



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:00 AM UTC

PDB ID : 1ETR / pdb_00001etr
Title : REFINED 2.3 ANGSTROMS X-RAY CRYSTAL STRUCTURE OF BOVINE THROMBIN COMPLEXES FORMED WITH THE BENZAMIDINE AND ARGININE-BASED THROMBIN INHIBITORS NAPAP, 4-TAPAP AND MQPA: A STARTING POINT FOR IMPROVING ANTITHROMBOTICS
Authors : Bode, W.; Brandstetter, H.
Deposited on : 1992-07-06
Resolution : 2.20 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)
Xtrriage (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

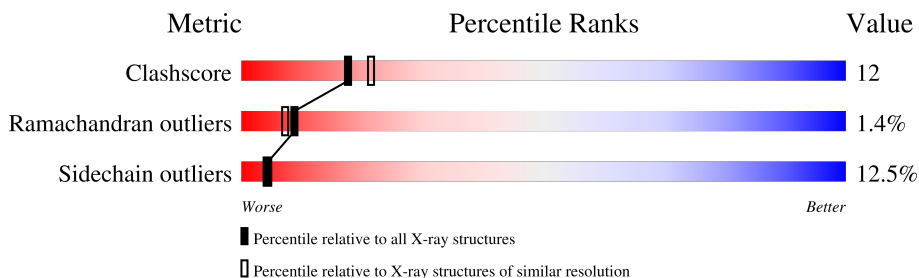
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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	L	49	
2	H	259	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3435 atoms, of which 864 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.

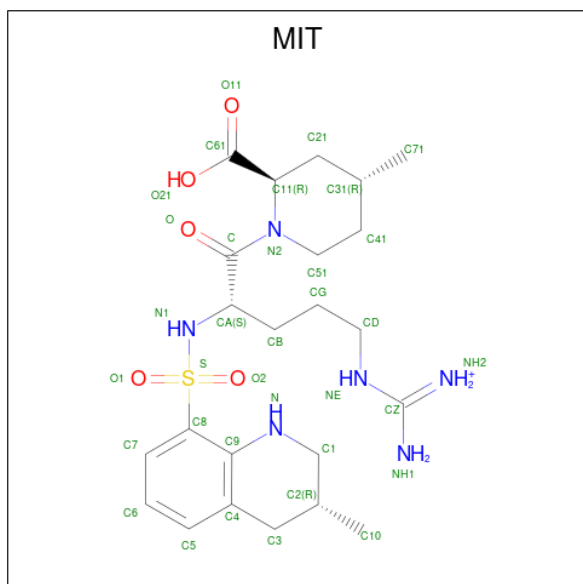
- Molecule 1 is a protein called EPSILON-THROMBIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	L	36	Total	C	H	N	O	S	141	0	0
			353	181	63	48	60	1			

- Molecule 2 is a protein called EPSILON-THROMBIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	H	259	Total	C	H	N	O	S	647	0	0
			2584	1337	490	376	369	12			

- Molecule 3 is amino{[(4S)-5-[(2R,4R)-2-carboxy-4-methylpiperidin-1-yl]-4-({[(3R)-3-methyl-1,2,3,4-tetrahydroquinolin-8-yl]sulfonyl}amino)-5-oxopentyl]amino}methaniminium (CCD ID: MIT) (formula: C₂₃H₃₇N₆O₅S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	H	1	Total	C	H	N	O	S	3	0
			42	23	7	6	5	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	17	Total 51	H 34	O 17	34	0
4	H	135	Total 405	H 270	O 135	270	0

3 Residue-property plots [i](#)

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

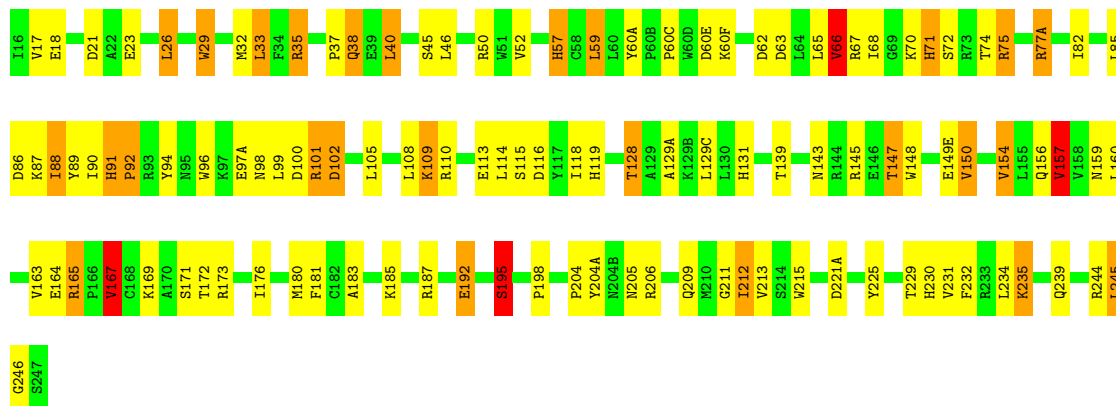
• Molecule 1: EPSILON-THROMBIN

Chain L:  51% 20% • 27%



• Molecule 2: EPSILON-THROMBIN

Chain H:  56% 33% 10% •



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	87.71Å 87.71Å 102.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.175 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3435	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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: MIT

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	L	0.76	0/294	1.73	7/390 (1.8%)
2	H	1.09	10/2148 (0.5%)	1.91	56/2905 (1.9%)
All	All	1.06	10/2442 (0.4%)	1.89	63/3295 (1.9%)

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
2	H	0	7

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	230	HIS	CD2-NE2	-6.88	1.30	1.37
2	H	131	HIS	CD2-NE2	-6.27	1.30	1.37
2	H	91	HIS	CD2-NE2	-6.02	1.31	1.37
2	H	71	HIS	CG-ND1	-5.97	1.31	1.38
2	H	57	HIS	CD2-NE2	-5.96	1.31	1.37
2	H	91	HIS	CG-ND1	-5.61	1.32	1.38
2	H	119	HIS	CG-ND1	-5.55	1.32	1.38
2	H	71	HIS	CD2-NE2	-5.36	1.31	1.37
2	H	119	HIS	CD2-NE2	-5.34	1.31	1.37
2	H	230	HIS	CG-ND1	-5.04	1.32	1.38

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	102	ASP	CA-CB-CG	9.64	122.24	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	157	VAL	CB-CA-C	-8.95	97.20	110.81
2	H	131	HIS	CB-CG-CD2	-8.82	119.74	131.20
2	H	230	HIS	ND1-CG-CD2	8.75	114.85	106.10
2	H	119	HIS	ND1-CG-CD2	8.60	114.70	106.10
2	H	165	ARG	CA-C-O	-7.90	111.36	118.63
2	H	21	ASP	CA-CB-CG	7.67	120.27	112.60
1	L	13	GLN	OE1-CD-NE2	-7.57	115.03	122.60
2	H	85	LEU	N-CA-C	-7.55	97.37	109.76
2	H	77(A)	ARG	N-CA-C	7.44	120.04	111.11
2	H	71	HIS	ND1-CG-CD2	7.33	113.43	106.10
2	H	26	LEU	N-CA-C	7.30	119.14	111.03
2	H	131	HIS	CB-CG-ND1	7.01	133.22	122.70
1	L	14	ASP	CA-CB-CG	6.96	119.56	112.60
2	H	35	ARG	CB-CG-CD	-6.80	95.66	111.30
2	H	156	GLN	OE1-CD-NE2	-6.75	115.85	122.60
1	L	11	GLN	OE1-CD-NE2	-6.74	115.86	122.60
2	H	245	LEU	N-CA-C	-6.71	99.10	109.24
2	H	159	ASN	CA-CB-CG	-6.71	105.89	112.60
2	H	129(A)	ALA	N-CA-C	6.53	119.23	111.33
2	H	119	HIS	CA-C-N	6.52	126.49	120.03
2	H	119	HIS	C-N-CA	6.52	126.49	120.03
2	H	230	HIS	CB-CG-CD2	-6.41	122.87	131.20
2	H	159	ASN	CB-CG-ND2	6.35	125.92	116.40
2	H	181	PHE	CA-CB-CG	6.34	120.14	113.80
2	H	91	HIS	CA-C-N	6.32	126.07	119.05
2	H	91	HIS	C-N-CA	6.32	126.07	119.05
2	H	98	ASN	OD1-CG-ND2	-6.24	116.36	122.60
2	H	119	HIS	CG-CD2-NE2	-6.19	101.01	107.20
2	H	159	ASN	OD1-CG-ND2	-6.14	116.46	122.60
2	H	143	ASN	CA-CB-CG	6.14	118.74	112.60
2	H	171	SER	CA-CB-OG	6.03	123.15	111.10
2	H	230	HIS	CG-CD2-NE2	-5.85	101.35	107.20
2	H	160	LEU	CA-C-N	5.77	125.53	119.76
2	H	160	LEU	C-N-CA	5.77	125.53	119.76
2	H	62	ASP	CA-CB-CG	5.66	118.26	112.60
2	H	91	HIS	ND1-CG-CD2	5.61	111.71	106.10
2	H	192	GLU	N-CA-C	-5.59	102.21	110.48
1	L	14(G)	PHE	CA-CB-CG	-5.43	108.37	113.80
2	H	59	LEU	N-CA-C	-5.36	105.80	112.93
2	H	167	VAL	CB-CA-C	5.35	118.92	111.92
2	H	235	LYS	CA-C-O	-5.35	114.88	120.55
2	H	173	ARG	N-CA-C	-5.30	106.66	113.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	239	GLN	CG-CD-NE2	-5.28	108.48	116.40
1	L	11	GLN	CG-CD-NE2	5.25	124.27	116.40
2	H	29	TRP	CG-CD2-CE3	5.25	139.15	133.90
2	H	157	VAL	CG1-CB-CG2	5.24	122.32	110.80
2	H	57	HIS	CB-CG-CD2	-5.22	124.42	131.20
2	H	221(A)	ASP	CA-CB-CG	5.22	117.82	112.60
1	L	13	GLN	CG-CD-NE2	5.21	124.21	116.40
2	H	157	VAL	CA-CB-CG1	-5.20	101.57	110.40
2	H	195	SER	CA-C-O	-5.18	115.84	121.44
2	H	92	PRO	N-CA-C	5.17	120.71	113.53
1	L	13	GLN	CA-CB-CG	5.14	124.39	114.10
2	H	231	VAL	CB-CA-C	-5.14	105.31	112.04
2	H	157	VAL	N-CA-C	5.12	115.96	108.53
2	H	185	LYS	CA-C-O	-5.10	116.01	120.60
2	H	119	HIS	CA-C-O	-5.09	114.04	119.13
2	H	17	VAL	N-CA-C	-5.05	101.18	108.71
2	H	66	VAL	CB-CA-C	5.05	118.26	110.28
2	H	148	TRP	CE2-CD2-CG	-5.02	101.17	107.20
2	H	167	VAL	CA-C-O	-5.01	114.71	120.83
2	H	88	ILE	N-CA-CB	-5.01	105.01	111.82

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	147	THR	Peptide
2	H	149(E)	GLU	Peptide
2	H	204(A)	TYR	Sidechain
2	H	225	TYR	Sidechain
2	H	245	LEU	Peptide
2	H	246	GLY	Peptide
2	H	63	ASP	Peptide

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	L	290	63	276	4	0
2	H	2094	490	2097	51	0
3	H	35	7	36	2	0
4	H	135	270	0	8	0
4	L	17	34	0	0	0
All	All	2571	864	2409	54	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 (54) 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)
2:H:212:ILE:HG13	2:H:229:THR:HB	1.59	0.84
2:H:90:ILE:HG13	4:H:654:HOH:O	1.94	0.68
2:H:209:GLN:NE2	4:H:647:HOH:O	2.27	0.66
2:H:145:ARG:HG2	2:H:150:VAL:HG11	1.76	0.66
2:H:212:ILE:HG12	4:H:647:HOH:O	1.96	0.65
2:H:68:ILE:HG22	2:H:118:ILE:HG12	1.77	0.64
2:H:211:GLY:HA2	2:H:229:THR:O	2.02	0.60
2:H:67:ARG:HG2	2:H:82:ILE:HG13	1.84	0.59
2:H:215:TRP:HB2	3:H:1:MIT:H1	1.84	0.59
2:H:164:GLU:HG2	2:H:167:VAL:HG13	1.85	0.57
2:H:70:LYS:HE3	2:H:72:SER:O	2.05	0.57
2:H:18:GLU:HG3	2:H:187:ARG:HB2	1.88	0.55
2:H:87:LYS:HD3	2:H:89:TYR:OH	2.06	0.55
2:H:165:ARG:HD3	2:H:169:LYS:NZ	2.24	0.53
2:H:128:THR:HG22	2:H:129(C):LEU:HD22	1.91	0.53
1:L:14(D):LYS:O	1:L:14(H):GLU:HB2	2.10	0.52
2:H:101:ARG:HH11	2:H:101:ARG:HB2	1.75	0.52
2:H:172:THR:HG21	2:H:176:ILE:HD11	1.92	0.52
2:H:72:SER:HB3	2:H:75:ARG:HH21	1.75	0.51
2:H:38:GLN:HB3	4:H:626:HOH:O	2.11	0.50
1:L:7:PHE:HA	1:L:12:VAL:HG23	1.94	0.50
2:H:45:SER:HB3	2:H:198:PRO:HG3	1.94	0.49
2:H:99:LEU:HD11	3:H:1:MIT:H41	1.95	0.49
2:H:29:TRP:O	2:H:45:SER:HA	2.12	0.49
2:H:101:ARG:HG2	2:H:234:LEU:HD21	1.96	0.48
2:H:172:THR:HB	4:H:659:HOH:O	2.14	0.47
2:H:139:THR:CG2	2:H:157:VAL:HG13	2.46	0.46
2:H:59:LEU:O	2:H:60(F):LYS:HG2	2.15	0.45
2:H:86:ASP:HB2	2:H:109:LYS:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:163:VAL:HG23	2:H:183:ALA:HA	1.98	0.45
1:L:7:PHE:HA	1:L:12:VAL:CG2	2.47	0.45
1:L:3:LEU:HG	2:H:206:ARG:HG2	1.99	0.45
2:H:232:PHE:O	2:H:235:LYS:HB2	2.18	0.44
2:H:32:MET:SD	2:H:70:LYS:HD3	2.57	0.44
2:H:33:LEU:O	2:H:40:LEU:HD22	2.18	0.44
2:H:71:HIS:CE1	2:H:154:VAL:HG13	2.54	0.43
2:H:23:GLU:O	2:H:26:LEU:HB2	2.19	0.43
2:H:35:ARG:HH11	2:H:35:ARG:HD2	1.68	0.43
2:H:195:SER:HA	2:H:213:VAL:HB	2.00	0.43
2:H:60(C):PRO:HD3	2:H:96:TRP:CE3	2.54	0.42
2:H:66:VAL:HG13	2:H:68:ILE:HD11	2.01	0.42
2:H:52:VAL:O	2:H:105:LEU:HD23	2.19	0.42
2:H:57:HIS:NE2	2:H:195:SER:HB2	2.33	0.42
2:H:99:LEU:O	2:H:102:ASP:HB2	2.20	0.42
2:H:139:THR:HG22	2:H:157:VAL:HG13	2.02	0.42
2:H:65:LEU:HD13	2:H:82:ILE:HG23	2.03	0.41
2:H:87:LYS:HG3	4:H:643:HOH:O	2.19	0.41
2:H:165:ARG:HD3	2:H:169:LYS:HZ3	1.83	0.41
2:H:204:PRO:C	4:H:604:HOH:O	2.63	0.41
2:H:101:ARG:HA	2:H:234:LEU:HD11	2.02	0.41
2:H:60(A):TYR:CE2	2:H:60(C):PRO:HB2	2.56	0.41
2:H:91:HIS:HA	2:H:92:PRO:HD2	1.85	0.41
2:H:77(A):ARG:HG2	4:H:594:HOH:O	2.20	0.40
2:H:94:TYR:HA	2:H:100:ASP:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	L	34/49 (69%)	26 (76%)	6 (18%)	2 (6%)	1 0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	257/259 (99%)	234 (91%)	21 (8%)	2 (1%)	16	16
All	All	291/308 (94%)	260 (89%)	27 (9%)	4 (1%)	9	7

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	1(B)	ALA
2	H	97(A)	GLU
1	L	11	GLN
2	H	60(E)	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	30/43 (70%)	28 (93%)	2 (7%)	15	17
2	H	226/226 (100%)	196 (87%)	30 (13%)	4	3
All	All	256/269 (95%)	224 (88%)	32 (12%)	4	4

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	6	LEU
1	L	11	GLN
2	H	33	LEU
2	H	37	PRO
2	H	38	GLN
2	H	40	LEU
2	H	46	LEU
2	H	50	ARG
2	H	66	VAL
2	H	74	THR
2	H	75	ARG
2	H	88	ILE

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Mol	Chain	Res	Type
2	H	101	ARG
2	H	108	LEU
2	H	109	LYS
2	H	110	ARG
2	H	113	GLU
2	H	114	LEU
2	H	115	SER
2	H	116	ASP
2	H	128	THR
2	H	147	THR
2	H	150	VAL
2	H	154	VAL
2	H	157	VAL
2	H	167	VAL
2	H	180	MET
2	H	192	GLU
2	H	195	SER
2	H	205	ASN
2	H	212	ILE
2	H	244	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	71	HIS
2	H	131	HIS
2	H	159	ASN
2	H	205	ASN
2	H	209	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	MIT	H	1	-	37,37,37	2.08	5 (13%)	48,53,53	2.57	13 (27%)

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	MIT	H	1	-	-	4/30/52/52	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	1	MIT	O1-S	8.58	1.53	1.43
3	H	1	MIT	O2-S	4.95	1.49	1.43
3	H	1	MIT	C8-S	4.42	1.83	1.77
3	H	1	MIT	S-N1	3.58	1.67	1.61
3	H	1	MIT	C21-C11	2.30	1.57	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1	MIT	O2-S-O1	-9.31	108.21	119.52
3	H	1	MIT	C8-S-N1	8.11	120.68	108.28
3	H	1	MIT	C5-C4-C9	4.66	124.28	117.53
3	H	1	MIT	C21-C31-C41	-4.33	102.80	109.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1	MIT	C9-C8-S	-4.05	114.13	121.72
3	H	1	MIT	CA-N1-S	-3.83	112.88	121.30
3	H	1	MIT	O2-S-N1	-3.44	100.64	106.88
3	H	1	MIT	C4-C3-C2	-2.62	108.10	113.63
3	H	1	MIT	CA-C-N2	2.48	123.57	118.71
3	H	1	MIT	CG-CD-NE	-2.45	105.33	112.20
3	H	1	MIT	C61-C11-N2	2.41	114.21	111.00
3	H	1	MIT	O1-S-C8	2.36	111.58	107.68
3	H	1	MIT	O21-C61-C11	2.27	120.74	113.34

There are no chirality outliers.

All (4) torsion outliers are listed below:

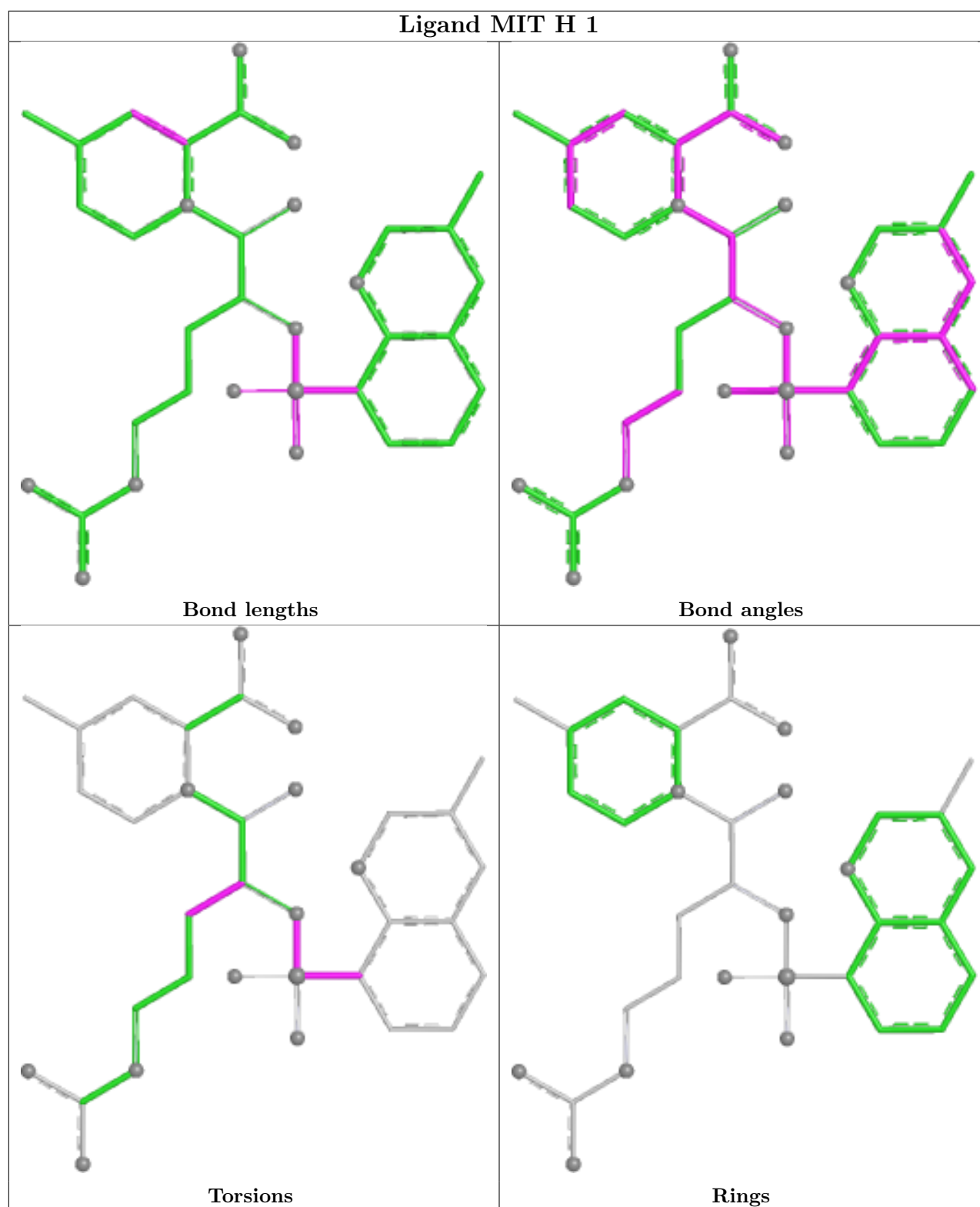
Mol	Chain	Res	Type	Atoms
3	H	1	MIT	N1-CA-CB-CG
3	H	1	MIT	C-CA-CB-CG
3	H	1	MIT	C9-C8-S-O1
3	H	1	MIT	CA-N1-S-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	MIT	2	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 ⓘ

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.