



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 12:09 am BST

PDB ID : 1ETT
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 : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

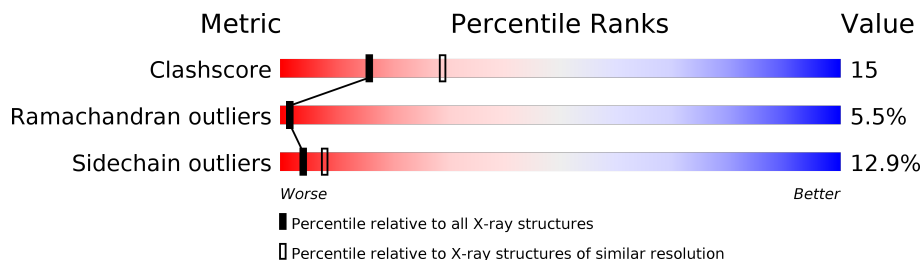
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

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 3512 atoms, of which 918 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.

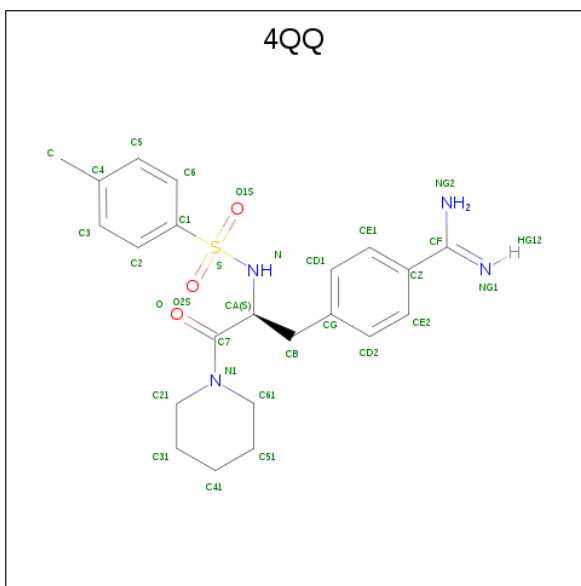
- Molecule 1 is a protein called EPSILON-THROMBIN.

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

- Molecule 2 is a protein called EPSILON-THROMBIN.

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

- Molecule 3 is 4-[(2S)-2-[[[(4-methylphenyl)sulfonyl]amino]-3-oxo-3-(piperidin-1-yl)propyl]benzene-1-carboximidamide (three-letter code: 4QQ) (formula: C₂₂H₂₈N₄O₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	S		
3	H	1	35	22	5	4	3	1	5	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	21	Total 63	H 42	O 21	42	0
4	H	159	Total 477	H 318	O 159	318	0

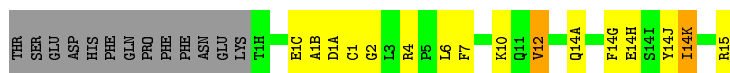
3 Residue-property plots [i](#)

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

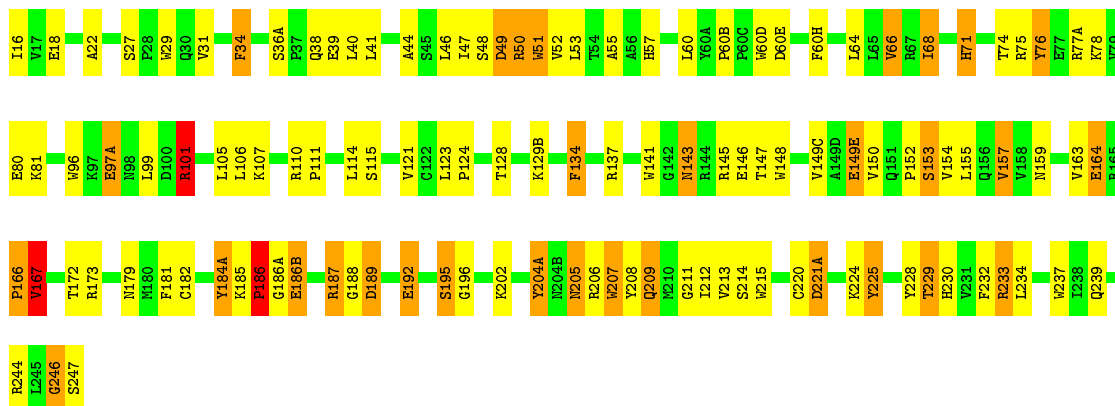
- Molecule 1: EPSILON-THROMBIN

Chain L: 



- Molecule 2: EPSILON-THROMBIN

Chain H: 



4 Data and refinement statistics

Xtrriage (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	88.55Å 88.55Å 102.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.167 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	3512	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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:
4QQ

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	0.75	0/294	1.48	2/390 (0.5%)
2	H	0.99	2/2148 (0.1%)	1.82	52/2905 (1.8%)
All	All	0.96	2/2442 (0.1%)	1.78	54/3295 (1.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	H	0	12

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	192	GLU	CD-OE2	-14.62	1.09	1.25
2	H	51	TRP	CG-CD2	-5.15	1.34	1.43

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	192	GLU	OE1-CD-OE2	-10.71	110.45	123.30
2	H	141	TRP	CD1-CG-CD2	10.51	114.71	106.30
2	H	192	GLU	CG-CD-OE1	9.55	137.39	118.30
2	H	51	TRP	CD1-CG-CD2	9.16	113.63	106.30
2	H	96	TRP	CD1-CG-CD2	8.69	113.25	106.30
2	H	164	GLU	CA-CB-CG	8.23	131.50	113.40
2	H	207	TRP	CD1-CG-CD2	8.16	112.83	106.30
2	H	148	TRP	CD1-CG-CD2	8.02	112.72	106.30
2	H	237	TRP	CD1-CG-CD2	7.98	112.69	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	145	ARG	NE-CZ-NH1	7.95	124.27	120.30
2	H	184(A)	TYR	CB-CG-CD2	-7.90	116.26	121.00
2	H	141	TRP	CE2-CD2-CG	-7.74	101.11	107.30
2	H	51	TRP	CE2-CD2-CG	-7.56	101.25	107.30
2	H	29	TRP	CE2-CD2-CG	-7.51	101.29	107.30
2	H	96	TRP	CE2-CD2-CG	-7.46	101.33	107.30
2	H	29	TRP	CD1-CG-CD2	7.39	112.22	106.30
2	H	148	TRP	CE2-CD2-CG	-7.39	101.39	107.30
2	H	60(D)	TRP	CD1-CG-CD2	7.18	112.05	106.30
2	H	237	TRP	CE2-CD2-CG	-7.12	101.60	107.30
2	H	215	TRP	CE2-CD2-CG	-6.90	101.78	107.30
2	H	141	TRP	CG-CD1-NE1	-6.82	103.28	110.10
2	H	204(A)	TYR	CB-CG-CD1	-6.82	116.91	121.00
2	H	101	ARG	NE-CZ-NH2	-6.68	116.96	120.30
2	H	207	TRP	CE2-CD2-CG	-6.55	102.06	107.30
2	H	215	TRP	CD1-CG-CD2	6.43	111.44	106.30
2	H	60(D)	TRP	CE2-CD2-CG	-6.42	102.17	107.30
2	H	207	TRP	CG-CD2-CE3	6.36	139.62	133.90
2	H	76	TYR	CB-CG-CD1	-6.10	117.34	121.00
2	H	186	PRO	N-CA-C	6.04	127.81	112.10
2	H	215	TRP	CG-CD2-CE3	5.98	139.28	133.90
2	H	51	TRP	CG-CD1-NE1	-5.90	104.20	110.10
2	H	221(A)	ASP	CB-CG-OD2	-5.89	113.00	118.30
1	L	1	CYS	CA-CB-SG	5.87	124.57	114.00
2	H	214	SER	CA-C-N	5.82	130.00	117.20
2	H	29	TRP	CB-CG-CD1	-5.75	119.53	127.00
2	H	225	TYR	CB-CG-CD1	-5.74	117.56	121.00
2	H	207	TRP	CG-CD1-NE1	-5.71	104.39	110.10
2	H	157	VAL	CB-CA-C	-5.62	100.72	111.40
2	H	215	TRP	CB-CG-CD1	-5.59	119.73	127.00
2	H	141	TRP	CB-CG-CD1	-5.49	119.86	127.00
2	H	134	PHE	CB-CG-CD2	-5.46	116.97	120.80
2	H	157	VAL	CA-CB-CG1	-5.37	102.85	110.90
2	H	237	TRP	CG-CD1-NE1	-5.35	104.75	110.10
2	H	187	ARG	NE-CZ-NH1	5.31	122.96	120.30
2	H	29	TRP	CG-CD2-CE3	5.30	138.67	133.90
2	H	46	LEU	N-CA-C	-5.29	96.71	111.00
2	H	209	GLN	CA-CB-CG	-5.18	101.99	113.40
2	H	247	SER	N-CA-C	-5.16	97.07	111.00
2	H	157	VAL	CG1-CB-CG2	5.13	119.10	110.90
2	H	39	GLU	CA-CB-CG	5.08	124.57	113.40
1	L	4	ARG	NE-CZ-NH1	-5.08	117.76	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	96	TRP	CG-CD1-NE1	-5.07	105.03	110.10
2	H	148	TRP	CG-CD1-NE1	-5.06	105.04	110.10
2	H	233	ARG	CA-CB-CG	5.03	124.45	113.40

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	101	ARG	Sidechain
2	H	110	ARG	Peptide
2	H	149(E)	GLU	Peptide
2	H	152	PRO	Peptide
2	H	166	PRO	Peptide
2	H	173	ARG	Peptide
2	H	184(A)	TYR	Sidechain
2	H	186(A)	GLY	Peptide
2	H	189	ASP	Peptide
2	H	204(A)	TYR	Sidechain
2	H	22	ALA	Peptide
2	H	246	GLY	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	8	0
2	H	2094	490	2097	58	0
3	H	30	5	0	0	0
4	H	159	318	0	14	0
4	L	21	42	0	1	0
All	All	2594	918	2373	63	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 15.

All (63) 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:31:VAL:HG13	2:H:68:ILE:HG13	1.67	0.77
1:L:7:PHE:HA	1:L:12:VAL:HG23	1.69	0.75
2:H:143:ASN:HA	2:H:150:VAL:O	1.92	0.68
2:H:57:HIS:NE2	2:H:195:SER:HB3	2.08	0.68
2:H:179:ASN:HA	2:H:233:ARG:HG2	1.75	0.67
2:H:211:GLY:HA2	2:H:229:THR:O	1.96	0.66
1:L:12:VAL:HB	4:L:109:HOH:O	1.97	0.64
2:H:172:THR:HA	2:H:224:LYS:NZ	2.14	0.63
2:H:18:GLU:HB2	2:H:188:GLY:HA2	1.81	0.62
1:L:14(J):TYR:HA	2:H:134:PHE:CE1	2.34	0.62
2:H:192:GLU:HB3	4:H:480:HOH:O	2.00	0.62
2:H:202:LYS:HE2	2:H:205:ASN:HD22	1.65	0.61
2:H:99:LEU:HG	4:H:525:HOH:O	2.01	0.60
2:H:213:VAL:HG22	2:H:228:TYR:HE1	1.66	0.59
2:H:202:LYS:HE2	2:H:205:ASN:ND2	2.19	0.58
1:L:6:LEU:HA	1:L:10:LYS:HD3	1.84	0.58
2:H:185:LYS:O	2:H:186(B):GLU:HB2	2.04	0.57
2:H:189:ASP:HB2	2:H:221(A):ASP:HB2	1.86	0.56
2:H:163:VAL:HG13	2:H:167:VAL:HG22	1.89	0.55
1:L:14(H):GLU:HA	1:L:14(K):ILE:CD1	2.37	0.55
2:H:137:ARG:HH21	2:H:159:ASN:HD21	1.53	0.54
2:H:124:PRO:HG2	2:H:232:PHE:HB2	1.90	0.53
2:H:101:ARG:HG2	2:H:234:LEU:HD21	1.91	0.52
2:H:163:VAL:CG1	2:H:167:VAL:HG22	2.40	0.52
2:H:81:LYS:HD2	4:H:419:HOH:O	2.11	0.51
2:H:51:TRP:CZ3	2:H:107:LYS:HB2	2.46	0.50
2:H:66:VAL:HG11	4:H:449:HOH:O	2.10	0.50
2:H:163:VAL:HG21	2:H:225:TYR:CD2	2.47	0.50
2:H:60:LEU:HG	2:H:60(B):PRO:HD3	1.94	0.49
2:H:124:PRO:HA	2:H:208:TYR:CD2	2.48	0.49
1:L:2:GLY:O	2:H:207:TRP:HD1	1.96	0.49
2:H:146:GLU:HG2	2:H:220:CYS:HB2	1.94	0.49
2:H:55:ALA:H	2:H:196:GLY:HA2	1.79	0.48
1:L:14(A):GLN:HG3	4:H:421:HOH:O	2.15	0.47
2:H:212:ILE:HG12	4:H:472:HOH:O	2.15	0.47
2:H:47:ILE:HG22	2:H:121:VAL:HG12	1.96	0.47
2:H:53:LEU:HA	2:H:105:LEU:HD23	1.97	0.47
2:H:57:HIS:CE1	2:H:195:SER:HB3	2.49	0.46
2:H:189:ASP:CB	2:H:221(A):ASP:HB2	2.46	0.45
2:H:18:GLU:HG3	2:H:187:ARG:HB2	1.99	0.45
1:L:14(G):PHE:CZ	2:H:202:LYS:HD3	2.51	0.45
2:H:16:ILE:N	4:H:458:HOH:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:124:PRO:HD3	2:H:209:GLN:O	2.17	0.45
2:H:181:PHE:HE1	2:H:230:HIS:HA	1.81	0.44
2:H:68:ILE:N	2:H:68:ILE:HD12	2.33	0.44
2:H:212:ILE:HD12	4:H:485:HOH:O	2.17	0.44
2:H:233:ARG:HG3	4:H:518:HOH:O	2.17	0.43
2:H:234:LEU:HG	4:H:518:HOH:O	2.18	0.43
2:H:48:SER:HG	2:H:51:TRP:HD1	1.66	0.43
2:H:77(A):ARG:HG3	2:H:78:LYS:N	2.32	0.43
2:H:206:ARG:HB2	4:H:451:HOH:O	2.18	0.43
2:H:60(H):PHE:CD2	2:H:64:LEU:HD21	2.53	0.43
2:H:50:ARG:HB3	2:H:51:TRP:CD1	2.54	0.43
2:H:213:VAL:HG22	2:H:228:TYR:CE1	2.51	0.42
2:H:75:ARG:HB3	4:H:420:HOH:O	2.18	0.42
2:H:41:LEU:HD13	2:H:60(H):PHE:CE2	2.55	0.42
2:H:44:ALA:HB1	2:H:52:VAL:CG1	2.50	0.41
2:H:71:HIS:HB2	4:H:425:HOH:O	2.20	0.41
2:H:163:VAL:HG21	2:H:225:TYR:CE2	2.55	0.41
2:H:49:ASP:O	2:H:111:PRO:HA	2.21	0.41
2:H:154:VAL:HG12	2:H:155:LEU:O	2.20	0.40
2:H:34:PHE:HZ	4:H:443:HOH:O	2.03	0.40
2:H:57:HIS:NE2	2:H:195:SER:CB	2.83	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%)	28 (82%)	4 (12%)	2 (6%)	1	1
2	H	257/259 (99%)	214 (83%)	29 (11%)	14 (5%)	2	2
All	All	291/308 (94%)	242 (83%)	33 (11%)	16 (6%)	2	2

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	1(B)	ALA
2	H	97(A)	GLU
2	H	76	TYR
2	H	149(E)	GLU
2	H	167	VAL
2	H	143	ASN
2	H	153	SER
2	H	186	PRO
2	H	49	ASP
2	H	71	HIS
2	H	149(C)	VAL
2	H	186(B)	GLU
1	L	1(C)	GLU
2	H	60(E)	ASP
2	H	166	PRO
2	H	246	GLY

5.3.2 Protein sidechains [i](#)

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%)	26 (87%)	4 (13%)	4 7
2	H	226/226 (100%)	197 (87%)	29 (13%)	4 8
All	All	256/269 (95%)	223 (87%)	33 (13%)	4 8

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1(A)	ASP
1	L	12	VAL
1	L	14(K)	ILE
1	L	15	ARG
2	H	27	SER
2	H	34	PHE
2	H	36(A)	SER

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Mol	Chain	Res	Type
2	H	38	GLN
2	H	40	LEU
2	H	50	ARG
2	H	66	VAL
2	H	68	ILE
2	H	74	THR
2	H	80	GLU
2	H	97(A)	GLU
2	H	106	LEU
2	H	114	LEU
2	H	115	SER
2	H	123	LEU
2	H	128	THR
2	H	129(B)	LYS
2	H	147	THR
2	H	153	SER
2	H	157	VAL
2	H	164	GLU
2	H	167	VAL
2	H	182	CYS
2	H	186	PRO
2	H	195	SER
2	H	205	ASN
2	H	229	THR
2	H	239	GLN
2	H	244	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	60(G)	ASN
2	H	159	ASN
2	H	179	ASN
2	H	205	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [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	4QQ	H	301	-	32,32,32	2.39	6 (18%)	41,45,45	2.44	11 (26%)

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	4QQ	H	301	-	-	2/27/35/35	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	301	4QQ	O1S-S	8.76	1.53	1.43
3	H	301	4QQ	C1-S	5.93	1.85	1.76
3	H	301	4QQ	O2S-S	4.96	1.49	1.43
3	H	301	4QQ	CZ-CF	-4.06	1.40	1.47
3	H	301	4QQ	S-N	3.00	1.66	1.61
3	H	301	4QQ	O-C7	2.12	1.26	1.22

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	H	301	4QQ	O2S-S-O1S	-9.48	107.89	119.55
3	H	301	4QQ	O2S-S-N	6.75	119.09	106.88
3	H	301	4QQ	C31-C21-N1	-4.39	101.99	110.66
3	H	301	4QQ	O1S-S-N	-4.12	99.43	106.88
3	H	301	4QQ	C1-S-N	3.44	112.54	107.78
3	H	301	4QQ	CA-C7-N1	-3.34	113.83	118.87
3	H	301	4QQ	C51-C61-N1	-2.75	105.23	110.66
3	H	301	4QQ	C51-C41-C31	-2.46	103.60	111.18
3	H	301	4QQ	C2-C1-S	2.26	122.23	119.77
3	H	301	4QQ	O-C7-CA	2.22	123.67	119.66
3	H	301	4QQ	C41-C31-C21	-2.06	107.20	111.19

There are no chirality outliers.

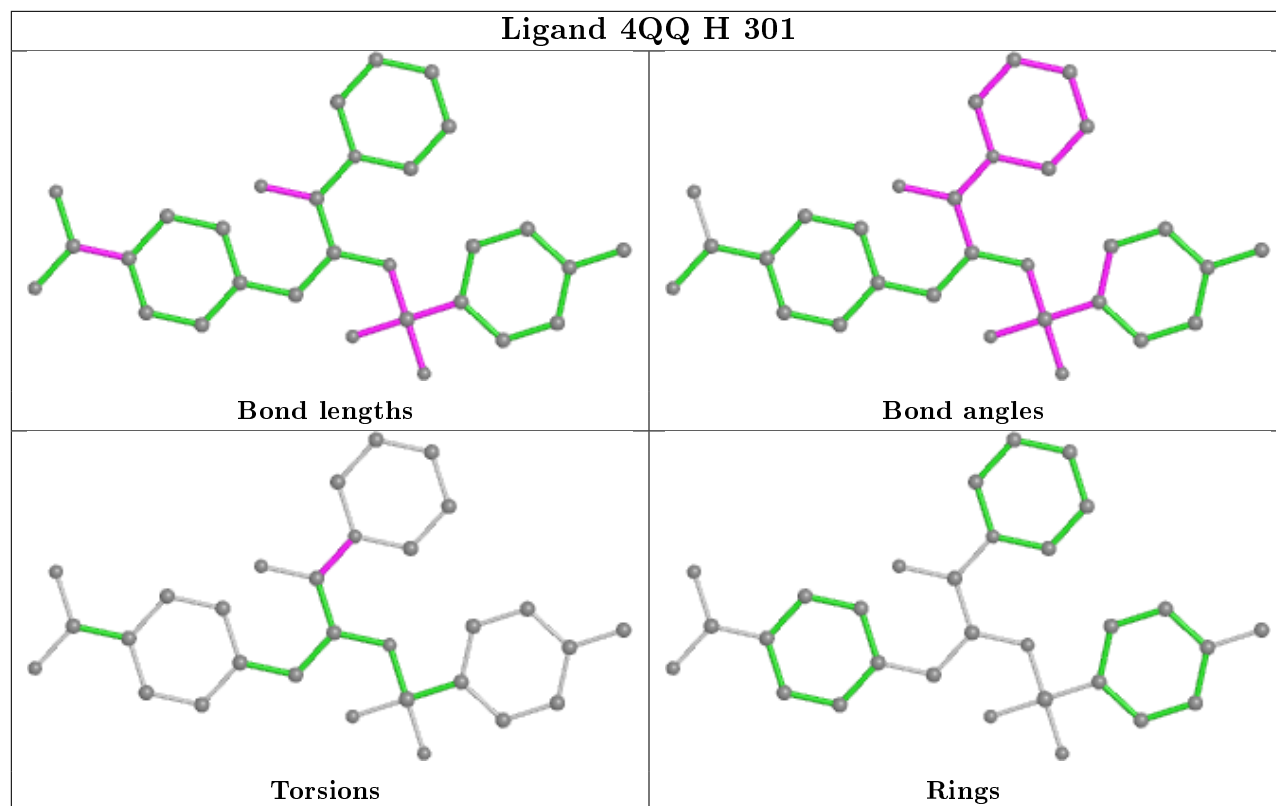
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	H	301	4QQ	O-C7-N1-C61
3	H	301	4QQ	CA-C7-N1-C61

There are no ring outliers.

No monomer is involved in short contacts.

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

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.