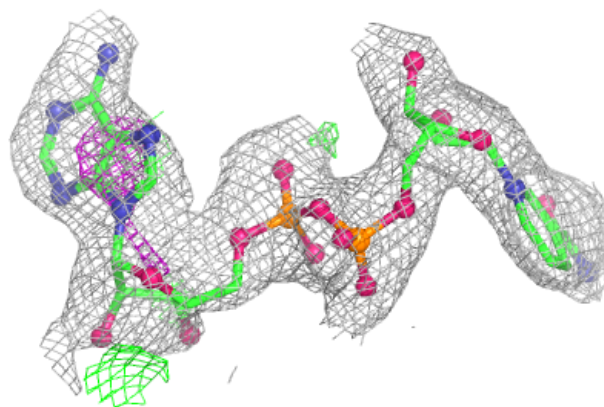


Electron density around FDA D 602:

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

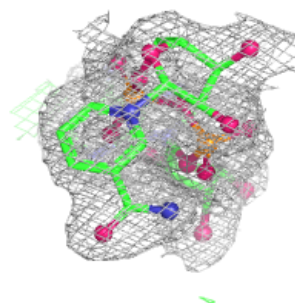
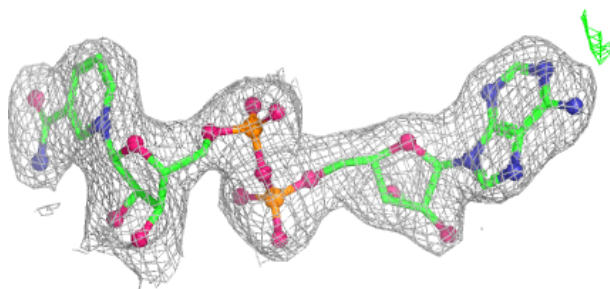
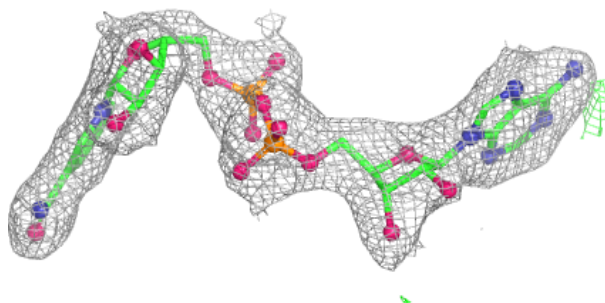
**Electron density around NAD C 601:**

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

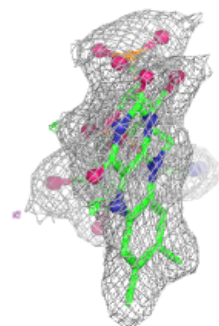
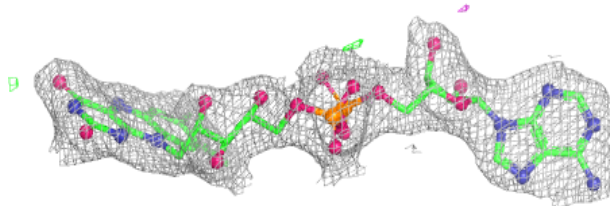
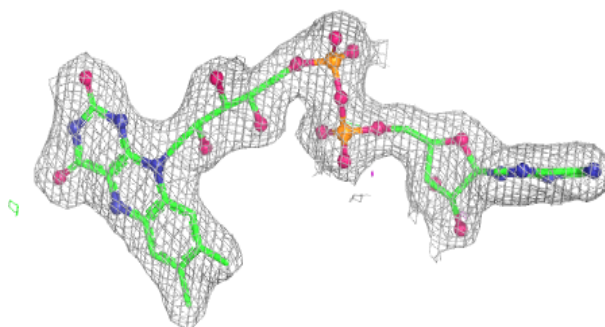


Electron density around NAD B 601:

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

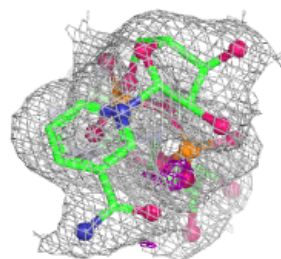
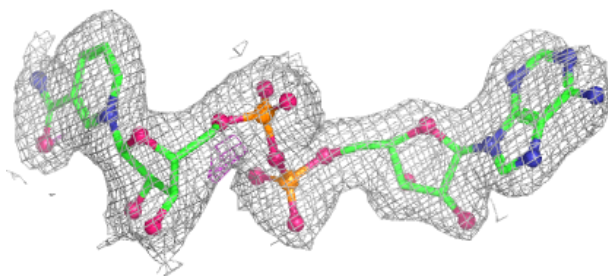
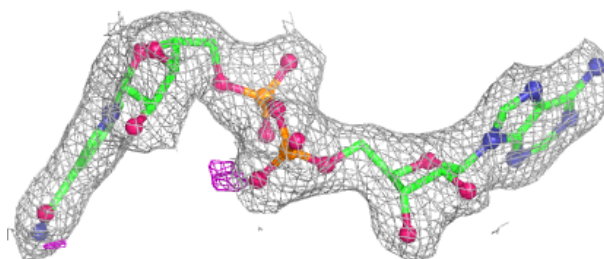
**Electron density around FDA C 602:**

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

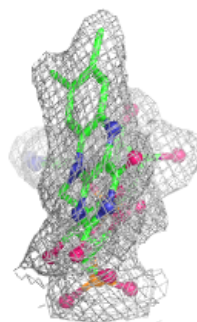
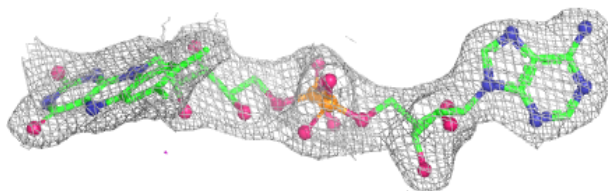


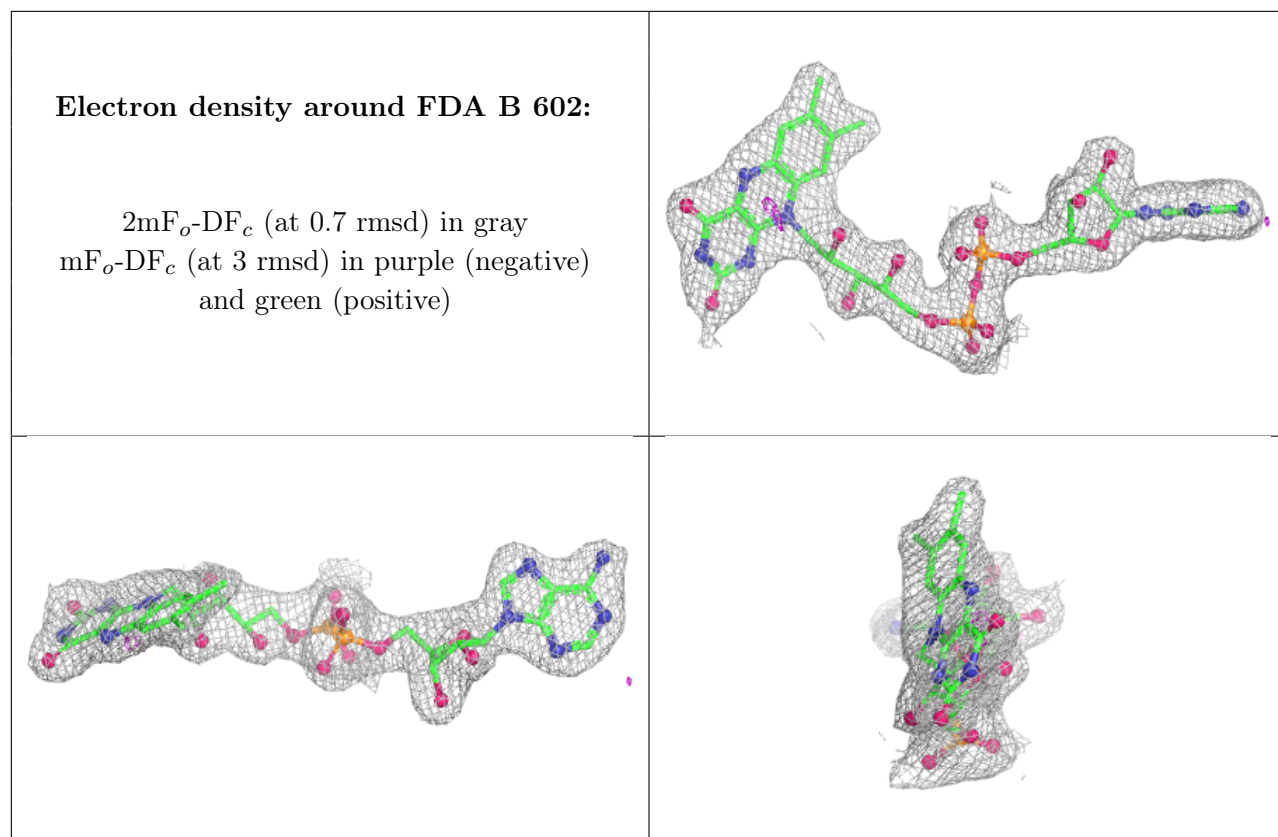
Electron density around NAD A 601:

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

**Electron density around FDA A 602:**

$2mF_o-DF_c$ (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.