



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:30 AM UTC

PDB ID : 1D5J / pdb_00001d5j
Title : CRYSTAL STRUCTURE OF MMP3 COMPLEXED WITH A THIAZEPINE
BASED INHIBITOR.
Authors : Almstead, N.G.; Bradley, R.S.; Pikul, S.; De, B.; Natchus, M.G.
Deposited on : 1999-10-07
Resolution : 2.60 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

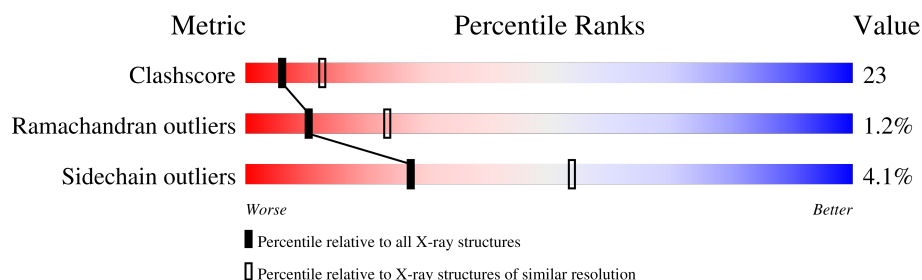
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)

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	173	 62% 30% 5% ••
1	B	173	 50% 38% 10% •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2869 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 STROMELYSIN-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	0	0
			1346	865	224	255	2			
1	B	173	Total	C	N	O	S	1	0	0
			1376	882	228	264	2			

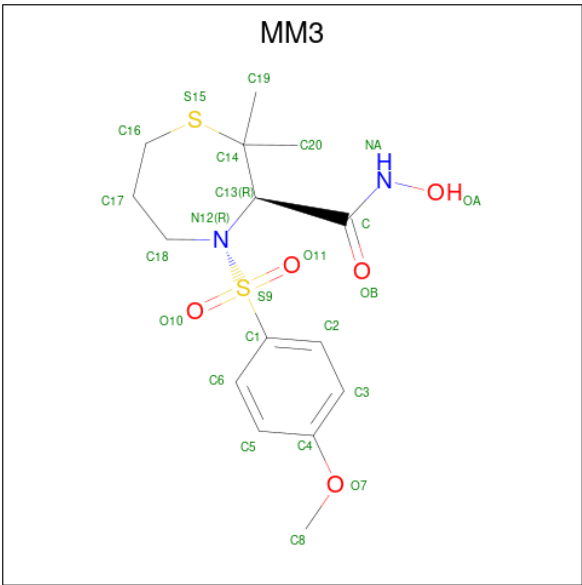
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Ca	0	0
			3	3		
3	B	3	Total	Ca	0	0
			3	3		

- Molecule 4 is N-HYDROXY-4-[(4-METHOXYLPHENYL)SULFONYL]-2,2-DIMETHYL-HEXAHYDRO-1,4-THIAZEPINE-3(S)-CARBOXAMIDE (CCD ID: MM3) (formula: C₁₅H₂₂N₂O₅S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			24	15	2	5	2		
4	B	1	Total	C	N	O	S	0	0
			24	15	2	5	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	49	Total	O	0	0
			49	49		
5	B	40	Total	O	0	0
			40	40		

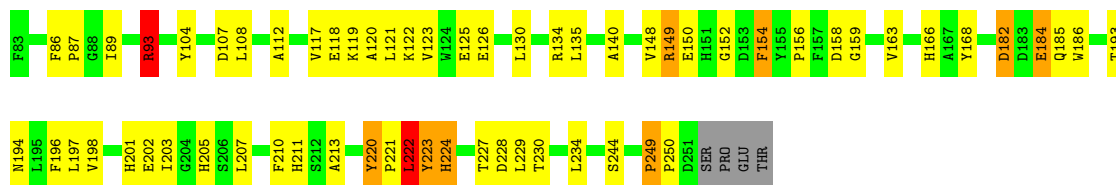
3 Residue-property plots [i](#)

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.

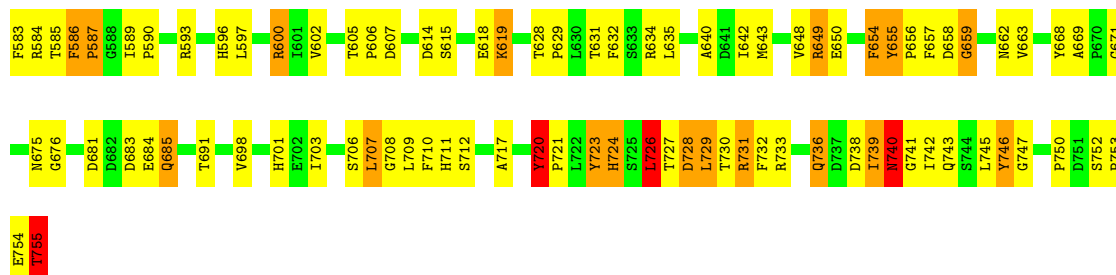
• Molecule 1: STROMELYSIN-1

Chain A: 



• Molecule 1: STROMELYSIN-1

Chain B: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	37.78Å 78.54Å 105.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.60	Depositor
% Data completeness (in resolution range)	92.2 (8.00-2.60)	Depositor
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 98.1	Depositor
R, R_{free}	0.237 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2869	wwPDB-VP
Average B, all atoms (Å ²)	18.0	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: ZN, MM3, CA

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	0.61	0/1390	1.58	34/1899 (1.8%)
1	B	0.62	0/1421	1.91	59/1941 (3.0%)
All	All	0.62	0/2811	1.75	93/3840 (2.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	2
All	All	0	5

There are no bond length outliers.

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	HIS	CA-CB-CG	-19.54	94.26	113.80
1	B	726	LEU	N-CA-C	-18.86	82.54	110.46
1	B	733	ARG	NE-CZ-NH2	-17.73	103.25	119.20
1	A	149	ARG	NE-CZ-NH1	-17.44	104.06	121.50
1	B	600	ARG	NE-CZ-NH2	-16.86	104.02	119.20
1	A	149	ARG	NE-CZ-NH2	16.28	133.85	119.20
1	B	733	ARG	NE-CZ-NH1	14.44	135.94	121.50
1	B	600	ARG	NE-CZ-NH1	14.02	135.52	121.50
1	A	93	ARG	NE-CZ-NH1	-13.78	107.72	121.50
1	B	731	ARG	NE-CZ-NH2	-13.51	107.04	119.20
1	B	736	GLN	CG-CD-NE2	-13.41	96.29	116.40
1	A	154	PHE	CA-CB-CG	-12.83	100.97	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	685	GLN	CG-CD-NE2	-12.78	97.23	116.40
1	B	740	ASN	CA-CB-CG	12.00	124.60	112.60
1	A	93	ARG	NE-CZ-NH2	11.82	129.84	119.20
1	B	710	PHE	CA-CB-CG	10.87	124.67	113.80
1	A	223	TYR	CA-CB-CG	-10.77	94.52	113.90
1	B	586	PHE	CB-CG-CD2	-10.34	103.13	120.70
1	B	746	TYR	CB-CG-CD1	-9.89	105.97	120.80
1	B	736	GLN	CG-CD-OE1	9.68	140.16	120.80
1	B	736	GLN	CB-CG-CD	-9.58	96.31	112.60
1	B	586	PHE	CB-CG-CD1	9.27	136.46	120.70
1	B	740	ASN	CB-CG-ND2	-9.27	102.50	116.40
1	B	723	TYR	CB-CG-CD2	-9.07	107.20	120.80
1	B	685	GLN	CG-CD-OE1	9.04	138.88	120.80
1	B	731	ARG	NE-CZ-NH1	8.96	130.47	121.50
1	B	701	HIS	CA-CB-CG	8.90	122.70	113.80
1	B	732	PHE	CB-CG-CD2	-8.83	105.69	120.70
1	B	710	PHE	CB-CG-CD1	-8.79	105.75	120.70
1	A	223	TYR	CB-CG-CD2	-8.62	107.87	120.80
1	B	685	GLN	CA-CB-CG	-8.51	97.08	114.10
1	B	701	HIS	CB-CG-CD2	-8.48	120.17	131.20
1	A	201	HIS	CA-CB-CG	8.20	122.00	113.80
1	B	723	TYR	CB-CG-CD1	8.17	133.05	120.80
1	B	602	VAL	N-CA-C	8.05	118.83	110.62
1	A	222	LEU	CA-C-N	-7.88	111.93	123.11
1	A	222	LEU	C-N-CA	-7.88	111.93	123.11
1	B	655	TYR	CB-CG-CD1	-7.84	109.04	120.80
1	B	746	TYR	CB-CG-CD2	7.84	132.56	120.80
1	B	728	ASP	N-CA-C	7.78	121.29	109.62
1	A	222	LEU	N-CA-C	7.77	127.35	110.80
1	B	727	THR	N-CA-CB	-7.71	99.95	110.67
1	B	732	PHE	CB-CG-CD1	7.52	133.49	120.70
1	B	733	ARG	CG-CD-NE	-7.47	95.56	112.00
1	A	159	GLY	CA-C-N	7.32	127.30	119.76
1	A	159	GLY	C-N-CA	7.32	127.30	119.76
1	B	710	PHE	CB-CG-CD2	7.25	133.03	120.70
1	B	607	ASP	N-CA-C	7.21	119.14	111.28
1	B	655	TYR	CB-CG-CD2	6.96	131.23	120.80
1	A	249	PRO	CA-C-N	6.93	128.50	119.84
1	A	249	PRO	C-N-CA	6.93	128.50	119.84
1	B	733	ARG	CA-CB-CG	6.92	127.94	114.10
1	B	723	TYR	CA-CB-CG	-6.88	101.52	113.90
1	B	733	ARG	CB-CG-CD	-6.72	95.84	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	720	TYR	CB-CG-CD2	-6.72	110.72	120.80
1	A	149	ARG	CG-CD-NE	6.69	126.71	112.00
1	B	600	ARG	CB-CG-CD	6.59	126.45	111.30
1	A	93	ARG	CD-NE-CZ	6.43	133.40	124.40
1	B	585	THR	N-CA-CB	6.38	122.35	110.87
1	A	223	TYR	CB-CG-CD1	6.34	130.31	120.80
1	B	669	ALA	CA-C-N	6.32	126.34	119.89
1	B	669	ALA	C-N-CA	6.32	126.34	119.89
1	A	149	ARG	CB-CG-CD	6.26	125.70	111.30
1	A	166	HIS	CB-CG-ND1	-6.20	113.41	122.70
1	B	754	GLU	CA-CB-CG	6.08	126.25	114.10
1	A	229	LEU	CA-C-N	6.07	128.42	120.28
1	A	229	LEU	C-N-CA	6.07	128.42	120.28
1	A	154	PHE	CB-CG-CD2	-6.05	110.41	120.70
1	A	120	ALA	N-CA-C	-5.97	104.85	111.36
1	B	634	ARG	N-CA-C	5.93	119.92	109.96
1	A	244	SER	N-CA-C	-5.88	104.87	111.28
1	A	93	ARG	CB-CG-CD	5.77	124.57	111.30
1	B	655	TYR	CA-C-N	5.76	126.64	120.13
1	B	655	TYR	C-N-CA	5.76	126.64	120.13
1	B	701	HIS	CB-CG-ND1	5.71	131.26	122.70
1	B	600	ARG	CG-CD-NE	5.64	124.41	112.00
1	B	739	ILE	N-CA-C	5.56	115.70	110.30
1	B	740	ASN	CB-CG-OD1	5.56	131.92	120.80
1	B	654	PHE	N-CA-C	5.50	118.04	111.71
1	B	600	ARG	CA-CB-CG	5.39	124.88	114.10
1	A	168	TYR	N-CA-C	5.34	118.44	109.95
1	A	220	TYR	N-CA-C	-5.29	102.78	110.08
1	A	213	ALA	N-CA-C	-5.29	106.33	112.89
1	A	134	ARG	N-CA-C	5.28	118.42	109.76
1	A	184	GLU	N-CA-C	-5.28	101.96	110.14
1	A	158	ASP	N-CA-C	5.28	119.94	111.81
1	A	166	HIS	CB-CG-CD2	5.26	138.04	131.20
1	B	600	ARG	CD-NE-CZ	5.19	131.67	124.40
1	B	755	THR	N-CA-C	5.15	125.42	111.00
1	B	659	GLY	CA-C-N	5.09	125.38	119.93
1	B	659	GLY	C-N-CA	5.09	125.38	119.93
1	B	707	LEU	N-CA-C	5.02	119.46	113.23
1	B	720	TYR	CB-CG-CD1	5.02	128.33	120.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	154	PHE	Sidechain
1	A	223	TYR	Sidechain
1	A	93	ARG	Sidechain
1	B	720	TYR	Sidechain
1	B	746	TYR	Sidechain

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	1346	0	1270	58	0
1	B	1376	0	1295	63	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	24	0	21	2	0
4	B	24	0	21	5	0
5	A	49	0	0	1	0
5	B	40	0	0	1	0
All	All	2869	0	2607	122	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 23.

All (122) 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:A:224:HIS:O	1:A:224:HIS:CD2	1.97	1.17
1:A:224:HIS:NE2	1:A:227:THR:HG21	1.67	1.09
1:A:228:ASP:HB2	1:A:230:THR:HG22	1.29	1.09
1:A:224:HIS:CD2	1:A:227:THR:HG23	1.86	1.08
1:A:224:HIS:O	1:A:224:HIS:HD2	1.32	1.02
1:A:224:HIS:NE2	1:A:227:THR:CG2	2.24	1.00
1:A:224:HIS:CD2	1:A:227:THR:CG2	2.56	0.89
1:B:736:GLN:HG3	1:B:740:ASN:OD1	1.77	0.85
1:A:220:TYR:C	1:A:222:LEU:H	1.84	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:VAL:HG13	1:A:234:LEU:HG	1.63	0.81
1:A:224:HIS:CE1	1:A:227:THR:HG21	2.18	0.78
1:B:586:PHE:HE1	1:B:709:LEU:O	1.69	0.74
1:A:224:HIS:CD2	1:A:224:HIS:C	2.57	0.73
1:B:586:PHE:HD1	1:B:708:GLY:O	1.71	0.72
1:B:589:ILE:HG22	1:B:589:ILE:O	1.87	0.72
1:B:663:VAL:HG13	4:B:901:MM3:H203	1.73	0.71
1:A:118:GLU:O	1:A:122:LYS:HG2	1.92	0.70
1:A:228:ASP:HB2	1:A:230:THR:CG2	2.17	0.70
1:A:108:LEU:HG	1:A:112:ALA:HB3	1.74	0.69
1:A:194:ASN:HD22	1:A:197:LEU:H	1.41	0.69
1:B:586:PHE:O	1:B:590:PRO:HD2	1.93	0.67
1:B:584:ARG:HD2	1:B:711:HIS:O	1.95	0.67
1:B:615:SER:O	1:B:619:LYS:HD2	1.96	0.66
1:B:663:VAL:HG11	5:B:45:HOH:O	1.95	0.65
1:B:703:ILE:O	1:B:707:LEU:HG	1.96	0.65
1:A:186:TRP:CZ3	1:A:193:THR:CG2	2.80	0.65
1:A:186:TRP:CZ3	1:A:193:THR:HG22	2.31	0.65
1:B:597:LEU:O	1:B:632:PHE:HA	1.96	0.64
1:B:731:ARG:HG3	1:B:731:ARG:O	1.96	0.64
1:B:659:GLY:HA2	1:B:683:ASP:OD2	1.98	0.64
1:B:628:THR:HG21	1:B:742:ILE:HG13	1.80	0.64
1:B:723:TYR:CD2	1:B:723:TYR:C	2.75	0.64
1:B:655:TYR:CD1	1:B:655:TYR:N	2.65	0.63
1:A:152:GLY:HA2	1:B:650:GLU:HG3	1.83	0.61
1:A:123:VAL:CG1	1:A:234:LEU:HG	2.28	0.61
1:B:586:PHE:CE1	1:B:709:LEU:O	2.53	0.61
1:A:220:TYR:HD2	1:A:222:LEU:HB2	1.65	0.61
1:B:736:GLN:HA	1:B:739:ILE:HD12	1.83	0.59
1:B:600:ARG:HD3	1:B:643:MET:HE1	1.85	0.59
1:B:649:ARG:HD3	1:B:657:PHE:O	2.03	0.58
1:A:224:HIS:HE2	1:A:227:THR:HG21	1.63	0.58
1:B:728:ASP:C	1:B:730:THR:H	2.11	0.57
1:A:203:ILE:O	1:A:207:LEU:HG	2.04	0.57
1:A:135:LEU:HD12	1:A:140:ALA:HB2	1.87	0.56
1:A:182:ASP:C	1:A:184:GLU:H	2.14	0.54
1:B:698:VAL:HG13	4:B:901:MM3:H51	1.89	0.54
1:A:220:TYR:O	1:A:222:LEU:N	2.40	0.54
1:A:122:LYS:O	1:A:126:GLU:HG3	2.08	0.53
1:B:586:PHE:HD1	1:B:708:GLY:C	2.16	0.53
1:A:202:GLU:O	1:A:205:HIS:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:ASN:HD21	1:A:196:PHE:HB3	1.72	0.53
1:B:583:PHE:HA	1:B:712:SER:OG	2.09	0.53
1:B:589:ILE:O	1:B:589:ILE:CG2	2.56	0.52
1:B:654:PHE:HB2	1:B:655:TYR:CE1	2.43	0.52
1:A:186:TRP:CZ3	1:A:193:THR:HG21	2.43	0.52
1:B:596:HIS:HD2	1:B:631:THR:OG1	1.93	0.51
1:B:635:LEU:HD12	1:B:640:ALA:HB2	1.94	0.50
1:B:614:ASP:O	1:B:618:GLU:HG3	2.11	0.50
1:B:724:HIS:ND1	1:B:724:HIS:C	2.70	0.50
1:B:755:THR:O	1:B:755:THR:HG23	2.12	0.50
1:B:663:VAL:HG22	4:B:901:MM3:S15	2.52	0.49
1:B:726:LEU:HD12	1:B:728:ASP:N	2.27	0.48
1:A:220:TYR:CD2	1:A:222:LEU:HB2	2.48	0.48
1:A:210:PHE:CG	1:A:211:HIS:N	2.83	0.47
1:B:668:TYR:O	1:B:706:SER:HA	2.14	0.47
1:B:681:ASP:O	1:B:684:GLU:HG2	2.15	0.47
1:A:108:LEU:HD22	1:A:186:TRP:O	2.15	0.46
1:B:648:VAL:O	1:B:649:ARG:C	2.56	0.46
1:B:662:ASN:O	1:B:663:VAL:C	2.57	0.46
1:B:726:LEU:HD11	1:B:729:LEU:HD13	1.97	0.46
1:A:194:ASN:ND2	1:A:197:LEU:H	2.09	0.46
1:A:89:ILE:O	1:A:89:ILE:HG22	2.15	0.46
1:A:150:GLU:OE2	1:A:156:PRO:HG3	2.16	0.46
1:A:220:TYR:HB3	1:A:222:LEU:HB2	1.97	0.46
1:B:717:ALA:HB1	1:B:738:ASP:OD1	2.15	0.46
1:B:743:GLN:HG2	1:B:747:GLY:C	2.40	0.46
1:B:709:LEU:HD21	1:B:741:GLY:C	2.40	0.46
1:B:587:PRO:O	1:B:590:PRO:HD3	2.17	0.45
1:B:739:ILE:O	1:B:742:ILE:HG22	2.16	0.45
1:A:117:VAL:O	1:A:121:LEU:HG	2.17	0.45
4:B:901:MM3:H161	4:B:901:MM3:H192	1.85	0.45
4:B:901:MM3:HNA	4:B:901:MM3:H202	1.82	0.45
1:A:125:GLU:HG3	1:A:130:LEU:O	2.16	0.45
1:A:148:VAL:HG12	1:A:149:ARG:HG2	1.98	0.45
1:B:655:TYR:HA	1:B:656:PRO:HD2	1.92	0.44
1:B:691:THR:O	1:B:691:THR:CG2	2.66	0.44
1:A:108:LEU:HG	1:A:112:ALA:CB	2.47	0.44
1:B:628:THR:HB	1:B:629:PRO:HD2	2.00	0.44
1:B:728:ASP:C	1:B:730:THR:N	2.73	0.44
1:B:654:PHE:CD1	1:B:654:PHE:N	2.85	0.43
1:A:194:ASN:ND2	1:A:196:PHE:HB3	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:707:LEU:HD23	1:B:707:LEU:HA	1.75	0.43
1:B:605:THR:HA	1:B:606:PRO:HD3	1.90	0.43
1:A:182:ASP:C	1:A:184:GLU:N	2.74	0.42
1:A:198:VAL:O	1:A:202:GLU:HG2	2.20	0.42
1:A:119:LYS:O	1:A:123:VAL:HG23	2.18	0.42
1:B:596:HIS:CD2	1:B:631:THR:OG1	2.72	0.42
1:B:628:THR:HB	1:B:629:PRO:CD	2.50	0.42
1:A:86:PHE:HB3	1:A:87:PRO:HD2	2.02	0.41
1:A:107:ASP:OD2	1:A:185:GLN:HA	2.20	0.41
1:A:222:LEU:O	1:A:222:LEU:HG	2.19	0.41
1:A:227:THR:C	1:A:228:ASP:O	2.60	0.41
1:B:642:ILE:HG23	1:B:676:GLY:O	2.20	0.41
1:B:597:LEU:HD12	1:B:597:LEU:N	2.36	0.41
1:A:222:LEU:HG	1:A:224:HIS:ND1	2.35	0.41
1:A:121:LEU:CD2	1:A:203:ILE:HD13	2.51	0.41
1:A:163:VAL:HG13	4:A:401:MM3:H203	2.02	0.41
1:A:224:HIS:HE2	1:A:227:THR:CG2	2.24	0.41
1:B:720:TYR:CD1	1:B:721:PRO:HD2	2.55	0.41
1:B:742:ILE:HD12	1:B:742:ILE:HA	1.90	0.41
1:A:123:VAL:CG1	1:A:234:LEU:CG	2.99	0.41
1:B:649:ARG:HA	1:B:657:PHE:O	2.21	0.41
1:B:681:ASP:HB3	1:B:684:GLU:HG2	2.02	0.41
1:A:104:TYR:HB3	5:A:405:HOH:O	2.19	0.41
1:A:121:LEU:HD23	1:A:203:ILE:HD13	2.02	0.40
1:A:249:PRO:HA	1:A:250:PRO:HD3	1.86	0.40
1:B:589:ILE:HD13	1:B:589:ILE:HG21	1.87	0.40
1:A:198:VAL:HG13	4:A:401:MM3:C5	2.51	0.40
1:B:752:SER:HA	1:B:753:PRO:HD3	1.87	0.40
1:B:658:ASP:O	1:B:659:GLY:C	2.65	0.40
1:A:86:PHE:HD1	1:A:87:PRO:HD2	1.87	0.40
1:B:671:GLY:C	1:B:675:ASN:HB2	2.46	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	167/173 (96%)	147 (88%)	18 (11%)	2 (1%)	10	23
1	B	171/173 (99%)	149 (87%)	20 (12%)	2 (1%)	10	23
All	All	338/346 (98%)	296 (88%)	38 (11%)	4 (1%)	10	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	LEU
1	A	221	PRO
1	B	649	ARG
1	B	750	PRO

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	143/147 (97%)	141 (99%)	2 (1%)	59	81
1	B	147/147 (100%)	137 (93%)	10 (7%)	14	32
All	All	290/294 (99%)	278 (96%)	12 (4%)	27	54

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	ARG
1	A	182	ASP
1	B	587	PRO
1	B	593	ARG
1	B	619	LYS
1	B	685	GLN
1	B	724	HIS
1	B	726	LEU
1	B	729	LEU

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Mol	Chain	Res	Type
1	B	740	ASN
1	B	745	LEU
1	B	755	THR

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

Mol	Chain	Res	Type
1	A	96	HIS
1	A	194	ASN
1	B	596	HIS
1	B	736	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 10 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	MM3	A	401	2	22,25,25	9.14	13 (59%)	27,37,37	5.02	16 (59%)
4	MM3	B	901	2	22,25,25	9.13	13 (59%)	27,37,37	3.99	18 (66%)

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
4	MM3	A	401	2	-	5/18/38/38	0/1/2/2
4	MM3	B	901	2	-	5/18/38/38	0/1/2/2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	401	MM3	O10-S9	32.51	1.78	1.43
4	B	901	MM3	O10-S9	31.35	1.77	1.43
4	B	901	MM3	O11-S9	25.51	1.70	1.43
4	A	401	MM3	O11-S9	23.81	1.68	1.43
4	B	901	MM3	S9-N12	8.00	1.75	1.63
4	A	401	MM3	S9-N12	7.57	1.74	1.63
4	A	401	MM3	OA-NA	4.94	1.51	1.40
4	A	401	MM3	C3-C2	4.66	1.46	1.38
4	A	401	MM3	C3-C4	4.45	1.47	1.38
4	A	401	MM3	C2-C1	4.39	1.45	1.38
4	B	901	MM3	C2-C1	4.38	1.45	1.38
4	B	901	MM3	C3-C4	4.31	1.46	1.38
4	A	401	MM3	C5-C4	4.11	1.46	1.38
4	B	901	MM3	C6-C5	3.99	1.45	1.38
4	B	901	MM3	C3-C2	3.86	1.45	1.38
4	A	401	MM3	C6-C1	3.81	1.44	1.38
4	A	401	MM3	C6-C5	3.73	1.44	1.38
4	B	901	MM3	C6-C1	3.68	1.44	1.38
4	B	901	MM3	C5-C4	3.67	1.45	1.38
4	B	901	MM3	C14-S15	-3.39	1.80	1.85
4	A	401	MM3	O7-C8	-3.15	1.34	1.42
4	B	901	MM3	O7-C8	-3.03	1.34	1.42
4	B	901	MM3	OA-NA	2.88	1.46	1.40
4	B	901	MM3	C16-S15	-2.44	1.80	1.81
4	A	401	MM3	C1-S9	2.34	1.79	1.76
4	A	401	MM3	C16-S15	2.07	1.83	1.81

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	MM3	C20-C14-C19	-14.88	96.79	111.17
4	A	401	MM3	C16-S15-C14	13.63	122.87	104.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	901	MM3	O10-S9-C1	-8.22	97.88	108.10
4	B	901	MM3	C20-C14-S15	-7.78	88.15	106.76
4	B	901	MM3	O11-S9-O10	-6.31	109.74	119.59
4	A	401	MM3	O11-S9-O10	-6.12	110.05	119.59
4	A	401	MM3	C2-C1-S9	-5.99	113.79	119.73
4	B	901	MM3	C18-N12-S9	-5.68	106.87	118.55
4	A	401	MM3	C1-S9-N12	5.60	117.13	107.36
4	B	901	MM3	C6-C1-C2	5.50	127.64	120.47
4	A	401	MM3	C6-C1-C2	5.34	127.42	120.47
4	B	901	MM3	C20-C14-C19	5.21	116.20	111.17
4	A	401	MM3	C5-C6-C1	-5.18	114.41	119.44
4	B	901	MM3	C1-S9-N12	4.99	116.06	107.36
4	B	901	MM3	O11-S9-N12	4.98	115.67	106.97
4	B	901	MM3	C6-C1-S9	-4.69	115.08	119.73
4	A	401	MM3	O11-S9-N12	-4.49	99.12	106.97
4	A	401	MM3	C18-N12-S9	-4.03	110.26	118.55
4	B	901	MM3	C2-C1-S9	-4.00	115.76	119.73
4	B	901	MM3	C3-C2-C1	-3.92	115.63	119.44
4	B	901	MM3	C16-S15-C14	3.82	109.61	104.44
4	A	401	MM3	O11-S9-C1	3.79	112.81	108.10
4	B	901	MM3	C5-C6-C1	-3.58	115.96	119.44
4	A	401	MM3	C8-O7-C4	3.25	124.47	117.50
4	A	401	MM3	C2-C3-C4	-2.96	116.35	119.73
4	B	901	MM3	C5-C4-C3	2.88	124.35	120.16
4	A	401	MM3	C13-C-NA	2.76	120.06	114.19
4	B	901	MM3	C19-C14-S15	-2.75	100.18	106.76
4	A	401	MM3	O10-S9-N12	2.62	111.55	106.97
4	A	401	MM3	C5-C4-C3	2.49	123.79	120.16
4	B	901	MM3	C13-C-NA	2.38	119.25	114.19
4	A	401	MM3	C16-C17-C18	-2.32	110.71	115.31
4	B	901	MM3	C17-C16-S15	-2.28	111.79	114.09
4	B	901	MM3	C16-C17-C18	-2.16	111.01	115.31

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	401	MM3	NA-C-C13-N12
4	A	401	MM3	C13-C-NA-OA
4	A	401	MM3	OB-C-NA-OA
4	B	901	MM3	C13-N12-S9-O11
4	B	901	MM3	NA-C-C13-N12

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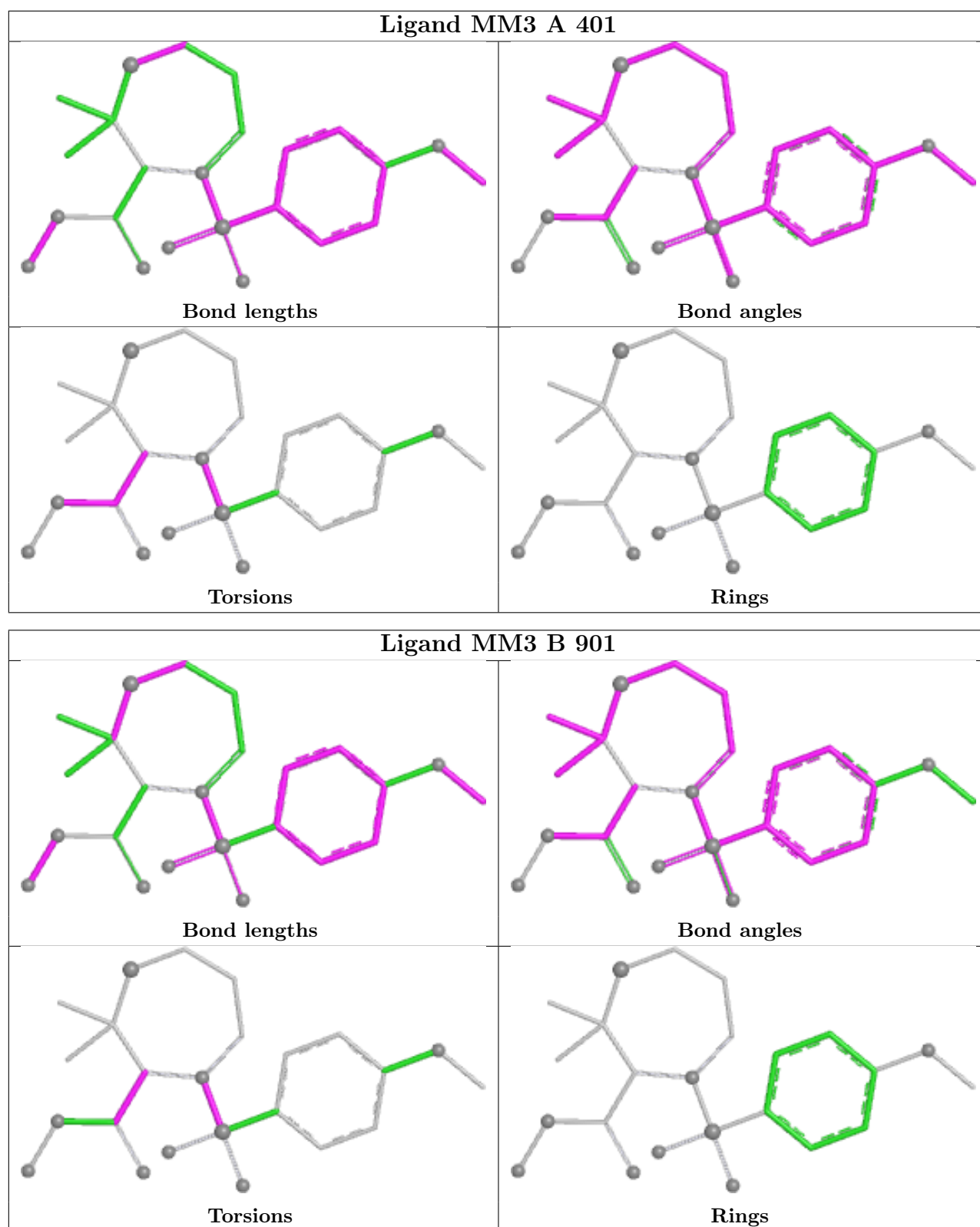
Mol	Chain	Res	Type	Atoms
4	B	901	MM3	C13-N12-S9-C1
4	A	401	MM3	C13-N12-S9-O11
4	A	401	MM3	OB-C-C13-N12
4	B	901	MM3	OB-C-C13-N12
4	B	901	MM3	C18-N12-S9-O11

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	MM3	2	0
4	B	901	MM3	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.



5.7 Other polymers ⓘ

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.