



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 10:14 PM UTC

PDB ID : 1KR2 / pdb_00001kr2
Title : CRYSTAL STRUCTURE OF HUMAN NMN/NAMN ADENYLYL TRANSFERASE COMPLEXED WITH TIAZOFURIN ADENINE DINUCLEOTIDE (TAD)
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Deposited on : 2002-01-08
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)	:	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

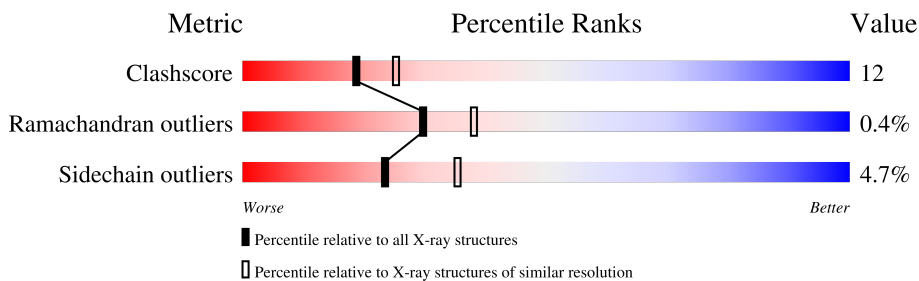
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	279	65% 16% • 16%
1	B	279	65% 16% • 16%
1	C	279	61% 18% • 16%
1	D	279	66% 15% • 16%
1	E	279	65% 16% • 17%
1	F	279	61% 19% • 17%

2 Entry composition ⓘ

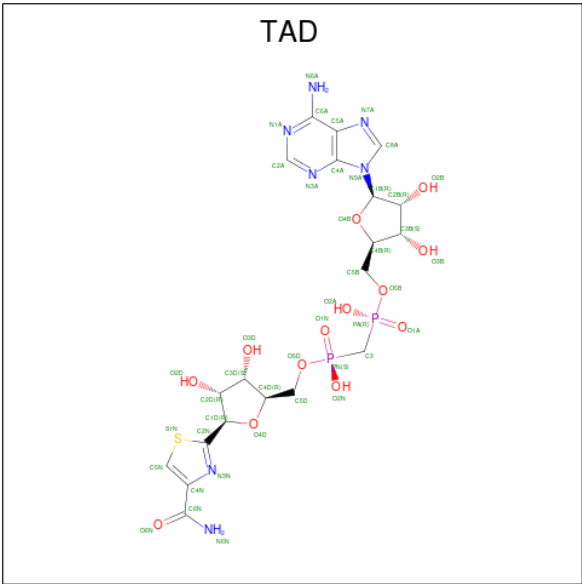
There are 3 unique types of molecules in this entry. The entry contains 12192 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1875	1195	331	343	6			
1	B	233	Total	C	N	O	S	0	0	0
			1875	1195	331	343	6			
1	C	233	Total	C	N	O	S	0	0	0
			1875	1195	331	343	6			
1	D	233	Total	C	N	O	S	0	0	0
			1875	1195	331	343	6			
1	E	232	Total	C	N	O	S	0	0	0
			1866	1190	330	340	6			
1	F	231	Total	C	N	O	S	0	0	0
			1861	1187	329	339	6			

- Molecule 2 is BETA-METHYLENE-THIAZOLE-4-CARBOXYAMIDE-ADENINE DINUCLEOTIDE (CCD ID: TAD) (formula: C₂₀H₂₇N₇O₁₃P₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 43	C 20	N 7	O 13	P 2	S 1	0	0
2	B	1	Total 43	C 20	N 7	O 13	P 2	S 1	0	0
2	C	1	Total 43	C 20	N 7	O 13	P 2	S 1	0	0
2	D	1	Total 43	C 20	N 7	O 13	P 2	S 1	0	0
2	E	1	Total 43	C 20	N 7	O 13	P 2	S 1	0	0
2	F	1	Total 43	C 20	N 7	O 13	P 2	S 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	155	Total 155	O 155	0	0
3	B	128	Total 128	O 128	0	0
3	C	168	Total 168	O 168	0	0
3	D	103	Total 103	O 103	0	0
3	E	83	Total 83	O 83	0	0
3	F	70	Total 70	O 70	0	0

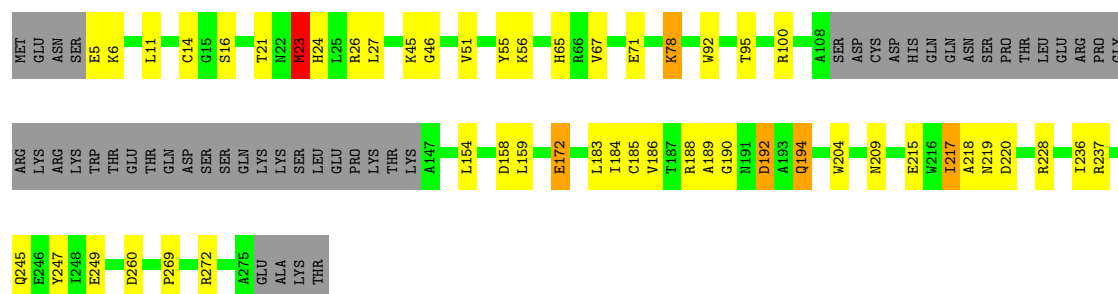
3 Residue-property plots

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

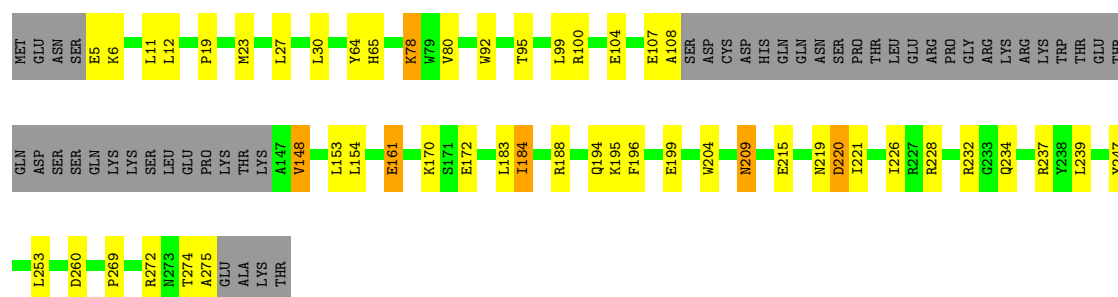
• Molecule 1: NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE

Chain A: 



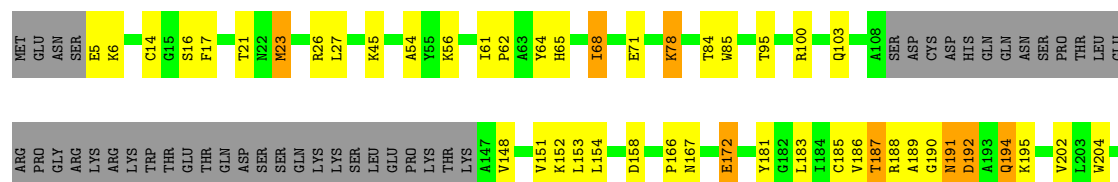
• Molecule 1: NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE

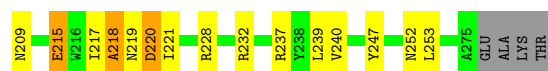
Chain B: 



• Molecule 1: NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE

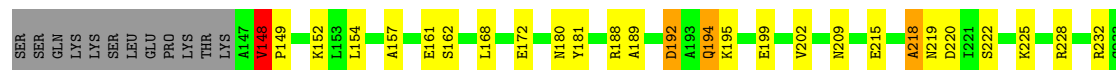
Chain C: 





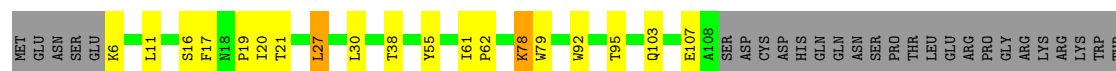
• Molecule 1: NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE

Chain D: 66% 15% 16%



• Molecule 1: NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE

Chain E: 65% 16% 17%



• Molecule 1: NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE

Chain F: 61% 19% 17%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	131.48Å 89.39Å 137.75Å 90.00° 117.42° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.30)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.203 , 0.231	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12192	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

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: TAD

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	0.83	1/1912 (0.1%)	1.04	5/2588 (0.2%)
1	B	0.84	3/1912 (0.2%)	1.09	9/2588 (0.3%)
1	C	0.84	1/1912 (0.1%)	1.07	9/2588 (0.3%)
1	D	0.80	2/1912 (0.1%)	1.10	8/2588 (0.3%)
1	E	0.66	0/1903	1.06	12/2576 (0.5%)
1	F	0.68	1/1898 (0.1%)	1.05	4/2569 (0.2%)
All	All	0.78	8/11449 (0.1%)	1.07	47/15497 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	23	MET	SD-CE	5.98	1.94	1.79
1	F	23	MET	SD-CE	5.73	1.93	1.79
1	B	23	MET	SD-CE	5.61	1.93	1.79
1	C	202	VAL	CA-CB	5.56	1.61	1.54
1	A	23	MET	CG-SD	5.46	1.94	1.80
1	B	226	ILE	CA-CB	5.19	1.60	1.54
1	B	80	VAL	CA-CB	5.15	1.60	1.54
1	D	51	VAL	CA-CB	5.12	1.60	1.54

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	253	LEU	N-CA-C	15.25	127.38	111.07
1	E	253	LEU	N-CA-C	13.32	125.32	111.07
1	D	253	LEU	N-CA-C	11.65	123.97	111.28
1	C	192	ASP	N-CA-C	-8.54	100.22	111.24
1	D	148	VAL	N-CA-C	8.34	126.90	108.88
1	F	192	ASP	N-CA-C	-8.06	102.18	110.97
1	F	253	LEU	N-CA-C	7.65	119.62	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	148	VAL	CA-C-N	7.26	127.31	119.90
1	E	148	VAL	C-N-CA	7.26	127.31	119.90
1	E	239	LEU	N-CA-C	-7.15	104.82	113.97
1	E	220	ASP	N-CA-C	7.03	125.78	110.80
1	A	192	ASP	N-CA-C	-7.01	103.57	111.14
1	D	148	VAL	CA-C-N	6.79	126.76	119.76
1	D	148	VAL	C-N-CA	6.79	126.76	119.76
1	C	191	ASN	N-CA-C	-6.61	105.19	113.18
1	C	56	LYS	N-CA-C	6.44	120.71	112.86
1	C	187	THR	N-CA-C	6.42	118.31	108.42
1	B	239	LEU	N-CA-C	-6.27	105.94	113.97
1	B	220	ASP	N-CA-C	6.24	119.31	109.39
1	B	148	VAL	CA-C-N	5.93	125.94	119.89
1	B	148	VAL	C-N-CA	5.93	125.94	119.89
1	B	272	ARG	N-CA-C	5.93	117.74	111.28
1	A	14	CYS	N-CA-C	-5.88	101.75	110.46
1	D	253	LEU	CA-C-N	-5.84	114.45	122.93
1	D	253	LEU	C-N-CA	-5.84	114.45	122.93
1	A	56	LYS	N-CA-C	5.84	119.99	112.86
1	A	55	TYR	N-CA-C	-5.72	105.19	111.82
1	C	239	LEU	N-CA-C	-5.70	106.48	113.50
1	E	20	ILE	N-CA-C	-5.65	102.29	109.80
1	B	170	LYS	N-CA-C	-5.58	102.02	110.28
1	D	218	ALA	N-CA-C	5.49	118.28	110.10
1	F	103	GLN	N-CA-C	-5.37	105.07	111.03
1	C	14	CYS	N-CA-C	-5.34	102.14	110.42
1	E	55	TYR	N-CA-C	-5.33	105.55	111.36
1	F	170	LYS	N-CA-C	-5.31	101.78	109.59
1	E	244	VAL	N-CA-C	-5.29	105.38	110.72
1	E	184	ILE	N-CA-C	-5.28	99.55	107.37
1	D	180	ASN	N-CA-C	5.25	119.96	113.50
1	A	272	ARG	N-CA-C	5.18	117.83	111.82
1	E	38	THR	N-CA-C	-5.17	106.23	112.54
1	C	148	VAL	CA-C-N	5.16	125.15	119.89
1	C	148	VAL	C-N-CA	5.16	125.15	119.89
1	B	148	VAL	N-CA-C	5.15	112.96	107.76
1	C	253	LEU	N-CA-C	5.14	116.88	111.28
1	E	148	VAL	N-CA-C	5.05	112.86	107.76
1	B	184	ILE	N-CA-C	-5.04	99.27	107.28
1	E	79	TRP	N-CA-C	5.03	121.11	114.12

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	1875	0	1903	59	0
1	B	1875	0	1903	43	0
1	C	1875	0	1903	65	0
1	D	1875	0	1903	40	0
1	E	1866	0	1897	33	0
1	F	1861	0	1892	44	0
2	A	43	0	25	1	0
2	B	43	0	25	1	0
2	C	43	0	25	1	0
2	D	43	0	25	5	0
2	E	43	0	25	2	0
2	F	43	0	25	2	0
3	A	155	0	0	14	0
3	B	128	0	0	11	0
3	C	168	0	0	18	0
3	D	103	0	0	6	0
3	E	83	0	0	4	0
3	F	70	0	0	4	0
All	All	12192	0	11551	273	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 12.

All (273) 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:C:23:MET:HB3	3:C:941:HOH:O	1.29	1.27
1:D:78:LYS:H	1:D:78:LYS:HD2	1.11	1.10
1:B:78:LYS:H	1:B:78:LYS:HD2	0.98	1.09
1:B:78:LYS:H	1:B:78:LYS:CD	1.69	1.03
1:E:219:ASN:HA	1:F:219:ASN:OD1	1.60	1.00
1:D:78:LYS:H	1:D:78:LYS:CD	1.76	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:183:LEU:HB3	3:C:977:HOH:O	1.64	0.96
1:B:78:LYS:HD2	1:B:78:LYS:N	1.80	0.95
1:C:78:LYS:H	1:C:78:LYS:HD2	1.28	0.94
1:C:192:ASP:HA	3:C:939:HOH:O	1.67	0.92
1:A:183:LEU:HB3	3:A:891:HOH:O	1.70	0.91
1:C:158:ASP:HB2	3:C:933:HOH:O	1.70	0.90
1:F:78:LYS:H	1:F:78:LYS:HD2	1.35	0.89
1:A:23:MET:HA	1:A:23:MET:HE2	1.56	0.88
1:D:78:LYS:HD2	1:D:78:LYS:N	1.88	0.88
1:F:78:LYS:H	1:F:78:LYS:CD	1.87	0.86
1:A:78:LYS:H	1:A:78:LYS:HD2	1.38	0.85
1:A:236:ILE:HG12	3:A:959:HOH:O	1.76	0.84
1:A:78:LYS:H	1:A:78:LYS:CD	1.90	0.83
1:A:194:GLN:HE21	1:A:194:GLN:HA	1.43	0.82
1:A:23:MET:HE3	1:A:219:ASN:HB3	1.60	0.82
2:F:814:TAD:H31	3:F:881:HOH:O	1.80	0.81
1:A:21:THR:H	1:A:24:HIS:HD2	1.27	0.81
1:C:78:LYS:H	1:C:78:LYS:CD	1.83	0.79
1:C:194:GLN:HA	1:C:194:GLN:HE21	1.48	0.79
1:D:23:MET:HG2	3:D:901:HOH:O	1.82	0.79
1:C:95:THR:H	2:C:811:TAD:H61N	1.30	0.78
1:C:5:GLU:HG3	1:C:6:LYS:H	1.48	0.78
1:F:78:LYS:HD2	1:F:78:LYS:N	1.98	0.77
1:C:189:ALA:HB1	1:C:192:ASP:HB2	1.67	0.77
1:C:188:ARG:HD2	1:C:218:ALA:HA	1.65	0.76
1:F:189:ALA:HB1	1:F:192:ASP:HB2	1.68	0.76
1:B:219:ASN:HB3	3:B:853:HOH:O	1.87	0.75
1:B:234:GLN:HG2	3:B:922:HOH:O	1.84	0.75
1:C:78:LYS:HD2	1:C:78:LYS:N	2.00	0.74
1:C:5:GLU:HB2	3:C:972:HOH:O	1.87	0.74
1:C:219:ASN:OD1	1:C:220:ASP:N	2.19	0.74
1:A:95:THR:H	2:A:809:TAD:H61N	1.36	0.74
1:C:5:GLU:HG2	3:C:966:HOH:O	1.87	0.73
1:A:245:GLN:O	1:A:249:GLU:HG2	1.89	0.73
1:C:217:ILE:HD12	3:C:935:HOH:O	1.89	0.73
1:D:30:LEU:HD11	1:D:215:GLU:CD	2.15	0.72
1:B:188:ARG:HD3	3:B:916:HOH:O	1.90	0.71
1:A:78:LYS:HD2	1:A:78:LYS:N	2.06	0.71
1:F:65:HIS:HD2	1:F:247:TYR:OH	1.73	0.71
1:B:65:HIS:HD2	1:B:247:TYR:OH	1.75	0.70
1:B:215:GLU:OE2	3:B:908:HOH:O	2.09	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:157:ALA:O	1:F:161:GLU:HG3	1.90	0.70
1:D:95:THR:H	2:D:812:TAD:H61N	1.40	0.69
1:D:202:VAL:HG23	3:D:845:HOH:O	1.92	0.69
1:F:268:ALA:HB3	1:F:269:PRO:HD3	1.74	0.69
1:F:92:TRP:CH2	1:F:269:PRO:HG3	2.26	0.69
1:E:78:LYS:H	1:E:78:LYS:HD2	1.57	0.69
1:A:71:GLU:CD	3:A:934:HOH:O	2.35	0.69
1:C:5:GLU:HG3	1:C:6:LYS:N	2.07	0.69
1:A:172:GLU:HB3	3:A:962:HOH:O	1.92	0.68
1:D:234:GLN:HG2	3:D:891:HOH:O	1.93	0.68
1:F:95:THR:H	2:F:814:TAD:H61N	1.42	0.68
1:E:11:LEU:HD22	1:E:184:ILE:HD12	1.75	0.68
1:A:23:MET:HE1	1:A:219:ASN:OD1	1.95	0.67
1:C:228:ARG:HD2	1:C:232:ARG:NH2	2.10	0.67
1:E:153:LEU:HB3	1:E:183:LEU:HD23	1.77	0.66
1:A:23:MET:HG3	3:A:932:HOH:O	1.94	0.66
1:C:219:ASN:O	1:C:220:ASP:CB	2.43	0.66
1:A:186:VAL:HG12	3:A:951:HOH:O	1.96	0.66
1:E:78:LYS:H	1:E:78:LYS:CD	2.07	0.66
1:B:95:THR:H	2:B:810:TAD:H61N	1.43	0.65
1:F:64:TYR:O	1:F:68:ILE:HG12	1.96	0.65
1:C:219:ASN:O	1:C:220:ASP:HB3	1.96	0.65
1:C:252:ASN:HB3	3:C:968:HOH:O	1.95	0.65
1:D:16:SER:H	2:D:812:TAD:H32	1.62	0.65
1:A:204:TRP:CD1	1:D:232:ARG:HD3	2.33	0.64
1:A:65:HIS:HD2	1:A:247:TYR:OH	1.81	0.64
1:A:194:GLN:HA	1:A:194:GLN:NE2	2.13	0.64
1:B:194:GLN:HE21	1:B:194:GLN:HA	1.63	0.64
1:E:220:ASP:HB3	3:E:837:HOH:O	1.96	0.64
1:A:185:CYS:HB2	3:A:891:HOH:O	1.98	0.63
1:F:188:ARG:HD3	3:F:838:HOH:O	1.97	0.63
1:B:275:ALA:HB2	3:B:917:HOH:O	1.97	0.63
1:B:64:TYR:HE1	3:B:926:HOH:O	1.82	0.63
1:C:183:LEU:HD13	3:C:977:HOH:O	1.98	0.63
1:D:194:GLN:HA	1:D:194:GLN:HE21	1.62	0.62
1:C:23:MET:SD	3:C:958:HOH:O	2.56	0.62
1:E:95:THR:H	2:E:813:TAD:H61N	1.46	0.62
1:A:92:TRP:CH2	1:A:269:PRO:HG3	2.35	0.62
1:E:228:ARG:HD2	1:E:232:ARG:NH2	2.14	0.62
1:E:92:TRP:CH2	1:E:269:PRO:HG3	2.35	0.62
1:C:189:ALA:CB	1:C:192:ASP:HB2	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ARG:O	1:B:104:GLU:HG3	2.01	0.61
1:C:54:ALA:HB3	3:C:945:HOH:O	2.00	0.60
1:A:23:MET:HA	1:A:23:MET:CE	2.27	0.60
1:C:5:GLU:CG	1:C:6:LYS:H	2.13	0.60
1:F:60:LEU:HD11	1:F:66:ARG:CZ	2.33	0.59
1:B:92:TRP:CH2	1:B:269:PRO:HB3	2.37	0.59
1:B:107:GLU:O	1:B:108:ALA:HB2	2.03	0.59
1:F:173:ASP:O	1:F:177:ILE:HG13	2.02	0.59
1:A:23:MET:HB3	3:A:932:HOH:O	2.02	0.59
1:C:194:GLN:HA	1:C:194:GLN:NE2	2.18	0.59
1:D:78:LYS:HD3	3:D:886:HOH:O	2.02	0.59
1:A:5:GLU:HG3	1:A:6:LYS:HG3	1.85	0.58
1:C:65:HIS:HD2	1:C:247:TYR:OH	1.85	0.58
1:B:5:GLU:HG3	1:B:6:LYS:N	2.18	0.58
1:F:100:ARG:O	1:F:104:GLU:HG3	2.04	0.58
1:A:217:ILE:HG22	1:A:219:ASN:ND2	2.18	0.58
1:F:92:TRP:CZ2	1:F:269:PRO:HG3	2.39	0.57
1:F:247:TYR:CE2	1:F:253:LEU:HD11	2.39	0.57
1:E:237:ARG:NE	3:E:890:HOH:O	2.37	0.57
1:F:11:LEU:CD2	1:F:184:ILE:HD12	2.33	0.57
1:F:189:ALA:CB	1:F:192:ASP:HB2	2.34	0.57
1:A:23:MET:HE3	1:A:219:ASN:CB	2.34	0.57
1:C:189:ALA:O	1:C:192:ASP:HB3	2.05	0.57
1:B:209:ASN:ND2	3:B:936:HOH:O	2.36	0.56
1:B:78:LYS:CD	1:B:78:LYS:N	2.49	0.56
1:C:219:ASN:HA	1:D:219:ASN:OD1	2.06	0.56
1:E:6:LYS:HE2	3:E:893:HOH:O	2.06	0.56
1:F:170:LYS:O	1:F:173:ASP:HB2	2.05	0.56
1:C:192:ASP:CA	3:C:939:HOH:O	2.38	0.56
1:E:78:LYS:HD2	1:E:78:LYS:N	2.21	0.56
1:B:153:LEU:HD23	1:B:183:LEU:CD2	2.36	0.55
1:B:194:GLN:HA	1:B:194:GLN:NE2	2.21	0.55
1:B:78:LYS:HD3	3:B:894:HOH:O	2.06	0.55
1:D:26:ARG:HG2	3:D:895:HOH:O	2.06	0.55
1:D:189:ALA:HB1	1:D:192:ASP:HB2	1.88	0.55
1:E:92:TRP:CZ2	1:E:269:PRO:HG3	2.42	0.55
1:A:189:ALA:O	1:A:192:ASP:HB3	2.06	0.55
1:A:71:GLU:CG	3:A:934:HOH:O	2.55	0.55
1:A:23:MET:CE	1:A:219:ASN:CB	2.85	0.55
1:D:228:ARG:HD2	1:D:232:ARG:NH2	2.22	0.55
1:B:65:HIS:CD2	1:B:247:TYR:OH	2.58	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:218:ALA:HB3	1:F:220:ASP:OD1	2.07	0.54
1:C:221:ILE:HG22	1:C:221:ILE:O	2.06	0.54
1:C:221:ILE:HD11	1:D:219:ASN:ND2	2.23	0.54
1:E:152:LYS:HD3	1:E:181:TYR:O	2.07	0.54
1:B:107:GLU:HG2	1:B:148:VAL:HG11	1.90	0.54
1:B:5:GLU:HG3	1:B:6:LYS:H	1.71	0.53
1:B:12:LEU:HD23	1:B:99:LEU:HD23	1.90	0.53
1:D:23:MET:HE1	2:D:812:TAD:N7A	2.24	0.53
1:C:153:LEU:HD23	1:C:183:LEU:CD2	2.38	0.53
1:A:21:THR:N	1:A:24:HIS:HD2	2.02	0.53
1:A:67:VAL:O	1:A:71:GLU:HG3	2.08	0.53
1:B:204:TRP:CD1	1:F:232:ARG:HD3	2.42	0.53
1:D:157:ALA:O	1:D:161:GLU:HG3	2.08	0.53
1:C:190:GLY:O	1:C:194:GLN:HG2	2.08	0.53
1:B:194:GLN:HE21	1:B:194:GLN:CA	2.21	0.53
1:A:23:MET:CE	1:A:219:ASN:HB3	2.33	0.52
1:C:194:GLN:HE21	1:C:194:GLN:CA	2.16	0.52
1:F:61:ILE:HB	1:F:62:PRO:CD	2.39	0.52
1:C:64:TYR:CE2	1:C:68:ILE:HD11	2.44	0.52
1:F:65:HIS:CD2	1:F:247:TYR:OH	2.57	0.52
1:B:30:LEU:HD11	1:B:215:GLU:CD	2.34	0.52
1:A:158:ASP:HB2	3:A:824:HOH:O	2.10	0.51
1:D:16:SER:O	1:D:66:ARG:HD2	2.09	0.51
1:F:11:LEU:HD22	1:F:184:ILE:HD12	1.92	0.51
1:C:232:ARG:HD3	1:F:204:TRP:CD1	2.45	0.51
1:D:23:MET:CE	2:D:812:TAD:N7A	2.73	0.51
1:A:65:HIS:CD2	1:A:247:TYR:OH	2.64	0.51
1:C:188:ARG:HD3	3:C:970:HOH:O	2.11	0.51
1:D:5:GLU:N	1:D:5:GLU:CD	2.68	0.51
1:A:21:THR:H	1:A:24:HIS:CD2	2.18	0.50
1:B:11:LEU:CD2	1:B:184:ILE:HD12	2.41	0.50
1:A:5:GLU:HG3	1:A:6:LYS:N	2.26	0.50
1:E:61:ILE:HB	1:E:62:PRO:HD2	1.93	0.50
1:D:195:LYS:O	1:D:199:GLU:HG3	2.12	0.50
1:A:45:LYS:HG2	1:A:46:GLY:N	2.26	0.50
1:C:153:LEU:HD11	3:C:956:HOH:O	2.11	0.49
1:A:188:ARG:NH1	1:A:220:ASP:OD2	2.40	0.49
1:D:152:LYS:HD3	1:D:181:TYR:O	2.13	0.49
1:D:194:GLN:HE21	1:D:194:GLN:CA	2.24	0.49
2:D:812:TAD:H4B	3:D:904:HOH:O	2.11	0.49
1:A:237:ARG:NH1	1:B:78:LYS:HE3	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:LEU:HD11	1:B:215:GLU:OE2	2.13	0.49
1:B:219:ASN:HD21	1:B:221:ILE:HD11	1.77	0.49
1:C:45:LYS:NZ	3:C:900:HOH:O	2.36	0.49
1:C:220:ASP:CG	1:C:220:ASP:O	2.56	0.49
1:F:22:ASN:O	1:F:26:ARG:HB2	2.13	0.49
1:F:61:ILE:HB	1:F:62:PRO:HD2	1.95	0.49
1:C:188:ARG:HD2	1:C:218:ALA:CA	2.41	0.49
1:F:16:SER:O	1:F:17:PHE:C	2.56	0.48
1:D:16:SER:O	1:D:17:PHE:C	2.56	0.48
1:A:100:ARG:HD2	3:A:882:HOH:O	2.13	0.48
1:A:194:GLN:HE21	1:A:194:GLN:CA	2.11	0.48
1:A:71:GLU:CB	3:A:934:HOH:O	2.61	0.48
1:E:11:LEU:CD2	1:E:184:ILE:HD12	2.44	0.48
1:A:23:MET:HE2	1:A:23:MET:CA	2.37	0.47
1:C:186:VAL:O	1:C:187:THR:HB	2.14	0.47
1:D:65:HIS:HD2	1:D:247:TYR:OH	1.96	0.47
1:C:185:CYS:HB2	3:C:977:HOH:O	2.13	0.47
1:C:219:ASN:HA	1:D:219:ASN:HA	1.96	0.47
1:E:159:LEU:C	1:E:159:LEU:HD23	2.40	0.47
1:F:45:LYS:NZ	3:F:878:HOH:O	2.47	0.47
1:C:103:GLN:OE1	1:C:151:VAL:HG12	2.14	0.47
1:E:16:SER:O	1:E:17:PHE:C	2.57	0.47
1:A:16:SER:HA	1:A:51:VAL:HG12	1.97	0.47
1:A:228:ARG:NH2	3:A:960:HOH:O	2.45	0.47
1:B:219:ASN:ND2	3:B:823:HOH:O	2.47	0.47
1:D:188:ARG:HD3	1:D:220:ASP:OD2	2.15	0.47
1:E:107:GLU:HG2	1:E:148:VAL:HG11	1.96	0.47
1:F:152:LYS:NZ	1:F:209:ASN:HD21	2.13	0.47
1:C:153:LEU:CD1	3:C:956:HOH:O	2.62	0.46
1:A:189:ALA:HB1	1:A:192:ASP:HB2	1.96	0.46
1:F:107:GLU:O	1:F:108:ALA:HB2	2.15	0.46
1:A:23:MET:SD	1:A:219:ASN:HB2	2.55	0.46
1:D:83:ASP:CG	1:D:102:HIS:HE1	2.24	0.46
1:A:217:ILE:HG22	1:A:219:ASN:HD21	1.81	0.46
1:F:28:PHE:HB3	1:F:79:TRP:CH2	2.50	0.46
1:E:209:ASN:HD22	1:E:209:ASN:HA	1.56	0.46
1:C:64:TYR:O	1:C:68:ILE:HG13	2.15	0.46
1:B:107:GLU:O	1:B:108:ALA:CB	2.63	0.46
1:B:161:GLU:HG3	1:B:196:PHE:CD1	2.51	0.45
1:D:148:VAL:HA	1:D:149:PRO:HD3	1.70	0.45
1:A:190:GLY:O	1:A:194:GLN:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:THR:HA	1:C:240:VAL:HG12	1.98	0.45
1:C:61:ILE:HB	1:C:62:PRO:CD	2.45	0.45
1:E:61:ILE:HB	1:E:62:PRO:CD	2.46	0.45
1:A:11:LEU:CD2	1:A:184:ILE:HD12	2.47	0.45
1:C:166:PRO:O	1:C:167:ASN:HB2	2.17	0.45
1:B:228:ARG:HD2	1:B:232:ARG:NH2	2.32	0.44
1:D:61:ILE:HB	1:D:62:PRO:CD	2.47	0.44
1:A:5:GLU:HG3	1:A:6:LYS:H	1.82	0.44
1:C:100:ARG:HD2	3:C:929:HOH:O	2.17	0.44
1:F:267:LEU:O	1:F:268:ALA:C	2.60	0.44
1:E:103:GLN:O	1:E:107:GLU:HG3	2.18	0.44
1:E:152:LYS:NZ	1:E:209:ASN:HD21	2.16	0.44
1:D:222:SER:OG	1:D:225:LYS:HG3	2.18	0.43
1:F:152:LYS:HD3	1:F:181:TYR:O	2.18	0.43
1:C:188:ARG:CD	1:C:218:ALA:HA	2.44	0.43
1:F:247:TYR:HE2	1:F:253:LEU:HD11	1.81	0.43
1:A:78:LYS:HE3	1:B:237:ARG:HH12	1.84	0.43
1:C:189:ALA:O	1:C:192:ASP:CB	2.67	0.43
1:E:189:ALA:O	1:E:192:ASP:HB2	2.19	0.43
1:F:26:ARG:HD2	1:F:26:ARG:HA	1.74	0.43
1:E:242:ASP:HB3	3:E:890:HOH:O	2.19	0.42
1:C:191:ASN:HA	1:C:194:GLN:HB2	2.01	0.42
1:E:153:LEU:HD23	1:E:183:LEU:HD21	2.01	0.42
1:E:27:LEU:HD11	2:E:813:TAD:C2A	2.49	0.42
1:F:151:VAL:HG13	1:F:151:VAL:O	2.19	0.42
1:F:153:LEU:HB3	1:F:183:LEU:HD23	2.00	0.42
1:F:237:ARG:O	1:F:238:TYR:HB2	2.19	0.42
1:A:215:GLU:CD	3:A:951:HOH:O	2.62	0.42
1:C:172:GLU:CD	1:C:172:GLU:H	2.28	0.42
1:D:30:LEU:HD11	1:D:215:GLU:OE2	2.18	0.42
1:E:148:VAL:HA	1:E:149:PRO:HD3	1.78	0.42
1:C:215:GLU:O	1:C:215:GLU:HG3	2.20	0.42
1:F:23:MET:HG2	3:F:821:HOH:O	2.19	0.42
1:A:219:ASN:CG	1:B:219:ASN:OD1	2.63	0.41
1:D:57:LYS:HA	1:D:267:LEU:CD1	2.50	0.41
1:E:268:ALA:HB3	1:E:269:PRO:HD3	2.02	0.41
1:A:26:ARG:HD2	1:A:26:ARG:HA	1.81	0.41
1:A:188:ARG:HD2	1:A:218:ALA:HA	2.02	0.41
1:B:228:ARG:HD3	3:B:824:HOH:O	2.20	0.41
1:C:191:ASN:O	1:C:195:LYS:HB2	2.20	0.41
1:E:174:ILE:O	1:E:178:VAL:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ASN:ND2	1:B:221:ILE:HD11	2.36	0.41
1:D:188:ARG:HD2	1:D:218:ALA:HB1	2.03	0.41
1:F:67:VAL:O	1:F:71:GLU:HG3	2.20	0.41
1:B:232:ARG:HD3	1:C:204:TRP:CD1	2.56	0.41
1:C:26:ARG:HD2	1:C:26:ARG:HA	1.77	0.41
1:F:103:GLN:OE1	1:F:151:VAL:HG12	2.21	0.41
1:A:218:ALA:C	1:A:219:ASN:HD22	2.29	0.41
1:C:16:SER:O	1:C:17:PHE:C	2.63	0.41
1:C:84:THR:O	1:C:85:TRP:C	2.64	0.41
1:C:237:ARG:NH1	1:D:78:LYS:HE3	2.35	0.41
1:D:162:SER:HB2	1:D:168:LEU:HD23	2.03	0.41
1:D:266:ILE:O	1:D:267:LEU:C	2.64	0.40
1:F:32:LYS:HD2	1:F:79:TRP:CE2	2.57	0.40
1:A:189:ALA:HB1	1:A:192:ASP:CB	2.51	0.40
1:E:21:THR:HA	1:E:240:VAL:HG12	2.03	0.40
1:E:30:LEU:HD23	1:E:30:LEU:HA	1.92	0.40
1:B:188:ARG:NH1	3:B:916:HOH:O	2.53	0.40
1:C:152:LYS:NZ	1:C:181:TYR:O	2.52	0.40
1:A:159:LEU:C	1:A:159:LEU:HD23	2.47	0.40
1:B:195:LYS:O	1:B:199:GLU:HG3	2.21	0.40
1:D:194:GLN:HA	1:D:194:GLN:NE2	2.34	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	229/279 (82%)	223 (97%)	6 (3%)	0	100	100
1	B	229/279 (82%)	221 (96%)	8 (4%)	0	100	100
1	C	229/279 (82%)	219 (96%)	8 (4%)	2 (1%)	14	17
1	D	229/279 (82%)	218 (95%)	10 (4%)	1 (0%)	30	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	228/279 (82%)	221 (97%)	6 (3%)	1 (0%)	30	38
1	F	227/279 (81%)	220 (97%)	6 (3%)	1 (0%)	30	38
All	All	1371/1674 (82%)	1322 (96%)	44 (3%)	5 (0%)	30	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	220	ASP
1	D	148	VAL
1	E	220	ASP
1	F	218	ALA
1	C	218	ALA

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	204/248 (82%)	195 (96%)	9 (4%)	25	38
1	B	204/248 (82%)	194 (95%)	10 (5%)	22	34
1	C	204/248 (82%)	194 (95%)	10 (5%)	22	34
1	D	204/248 (82%)	195 (96%)	9 (4%)	25	38
1	E	203/248 (82%)	193 (95%)	10 (5%)	22	34
1	F	203/248 (82%)	193 (95%)	10 (5%)	22	34
All	All	1222/1488 (82%)	1164 (95%)	58 (5%)	23	35

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	MET
1	A	27	LEU
1	A	78	LYS
1	A	154	LEU
1	A	172	GLU

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Mol	Chain	Res	Type
1	A	194	GLN
1	A	209	ASN
1	A	217	ILE
1	A	260	ASP
1	B	19	PRO
1	B	27	LEU
1	B	78	LYS
1	B	154	LEU
1	B	161	GLU
1	B	172	GLU
1	B	209	ASN
1	B	220	ASP
1	B	260	ASP
1	B	274	THR
1	C	23	MET
1	C	27	LEU
1	C	68	ILE
1	C	71	GLU
1	C	78	LYS
1	C	154	LEU
1	C	172	GLU
1	C	194	GLN
1	C	209	ASN
1	C	215	GLU
1	D	5	GLU
1	D	78	LYS
1	D	148	VAL
1	D	154	LEU
1	D	172	GLU
1	D	192	ASP
1	D	194	GLN
1	D	209	ASN
1	D	260	ASP
1	E	19	PRO
1	E	27	LEU
1	E	78	LYS
1	E	154	LEU
1	E	172	GLU
1	E	191	ASN
1	E	207	ARG
1	E	209	ASN
1	E	252	ASN

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Mol	Chain	Res	Type
1	E	260	ASP
1	F	19	PRO
1	F	27	LEU
1	F	78	LYS
1	F	148	VAL
1	F	154	LEU
1	F	159	LEU
1	F	172	GLU
1	F	188	ARG
1	F	191	ASN
1	F	209	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	65	HIS
1	A	191	ASN
1	A	194	GLN
1	A	206	HIS
1	A	209	ASN
1	A	214	ASN
1	A	219	ASN
1	A	271	GLN
1	B	65	HIS
1	B	76	ASN
1	B	194	GLN
1	B	209	ASN
1	C	65	HIS
1	C	194	GLN
1	C	209	ASN
1	C	251	HIS
1	D	65	HIS
1	D	76	ASN
1	D	102	HIS
1	D	176	GLN
1	D	194	GLN
1	D	209	ASN
1	D	234	GLN
1	E	76	ASN
1	E	194	GLN
1	E	206	HIS

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Mol	Chain	Res	Type
1	E	209	ASN
1	E	214	ASN
1	E	245	GLN
1	F	22	ASN
1	F	65	HIS
1	F	102	HIS
1	F	176	GLN
1	F	194	GLN
1	F	209	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	TAD	C	811	-	42,47,47	1.29	6 (14%)	62,72,72	2.15	17 (27%)
2	TAD	E	813	-	42,47,47	1.31	6 (14%)	62,72,72	2.37	17 (27%)
2	TAD	F	814	-	42,47,47	1.32	7 (16%)	62,72,72	2.25	16 (25%)
2	TAD	A	809	-	42,47,47	1.34	5 (11%)	62,72,72	2.09	17 (27%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	TAD	B	810	-	42,47,47	1.35	7 (16%)	62,72,72	2.32	20 (32%)
2	TAD	D	812	-	42,47,47	1.29	5 (11%)	62,72,72	2.35	22 (35%)

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	TAD	C	811	-	-	9/30/62/62	0/5/5/5
2	TAD	E	813	-	-	9/30/62/62	0/5/5/5
2	TAD	F	814	-	-	9/30/62/62	0/5/5/5
2	TAD	A	809	-	-	10/30/62/62	0/5/5/5
2	TAD	B	810	-	-	9/30/62/62	0/5/5/5
2	TAD	D	812	-	-	7/30/62/62	0/5/5/5

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	813	TAD	PN-O2N	-3.44	1.48	1.56
2	A	809	TAD	PN-O2N	-3.42	1.48	1.56
2	C	811	TAD	PN-O2N	-3.41	1.48	1.56
2	C	811	TAD	C5A-N7A	-3.32	1.33	1.39
2	D	812	TAD	C5A-N7A	-3.11	1.33	1.39
2	F	814	TAD	PN-O2N	-2.97	1.49	1.56
2	B	810	TAD	C5A-N7A	-2.97	1.33	1.39
2	E	813	TAD	C5A-N7A	-2.96	1.33	1.39
2	F	814	TAD	C5A-N7A	-2.93	1.33	1.39
2	E	813	TAD	C1B-N9A	2.93	1.54	1.46
2	A	809	TAD	C5A-N7A	-2.77	1.34	1.39
2	F	814	TAD	O4B-C1B	2.77	1.48	1.42
2	B	810	TAD	PA-O1A	-2.76	1.45	1.51
2	B	810	TAD	PN-O2N	-2.59	1.50	1.56
2	D	812	TAD	PA-O2A	-2.55	1.50	1.56
2	D	812	TAD	C2D-C1D	-2.55	1.50	1.53
2	A	809	TAD	PN-O1N	-2.47	1.45	1.51
2	E	813	TAD	PN-O1N	-2.46	1.45	1.51
2	D	812	TAD	PN-O2N	-2.42	1.50	1.56
2	C	811	TAD	PA-O2A	-2.39	1.50	1.56
2	B	810	TAD	C2N-N3N	2.38	1.35	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	809	TAD	C4N-N3N	-2.34	1.33	1.38
2	C	811	TAD	C2N-S1N	-2.33	1.68	1.73
2	F	814	TAD	PN-O1N	-2.32	1.46	1.51
2	F	814	TAD	PA-O1A	-2.32	1.46	1.51
2	B	810	TAD	C2D-C1D	-2.28	1.50	1.53
2	B	810	TAD	PN-O1N	-2.24	1.46	1.51
2	E	813	TAD	O4B-C1B	2.23	1.47	1.42
2	D	812	TAD	PN-O1N	-2.20	1.46	1.51
2	E	813	TAD	C4A-N3A	2.19	1.38	1.34
2	B	810	TAD	O4D-C1D	2.18	1.46	1.43
2	F	814	TAD	C1B-N9A	2.16	1.52	1.46
2	F	814	TAD	C2D-C1D	-2.11	1.50	1.53
2	C	811	TAD	C4A-N3A	2.09	1.38	1.34
2	C	811	TAD	PN-O1N	-2.06	1.46	1.51
2	A	809	TAD	C1B-N9A	2.06	1.52	1.46

All (109) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	813	TAD	N3A-C4A-N9A	7.57	140.04	127.17
2	F	814	TAD	N3A-C4A-N9A	7.51	139.94	127.17
2	D	812	TAD	N3A-C4A-N9A	7.34	139.66	127.17
2	E	813	TAD	C5A-C4A-N3A	-7.29	116.67	126.72
2	C	811	TAD	N3A-C4A-N9A	7.15	139.33	127.17
2	B	810	TAD	N3A-C4A-N9A	7.10	139.24	127.17
2	F	814	TAD	C5A-C4A-N3A	-7.07	116.98	126.72
2	C	811	TAD	C5A-C4A-N3A	-6.87	117.26	126.72
2	D	812	TAD	C5A-C4A-N3A	-6.87	117.26	126.72
2	A	809	TAD	N3A-C4A-N9A	6.69	138.55	127.17
2	B	810	TAD	C5A-C4A-N3A	-6.54	117.71	126.72
2	A	809	TAD	C5A-C4A-N3A	-6.38	117.93	126.72
2	A	809	TAD	N3A-C2A-N1A	-5.61	120.09	128.58
2	F	814	TAD	N3A-C2A-N1A	-5.57	120.14	128.58
2	C	811	TAD	N3A-C2A-N1A	-5.56	120.17	128.58
2	E	813	TAD	N3A-C2A-N1A	-5.53	120.22	128.58
2	B	810	TAD	N3A-C2A-N1A	-5.37	120.45	128.58
2	D	812	TAD	N3A-C2A-N1A	-5.28	120.59	128.58
2	E	813	TAD	O6N-C6N-C4N	5.15	125.20	120.25
2	D	812	TAD	C6A-C5A-N7A	-4.63	123.17	132.09
2	B	810	TAD	O6N-C6N-C4N	4.61	124.68	120.25
2	C	811	TAD	C6A-C5A-N7A	-4.56	123.30	132.09
2	E	813	TAD	C4N-C6N-N6N	-4.47	111.73	116.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	813	TAD	C2A-N3A-C4A	4.31	122.36	111.83
2	D	812	TAD	C6A-C5A-C4A	4.30	123.05	117.18
2	B	810	TAD	C6A-C5A-N7A	-4.29	123.82	132.09
2	F	814	TAD	C2A-N3A-C4A	4.24	122.19	111.83
2	E	813	TAD	C6A-C5A-N7A	-4.21	123.98	132.09
2	C	811	TAD	C6A-C5A-C4A	4.15	122.84	117.18
2	F	814	TAD	C4N-C6N-N6N	-4.10	112.11	116.28
2	D	812	TAD	O4D-C4D-C5D	4.08	122.42	109.33
2	A	809	TAD	C6A-C5A-N7A	-4.07	124.24	132.09
2	D	812	TAD	O6N-C6N-C4N	4.02	124.11	120.25
2	F	814	TAD	O6N-C6N-C4N	4.01	124.10	120.25
2	F	814	TAD	C6A-C5A-N7A	-3.98	124.41	132.09
2	E	813	TAD	C6A-C5A-C4A	3.97	122.60	117.18
2	B	810	TAD	C6A-C5A-C4A	3.92	122.53	117.18
2	B	810	TAD	C3D-C2D-C1D	3.90	106.30	101.69
2	B	810	TAD	C2A-N3A-C4A	3.90	121.35	111.83
2	A	809	TAD	C2A-N3A-C4A	3.90	121.35	111.83
2	F	814	TAD	C6A-C5A-C4A	3.85	122.44	117.18
2	A	809	TAD	O6N-C6N-C4N	3.85	123.95	120.25
2	C	811	TAD	C2A-N3A-C4A	3.84	121.20	111.83
2	D	812	TAD	C2A-N3A-C4A	3.83	121.18	111.83
2	B	810	TAD	C4N-C6N-N6N	-3.71	112.50	116.28
2	A	809	TAD	C6A-C5A-C4A	3.62	122.12	117.18
2	B	810	TAD	O4D-C4D-C5D	3.41	120.25	109.33
2	C	811	TAD	O2A-PA-O1A	3.28	120.62	109.95
2	E	813	TAD	O4B-C4B-C5B	3.07	119.18	109.33
2	D	812	TAD	C4N-C6N-N6N	-3.05	113.17	116.28
2	E	813	TAD	O2B-C2B-C1B	2.94	120.22	110.10
2	D	812	TAD	C5N-S1N-C2N	-2.88	86.12	89.52
2	C	811	TAD	C4A-C5A-N7A	2.86	113.86	110.58
2	D	812	TAD	C4A-C5A-N7A	2.80	113.79	110.58
2	D	812	TAD	O2B-C2B-C1B	2.80	119.75	110.10
2	F	814	TAD	O4B-C4B-C5B	2.70	117.98	109.33
2	B	810	TAD	C4A-C5A-N7A	2.68	113.65	110.58
2	A	809	TAD	C4A-C5A-N7A	2.68	113.64	110.58
2	D	812	TAD	O2A-PA-O1A	2.65	118.58	109.95
2	F	814	TAD	O2B-C2B-C1B	2.64	119.21	110.10
2	B	810	TAD	O2B-C2B-C1B	2.61	119.08	110.10
2	E	813	TAD	O2A-PA-O1A	2.59	118.38	109.95
2	D	812	TAD	O4B-C4B-C5B	2.58	117.60	109.33
2	B	810	TAD	C5A-C4A-N9A	-2.55	103.04	105.81
2	D	812	TAD	S1N-C2N-N3N	2.54	117.50	114.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	812	TAD	C5A-C4A-N9A	-2.51	103.08	105.81
2	F	814	TAD	C5A-C4A-N9A	-2.51	103.08	105.81
2	F	814	TAD	O2A-PA-O1A	2.51	118.10	109.95
2	E	813	TAD	C4A-C5A-N7A	2.49	113.43	110.58
2	C	811	TAD	N6A-C6A-N1A	2.48	123.89	118.38
2	D	812	TAD	C3D-C2D-C1D	2.46	104.59	101.69
2	D	812	TAD	N6A-C6A-N1A	2.46	123.84	118.38
2	E	813	TAD	C3D-C2D-C1D	2.44	104.57	101.69
2	A	809	TAD	O2A-PA-O1A	2.44	117.88	109.95
2	A	809	TAD	C4N-C6N-N6N	-2.40	113.83	116.28
2	B	810	TAD	C4N-N3N-C2N	-2.40	108.17	110.65
2	A	809	TAD	O2B-C2B-C1B	2.36	118.22	110.10
2	C	811	TAD	S1N-C2N-N3N	2.33	117.23	114.19
2	E	813	TAD	C5A-C4A-N9A	-2.32	103.28	105.81
2	F	814	TAD	C4A-N9A-C8A	2.32	108.17	105.74
2	B	810	TAD	O2A-PA-O1A	2.32	117.49	109.95
2	B	810	TAD	N6A-C6A-N1A	2.31	123.52	118.38
2	F	814	TAD	N6A-C6A-N1A	2.28	123.46	118.38
2	B	810	TAD	C5N-S1N-C2N	-2.27	86.84	89.52
2	C	811	TAD	C4N-N3N-C2N	-2.26	108.31	110.65
2	F	814	TAD	C3D-C2D-C1D	2.25	104.34	101.69
2	A	809	TAD	C5N-S1N-C2N	-2.24	86.88	89.52
2	F	814	TAD	C4A-C5A-N7A	2.23	113.14	110.58
2	C	811	TAD	C2A-N1A-C6A	2.22	122.38	118.73
2	D	812	TAD	C4N-N3N-C2N	-2.21	108.36	110.65
2	C	811	TAD	O2B-C2B-C1B	2.21	117.72	110.10
2	A	809	TAD	C2A-N1A-C6A	2.21	122.36	118.73
2	C	811	TAD	C5A-C4A-N9A	-2.21	103.41	105.81
2	C	811	TAD	O4D-C4D-C5D	2.19	116.35	109.33
2	B	810	TAD	C2B-C1B-N9A	2.18	118.73	113.30
2	A	809	TAD	O4B-C4B-C5B	2.17	116.30	109.33
2	A	809	TAD	N6A-C6A-N1A	2.17	123.20	118.38
2	E	813	TAD	N6A-C6A-N1A	2.16	123.19	118.38
2	D	812	TAD	O2N-PN-O1N	2.16	116.98	109.95
2	B	810	TAD	O4B-C4B-C5B	2.12	116.11	109.33
2	D	812	TAD	C1B-N9A-C8A	-2.11	122.42	127.09
2	A	809	TAD	C5A-C4A-N9A	-2.11	103.52	105.81
2	B	810	TAD	C4A-N9A-C8A	2.10	107.94	105.74
2	C	811	TAD	O6N-C6N-C4N	2.09	122.26	120.25
2	A	809	TAD	O4D-C4D-C5D	2.09	116.03	109.33
2	E	813	TAD	C2B-C1B-N9A	2.07	118.46	113.30
2	C	811	TAD	O1A-PA-C3	2.05	114.45	109.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	812	TAD	O4D-C1D-C2D	2.02	107.94	105.15
2	E	813	TAD	C1B-N9A-C8A	-2.02	122.62	127.09

There are no chirality outliers.

All (53) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	809	TAD	C5D-O5D-PN-O1N
2	B	810	TAD	PN-C3-PA-O5B
2	B	810	TAD	PA-C3-PN-O1N
2	B	810	TAD	PA-C3-PN-O2N
2	B	810	TAD	C5D-O5D-PN-O1N
2	B	810	TAD	O4D-C1D-C2N-N3N
2	C	811	TAD	C5B-O5B-PA-O2A
2	C	811	TAD	PA-C3-PN-O1N
2	C	811	TAD	PA-C3-PN-O2N
2	C	811	TAD	C5D-O5D-PN-O1N
2	C	811	TAD	O4D-C1D-C2N-N3N
2	D	812	TAD	PN-C3-PA-O1A
2	D	812	TAD	PN-C3-PA-O2A
2	D	812	TAD	PN-C3-PA-O5B
2	E	813	TAD	PN-C3-PA-O5B
2	E	813	TAD	C5D-O5D-PN-O1N
2	E	813	TAD	O4D-C1D-C2N-N3N
2	F	814	TAD	C5D-O5D-PN-O1N
2	F	814	TAD	O4D-C1D-C2N-N3N
2	D	812	TAD	C3D-C4D-C5D-O5D
2	A	809	TAD	O4D-C1D-C2N-N3N
2	D	812	TAD	C5D-O5D-PN-O1N
2	A	809	TAD	PN-C3-PA-O2A
2	A	809	TAD	PA-C3-PN-O2N
2	B	810	TAD	PN-C3-PA-O2A
2	C	811	TAD	PN-C3-PA-O2A
2	E	813	TAD	PN-C3-PA-O2A
2	E	813	TAD	PA-C3-PN-O2N
2	F	814	TAD	PN-C3-PA-O2A
2	F	814	TAD	PA-C3-PN-O2N
2	D	812	TAD	O4D-C4D-C5D-O5D
2	A	809	TAD	C2D-C1D-C2N-N3N
2	B	810	TAD	C2D-C1D-C2N-N3N
2	C	811	TAD	C2D-C1D-C2N-N3N
2	D	812	TAD	C2D-C1D-C2N-N3N

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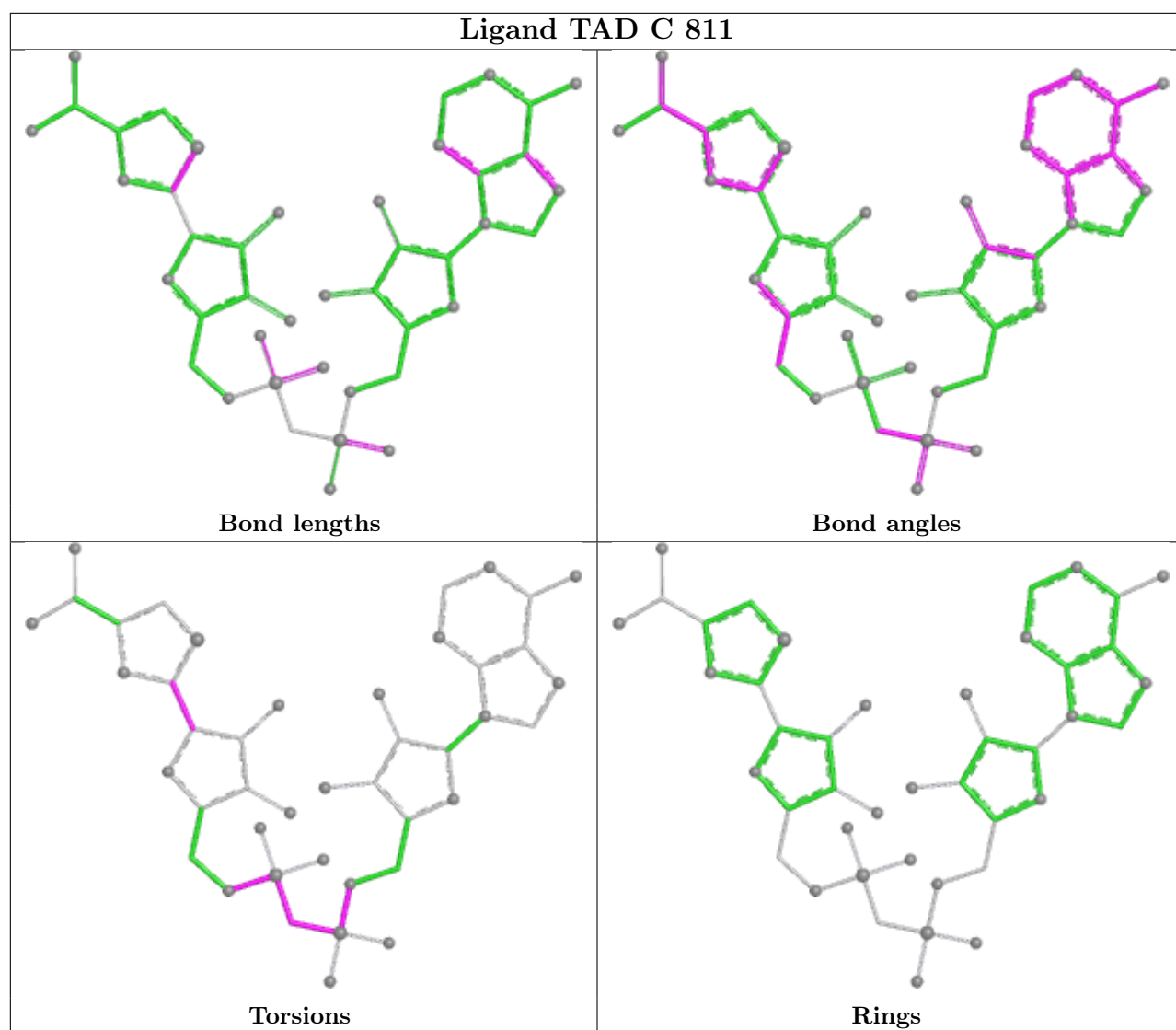
Mol	Chain	Res	Type	Atoms
2	E	813	TAD	C2D-C1D-C2N-N3N
2	F	814	TAD	C2D-C1D-C2N-N3N
2	A	809	TAD	O4D-C4D-C5D-O5D
2	A	809	TAD	C5B-O5B-PA-O2A
2	A	809	TAD	C5D-O5D-PN-O2N
2	C	811	TAD	C5D-O5D-PN-O2N
2	E	813	TAD	C5D-O5D-PN-O2N
2	F	814	TAD	C5B-O5B-PA-O2A
2	F	814	TAD	C5D-O5D-PN-O2N
2	A	809	TAD	C5B-O5B-PA-C3
2	E	813	TAD	C5B-O5B-PA-C3
2	F	814	TAD	C5B-O5B-PA-C3
2	B	810	TAD	O4D-C4D-C5D-O5D
2	A	809	TAD	C5B-O5B-PA-O1A
2	B	810	TAD	C5B-O5B-PA-O1A
2	C	811	TAD	C5B-O5B-PA-O1A
2	E	813	TAD	C5B-O5B-PA-O1A
2	F	814	TAD	C5B-O5B-PA-O1A

There are no ring outliers.

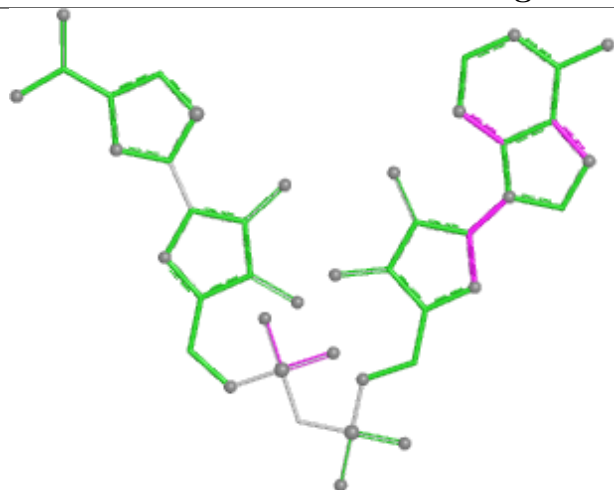
6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	811	TAD	1	0
2	E	813	TAD	2	0
2	F	814	TAD	2	0
2	A	809	TAD	1	0
2	B	810	TAD	1	0
2	D	812	TAD	5	0

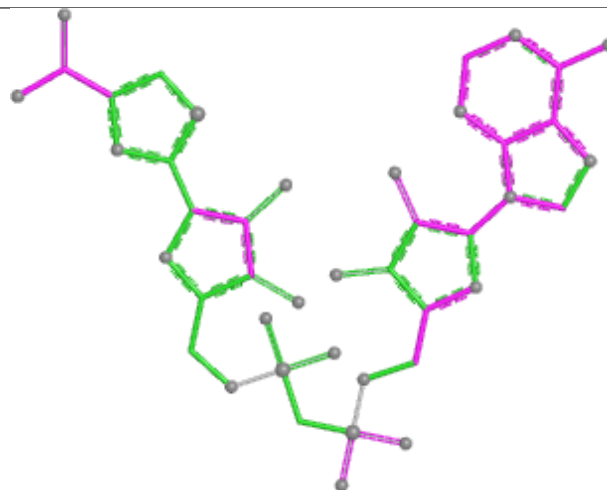
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



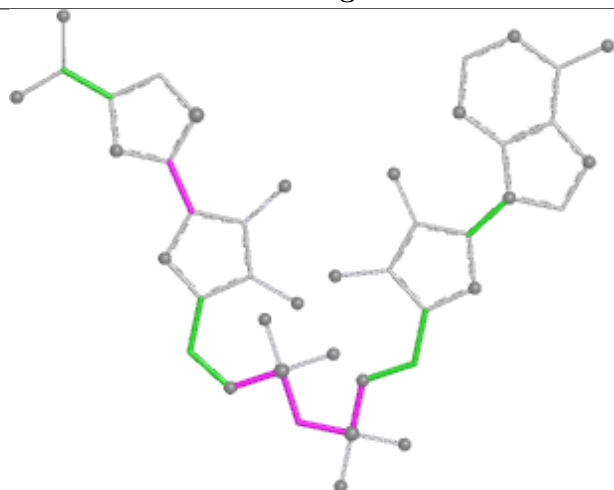
Ligand TAD E 813



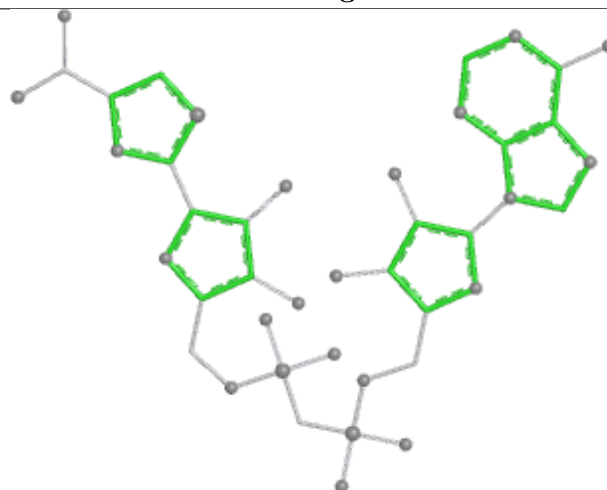
Bond lengths



Bond angles

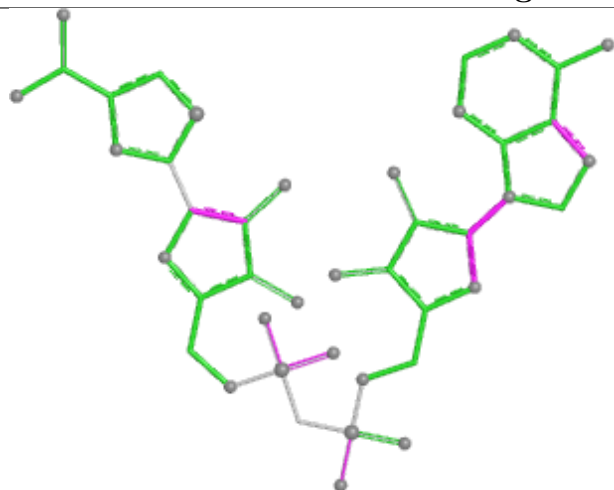


Torsions

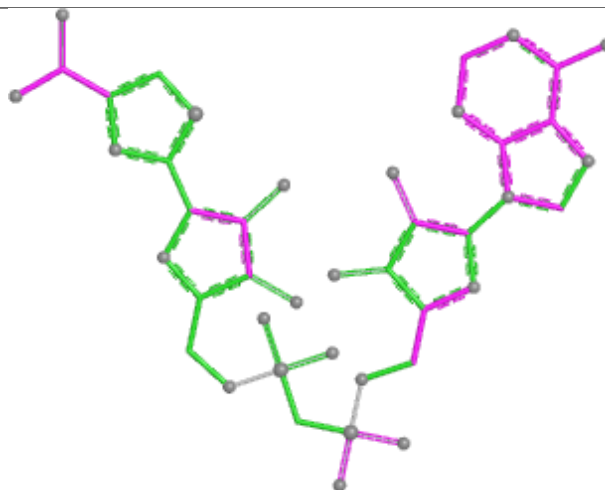


Rings

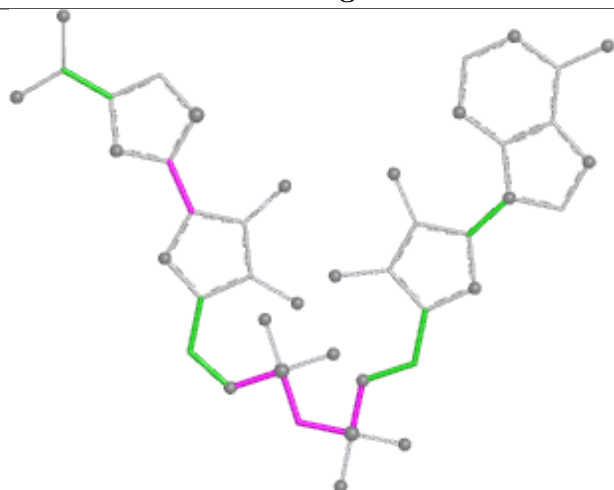
Ligand TAD F 814



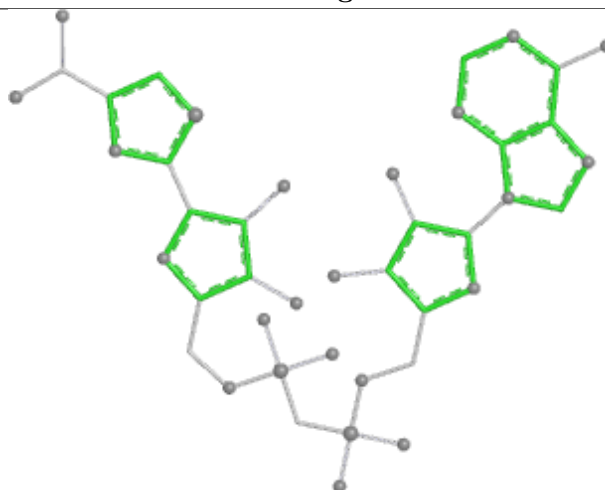
Bond lengths



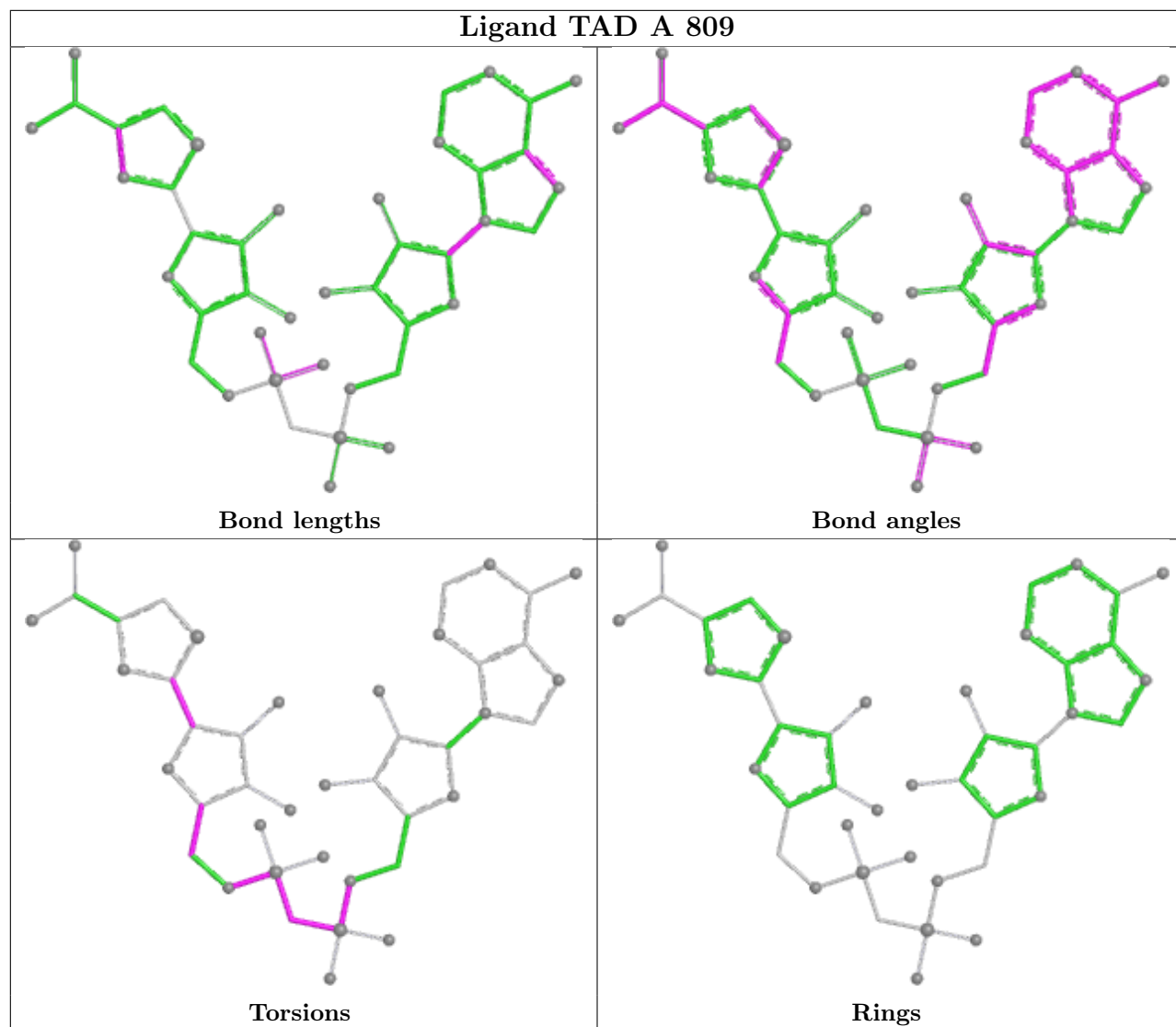
Bond angles

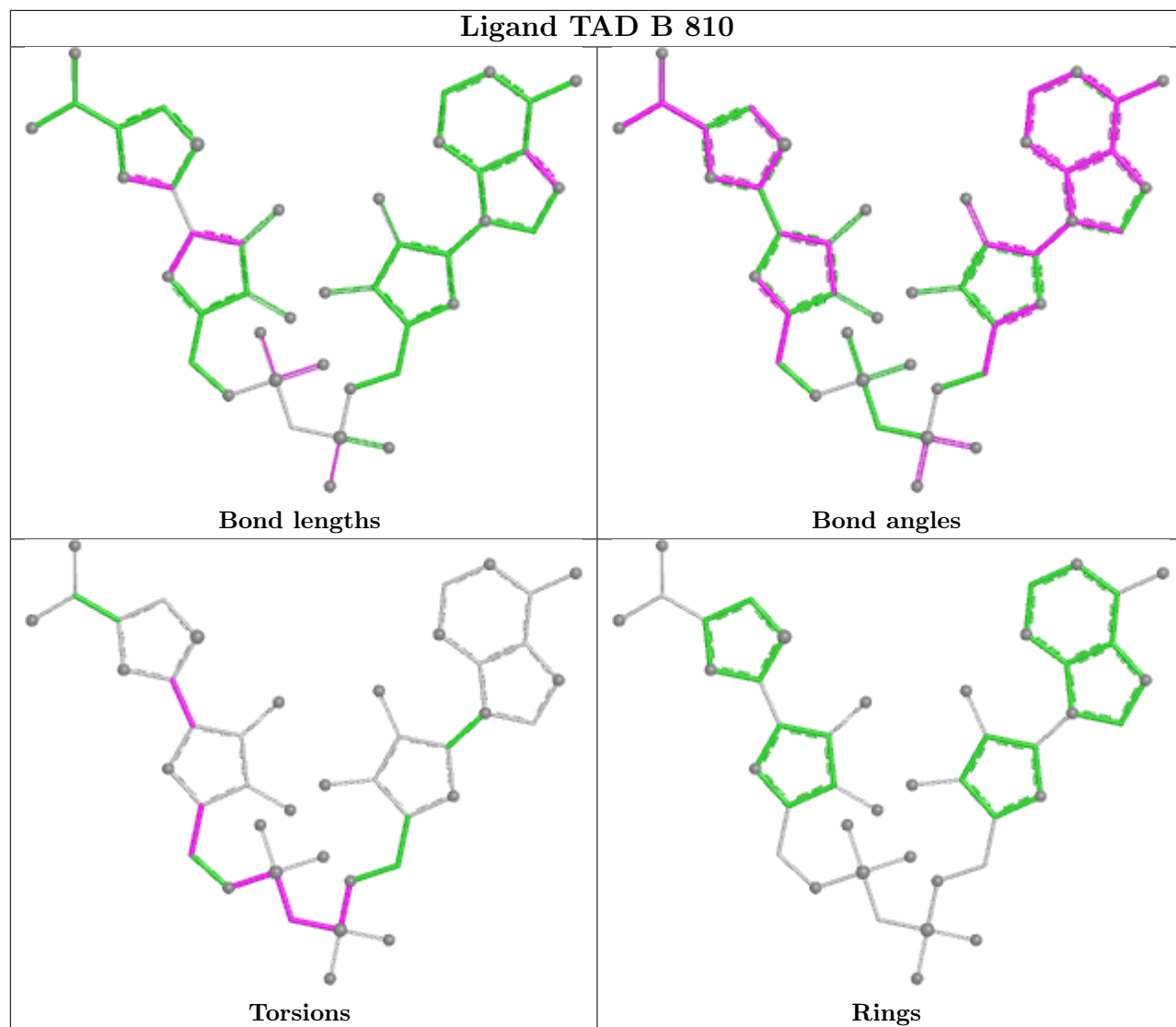


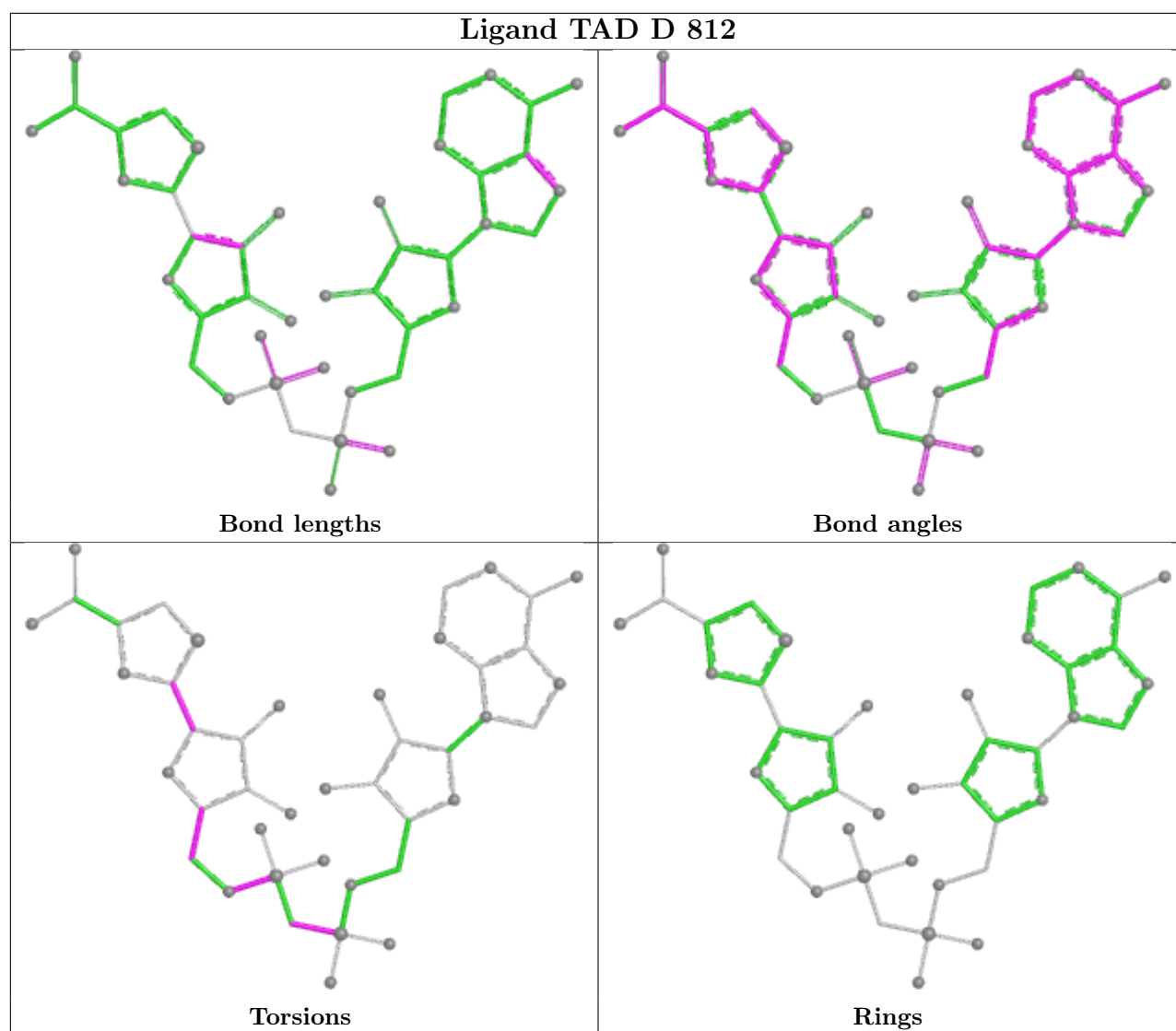
Torsions



Rings







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.