

## Supporting Information

### Emac – A Comparative Index for the Assessment of Macrocyclization Efficiency

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#### Contents

Part 1: Full data for literature survey (2009, 2010 and 2011) Page 1-20

Part 2: Full reference list Page 20-39

#### Part 1: Full data for literature survey (2009, 2010 and 2011)

2009										
Name	Journal	Paper type	Reaction	Cmpd no	Ring size	Yield	mmol	Vol (ml)	Conc (mM)	Emac
Abell	J. Org. Chem.	Methodology	RCM	1	16	48			10.00	6.044
Abell	J. Org. Chem.	Methodology	RCM	2	17	91			10.00	6.877
Abell	J. Org. Chem.	Methodology	RCM	3	18	100			10.00	7.000
Abell	J. Org. Chem.	Methodology	RCM	4	18	39			10.00	5.773
Abell	J. Org. Chem.	Methodology	RCM	5	19	33			10.00	5.556
Abell	J. Org. Chem.	Methodology	RCM	6	18	40			10.00	5.806
Abell	ACIE	Med Chem	RCM	1a	16	24			10.00	5.141
Abell	ACIE	Med Chem	RCM	2a	17	22			10.00	5.027
Abell	ACIE	Med Chem	RCM	3a	18	29			10.00	5.387
Abell	ACIE	Med Chem	RCM	4a	19	66			10.00	6.459
Altamura	BMCL	Med Chem	Lactam	18	24	42	0.270	55	4.909	5.561
Altamura	BMCL	Med Chem	Lactam	18	24	25	0.160	35	4.571	4.854
Baird	J. Org. Chem.	Tot Syn	RCM	19	15	86	0.099	20	4.950	6.498
Batenburg	Tetrahedron	Med Chem	Heck	3	17	15	0.080	2	40.000	5.130
Battistini	OBC	Med Chem	Lactam	14	13	66	4.510	1290	3.50	6.002
Beierle	ACIE	Med Chem	Huisgen	1	13	55	0.045	230	0.196	4.513
Beierle	ACIE	Med Chem	Huisgen	2	13	23			0.20	3.386
Beierle	ACIE	Med Chem	Huisgen	3	14	8	0.050	100	0.50	2.408
Beierle	ACIE	Med Chem	Huisgen	4	14	66			0.50	5.158
Bindl	Chem. Eur. J.	Tot Syn	RCAM	34	12	87	0.200	23	8.70	6.758
Bindl	Chem. Eur. J.	Tot Syn	RCAM	62	12	85	0.200	23	8.70	6.728
Bindl	Chem. Eur. J.	Tot Syn	RCAM	78	12	86	0.200	23	8.70	6.743
Bitar	Org. Lett.	Tot Syn	Alkylation	21	13	45			0.97	4.946
Blackburn	Tetrahedron	Tot Syn	Alkylation	28	12	85			50.00	7.487
Bourcet	Tet. Lett.	Tot Syn	RCM	8a	14	86			2.00	6.105
Bourcet	Tet. Lett.	Tot Syn	RCM	8b	14	83			2.00	6.058
Bowers	JACS	Med Chem	Lactam	10	16	53	0.015	15	1.00	5.173
Bowers	JACS	Med Chem	Lactam	30	16	62	0.110	100	1.10	5.419
Bowers	Org. Lett.	Med Chem	Lactam	9a	16	36	0.088	93.1	0.95	4.644
Bowers	Org. Lett.	Med Chem	Lactam	22a	16	40			1.00	4.806

<b>Bowers</b>	Org. Lett.	Med Chem	Lactam	29a	16	64			1.00	5.419
<b>Bowers</b>	Org. Lett.	Med Chem	Lactam	44	16	77			1.00	5.659
<b>Bowers</b>	Org. Lett.	Med Chem	Lactam	50	16	43			1.00	4.900
<b>Braddock</b>	J. Org. Chem.	Tot Syn	RCM	15	11	34	0.200	88	2.27	4.951
<b>Braddock</b>	J. Org. Chem.	Tot Syn	RCM	16	11	33	0.032	14	2.29	4.915
<b>Bruder</b>	OBC	Tot Syn	Mitsunobu	4	16	27			8.00	5.197
<b>Burns</b>	JACS	Tot Syn	Alkylation	7	11	79			2.00	5.994
<b>Calo</b>	Org. Lett.	Tot Syn	RCM	30	14	83	0.024	41	0.59	5.525
<b>Campbell</b>	Tet. Lett.	Methodology	Lactam	2	18	60			18.00	6.590
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3e	11	6			3.00	2.812
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3f	12	30			3.00	4.908
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3g	13	48			3.00	5.521
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3h	14	51			3.00	5.600
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3j	13	31			3.00	4.951
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3k	13	39			3.00	5.250
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3l	13	32			3.00	4.993
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3m	16	44			3.00	5.407
<b>Castagnolo</b>	EJOC	Med Chem	Guanidine synth	3n	16	42			3.00	5.347
<b>Chandrasekhar</b>	Tet. Lett.	Tot Syn	Lactam	1	19	60			0.05	4.033
<b>Chandrasekhar</b>	Tet. Lett.	Tot Syn	Lactam	2	19	58			0.05	3.989
<b>Cribiu</b>	ACIE	Tot Syn	RCM	24	15	76	0.067	130	0.52	5.355
<b>Crimmins</b>	Org. Lett.	Tot Syn	Lactone		20	74	0.004	303	0.014	3.752
<b>Custar</b>	JACS	Tot Syn	Prins	30	12	25	0.050	5	10.000	5.194
<b>Custar</b>	JACS	Tot Syn	Prins	41	12	25	0.050	5	10.000	5.194
<b>Czajkowska</b>	Tet. Lett.	Med Chem	RCM	7	22	84			0.20	5.074
<b>Dakas</b>	Chem. Eur. J.	Tot Syn	RCM	49	14	94	0.400	300	1.333	6.044
<b>Di Francesco</b>	J. Med. Chem.	Med Chem	RCM	12	18	55	0.190	20	9.500	6.199
<b>Di Francesco</b>	J. Med. Chem.	Med Chem	RCM	21a	18	55	0.370	7	52.86	6.944
<b>Dietrich</b>	Chem. Eur. J.	Tot Syn	Lactone	31	16	57	0.023	10	2.30	5.629
<b>Dietrich</b>	Chem. Eur. J.	Tot Syn	Lactone	37	16	83	0.320	358	0.89	5.709
<b>Dietrich</b>	Chem. Eur. J.	Tot Syn	Lactone	40	16	75	1.040	802	1.30	5.738
<b>Eissler</b>	Chem. Eur. J.	Tot Syn	RCM	22a	16	85	0.650	65	10.00	6.788
<b>Eissler</b>	Chem. Eur. J.	Tot Syn	RCM	22b	16	71	0.650	65	10.00	6.554
<b>Eissler</b>	Chem. Eur. J.	Tot Syn	RCM	22c	16	75	0.650	65	10.00	6.625
<b>Eissler</b>	Chem. Eur. J.	Tot Syn	RCM	22d	16	76	0.650	65	10.00	6.642
<b>Evans</b>	JACS	Tot Syn	Lactone	20	16	68			0.90	5.452
<b>Farina</b>	OPRD	Process	RCM	7	15	93			200.00	8.206
<b>Faure</b>	Org. Lett.	Tot Syn	Lactam	15	19	52	0.008	6	1.37	5.284
<b>Feng</b>	JACS	Methodology	Lactam	22c	77	70			4.10	6.148
<b>Feng</b>	JACS	Methodology	Lactam	23c	88	84			6.70	6.599
<b>Frein</b>	Org. Lett.	Tot Syn	Lactone	16	16	81			0.30	5.203
<b>Furstner</b>	ACIE	Tot Syn	RCM	67	26	72	0.016	20	0.80	5.475
<b>Furstner</b>	ACIE	Tot Syn	RCM	81	26	84			0.80	5.676
<b>Furstner</b>	ACIE	Tot Syn	RCM	95	26	91			0.80	5.780
<b>Furstner</b>	ACIE	Tot Syn	RCM	100	26	56			0.80	5.148
<b>Furstner</b>	ACIE	Tot Syn	RCM	110	26	54			0.80	5.100
<b>Furstner</b>	Chem. Eur. J.	Tot Syn	RCAM	27	14	84	0.169	599	0.28	5.223

<b>Furstner</b>	Chem. Eur. J.	Tot Syn	RCAM	37	14	89		0.28	5.295	
<b>Furstner</b>	Chem. Eur. J.	Tot Syn	RCAM	45	14	75		0.28	5.072	
<b>Furstner</b>	Chem. Eur. J.	Tot Syn	RCAM	53	14	93		0.28	5.353	
<b>Garfunkle</b>	JACS	Tot Syn	Larock Indole	21	16	89		0.85	5.778	
<b>Gentilucci</b>	J. Med. Chem.	Med Chem	Lactam	14	13	69	0.310	70	4.43	6.163
<b>Gesinski</b>	Org. Lett.	Methodology	Prins	13	18	51			50	6.822
<b>Gesinski</b>	Org. Lett.	Methodology	Prins	28	16	43			30	6.378
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	14a	15	88	0.047	15	3.13	6.329
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	14b	14	50			3.13	5.592
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	14c	13	89			3.13	6.344
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	14d	12	71			3.13	6.049
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	14e	12	52			3.13	5.644
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	14f	11	81			3.13	6.221
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	20a	14	96	0.140	90	1.56	6.139
<b>Ghosh</b>	J. Med. Chem.	Med Chem	RCM	21b	14	97	0.170	125	1.36	6.094
<b>Ghosh</b>	Org. Lett.	Tot Syn	Lactam	13	12	95	0.036	36	1.00	5.933
<b>Ghosh</b>	Org. Lett.	Tot Syn	Lactam	19	12	55	0.082	40.8	2.01	5.524
<b>Ghosh, P</b>	Org. Lett.	Tot Syn	Lactone	2	14	77	2.000	200	10.00	6.659
<b>Gilles</b>	J. Org. Chem.	Tot Syn	Lactam	26	25	45			0.80	4.863
<b>Goldring</b>	Tetrahedron	Tot Syn	Radical	42	12	80			2.40	6.089
<b>Gollner</b>	Chem. Eur. J.	Tot Syn	Lactone	45	18	75	0.091	122	0.74	5.496
<b>Gollner</b>	Chem. Eur. J.	Tot Syn	Lactone	49	18	35	0.047	221	0.21	3.957
<b>Gollner</b>	Chem. Eur. J.	Tot Syn	Lactone	62	18	32	0.020	140	0.14	3.666
<b>Grauer</b>	Chem. As. J.	Methodology	Ullmann	2	17	7	0.070	0.7	100.00	4.535
<b>Grauer</b>	Chem. As. J.	Methodology	Ullmann	4	15	40	0.180	7.2	25.00	6.204
<b>Grauer</b>	Chem. As. J.	Methodology	Ullmann	50	15	19	0.140	5.6	25.00	5.234
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	25	12	62			1.00	5.377
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	27	12	74			1.00	5.608
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	28	12	35			1.00	4.632
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	29	12	65			1.00	5.439
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	30	13	72			1.00	5.572
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	31	14	72			1.00	5.572
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	32	16	34			1.00	4.594
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	33	17	54			1.00	5.197
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	34	13	36			1.00	4.669
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	35	14	68			1.00	5.498
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	36	16	61			1.00	5.356
<b>Grimwood</b>	Tetrahedron	Methodology	RCM Enyne	37	18	67			1.00	5.478
<b>Guinchard</b>	Org. Lett.	Tot Syn	Lactone	22	12	85	0.076	102	0.74	5.658
<b>Han</b>	Org. Lett.	Methodology	Lactam	3b	14	73			600.00	8.368
<b>Han</b>	Org. Lett.	Methodology	Lactam	3a	14	80			600.00	8.487
<b>Han</b>	Org. Lett.	Methodology	Lactam	3c	14	54			600.00	7.975
<b>Han</b>	Org. Lett.	Methodology	Lactam	3d	14	60			600.00	8.113
<b>Harper</b>	J. Med. Chem.	Med Chem	RCM	10c	19	37	0.700	45	15.56	5.896
<b>Harper</b>	J. Med. Chem.	Med Chem	RCM	10d	19	27	0.700	45	15.56	5.486
<b>Harper</b>	J. Med. Chem.	Med Chem	RCM	10e	19	31	0.700	45	15.56	5.666
<b>Harper</b>	J. Med. Chem.	Med Chem	RCM	10f	19	78	0.700	45	15.56	6.868

Harper	J. Med. Chem.	Med Chem	RCM	10g	19	66	0.700	45	15.56	6.651
Harper	J. Med. Chem.	Med Chem	RCM	10l	19	12	0.700	45	15.56	4.429
Hioki	BMCL	Med Chem	Suzuki	16	18	39			5.00	5.472
Hjelmgaard	Org. Lett.	Med Chem	Lactam	8b	14	82			5.00	6.440
Hjelmgaard	Org. Lett.	Med Chem	Lactam	8c	21	63			5.00	6.097
Hjelmgaard	Org. Lett.	Med Chem	Lactam	8d	28	74			5.00	6.307
Hjelmgaard	Org. Lett.	Med Chem	Lactam	8e	35	70			5.00	6.234
Horne	ACIE	Med Chem	Huisgen	2	13	50	0.045	225	0.20	4.398
Houghton	J. Org. Chem.	Tot Syn	Alkylation	22	20	93			10.00	6.905
Katagiri	J. Org. Chem.	Methodology	Lactam	5	38	55			50.00	6.920
Katagiri	J. Org. Chem.	Methodology	Lactam	6	40	19			50.00	5.535
Kelly	Org. Lett.	Methodology	Huisgen	4a	11	58			2.00	5.591
Kelly	Org. Lett.	Methodology	Huisgen	5a	12	55			10.00	6.221
Kelly	Org. Lett.	Methodology	Huisgen	4b	12	67			2.00	5.779
Kelly	Org. Lett.	Methodology	Huisgen	5b	13	56			10.00	6.245
Kelly	Org. Lett.	Methodology	Huisgen	4c	11	45			2.00	5.261
Kelly	Org. Lett.	Methodology	Huisgen	5c	12	38			10.00	5.739
Kelly	Org. Lett.	Methodology	Huisgen	4d	12	58			2.00	5.591
Kelly	Org. Lett.	Methodology	Huisgen	5d	13	40			10.00	5.806
Kelly	Org. Lett.	Methodology	Huisgen	4e	11	60			2.00	5.635
Kelly	Org. Lett.	Methodology	Huisgen	5e	12	37			10.00	5.705
Kelly	Org. Lett.	Methodology	Huisgen	4f	12	54			2.00	5.498
Kelly	Org. Lett.	Methodology	Huisgen	5f	13	40			10.00	5.806
Kelly	Org. Lett.	Methodology	Huisgen	4g	12	61			2.00	5.657
Kelly	Org. Lett.	Methodology	Huisgen	5g	13	48			10.00	6.044
Kelly	Org. Lett.	Methodology	Huisgen	4h	12	64			2.00	5.720
Kelly	Org. Lett.	Methodology	Huisgen	5h	13	31			10.00	5.474
Kelly	Org. Lett.	Methodology	Huisgen	4i	12	48			2.00	5.345
Kelly	Org. Lett.	Methodology	Huisgen	5i	13	65			10.00	6.439
Kelly	Org. Lett.	Methodology	Huisgen	4j	12	71			2.00	5.855
Kelly	Org. Lett.	Methodology	Huisgen	5j	13	17			10.00	4.691
Kelly	Org. Lett.	Methodology	Huisgen	5j	13	46			2.00	5.289
Kelly	Org. Lett.	Methodology	Huisgen	4k	12	70			2.00	5.836
Kelly	Org. Lett.	Methodology	Huisgen	5k	13	62			10.00	6.377
Kim	ACIE	Tot Syn	Lactone	33	12	69	0.031	35	0.89	5.464
Kim	ACIE	Tot Syn	Lactone	22	15	90	0.070	15	4.67	6.532
Kojima	Tet. Lett.	Tot Syn	Lactone		22	90			1.50	6.039
Kong	ACIE	Tot Syn	Barbier	19	15	56			6.70	6.071
Kumar	EJOC	Med Chem	Lactam	21	18	58			1.00	5.290
Lee	Tetrahedron	Tot Syn	Lactam	4	27	76			10.00	6.642
LeFranc	ACIE	Tot Syn	Lactam	1	26	41			85.00	6.768
Ley	Chem. Eur. J.	Tot Syn	Alkylation	145	35	70			1.00	5.535
Li	JACS	Tot Syn	Heck	27	22	70			4.00	6.137
Li	Org. Lett.	Tot Syn	Lactam	5	13	87	0.105	105	1.00	5.819
Li	ACIE	Tot Syn	Alkylation	30	18	79	0.045	30	1.50	5.869
Li	J. Org. Chem.	Methodology	SNAr	6a	14	92			1.00	5.891
Li	J. Org. Chem.	Methodology	SNAr	6b	14	77			1.40	5.806

Li	J. Org. Chem.	Methodology	SNAr	6c	14	89		1.30	5.962	
Li	J. Org. Chem.	Methodology	SNAr	6d	14	88		1.10	5.875	
Li	J. Org. Chem.	Methodology	SNAr	6e	14	80		1.20	5.788	
Li	J. Org. Chem.	Methodology	SNAr	6f	14	85		1.00	5.788	
Li	J. Org. Chem.	Methodology	SNAr	6g	15	78		1.00	5.676	
Li	J. Org. Chem.	Methodology	SNAr	6h	15	70		1.20	5.614	
Li	Tetrahedron	Tot Syn	Lactam	1	19	76		1.00	5.642	
Li	Tetrahedron	Tot Syn	Lactam	2	19	75		1.00	5.625	
Li	Tetrahedron	Tot Syn	Lactam	19	19	73		1.00	5.590	
Liu	J. Org. Chem.	Tot Syn	Lactone	7.4	15	58	0.006	11.2	0.56	5.040
Llacer	Org. Lett.	Tot Syn	RCM	10	14	85		0.17	5.019	
Lu	Tetrahedron	Tot Syn	Lactone	15a	12	73	0.150	28.4	5.28	6.313
Lu	Tetrahedron	Tot Syn	Lactone	15b	12	78	0.410	72	5.69	6.432
Luo	JACS	Methodology	RCM	5h	12	99		5.00	6.686	
Luo	JACS	Methodology	RCM	5ia	12	70		5.00	6.234	
Luo	JACS	Methodology	RCM	5ib	14	53		5.00	5.872	
Luo	JACS	Methodology	RCM	5ja	14	74		5.00	6.307	
Luo	JACS	Methodology	RCM	5jb	14	77		5.00	6.358	
Luo	JACS	Methodology	RCM	5jc	15	86		5.00	6.502	
Luo	JACS	Methodology	RCM	5jd	15	97		5.00	6.659	
Luo	JACS	Methodology	RCM	5k	15	69		5.00	6.216	
Luo	JACS	Methodology	RCM	15ica	14	77		5.00	6.358	
Luo	JACS	Methodology	RCM	15icb	14	72		5.00	6.271	
Luo	JACS	Methodology	RCM	15icc	14	81		5.00	6.424	
Ma	JACS	Tot Syn	Lactam	32	17	90	0.872	35	24.91	7.259
Ma	JACS	Tot Syn	Lactam	57b	17	91	0.048	10	4.84	6.562
Ma	JACS	Tot Syn	Lactam	63	17	83	0.005	2	2.70	6.189
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	9	26	15		0.15	2.704	
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	10	31	15		0.15	2.704	
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	11	28	41		0.15	4.014	
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	12	32	38		0.15	3.915	
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	13	28	64		0.15	4.595	
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	14	28	64		0.15	4.595	
Madden	Chem. Comm.	Methodology	1,3-cycloaddition	15	28	94		0.15	5.095	
Magauer	ACIE	Tot Syn	Lactone	3	20	55		1.20	5.300	
Mahoney	J. Org. Chem.	Tot Syn	Lactam	24	21	94	0.260	9.9	26.26	7.339
Mandel	Org. Lett.	Tot Syn	RCM	22	18	43		1.00	4.900	
Mbere-Nguyen	Tetrahedron	Methodology	RCM	24	18	35		2.20	4.975	
McGowan	OBC	Tot Syn	Alkylation	42	12	68	0.190	1.5	126.67	7.600
McGowan	OBC	Tot Syn	Alkylation	62	12	45	0.083	1.6	51.88	6.675
Menche	J. Org. Chem.	Tot Syn	HWE	68	24	44	0.001	1.5	0.73	4.796
Menche	J. Org. Chem.	Tot Syn	Heck	2	24	60	0.004	2	1.82	5.593
Muri	J. Org. Chem.	Tot Syn	Lactone	33	14	55	0.059	27	2.19	5.561
Muri	J. Org. Chem.	Tot Syn	Lactone	46	14	78	0.214	40	5.35	6.405
Muthusamy	Tet. Lett.	Methodology	Rh O-H insertion	5a	16	75	1.000	52	19.23	6.909
Muthusamy	Tet. Lett.	Methodology	Rh O-H insertion	5b	15	78	1.000	52	19.23	6.960
Muthusamy	Tet. Lett.	Methodology	Rh O-H insertion	5c	17	70	1.000	52	19.23	6.819

<b>Muthusamy</b>	Tet. Lett.	Methodology	Rh O-H insertion	5d	19	72	1.000	52	19.23	6.856
<b>Muthusamy</b>	Tet. Lett.	Methodology	Rh O-H insertion	5e	21	65	1.000	52	19.23	6.723
<b>Nagano</b>	Chem. Eur. J.	Tot Syn	RCM	31	20	71	0.056	10	5.58	6.300
<b>Nagano</b>	Chem. Eur. J.	Tot Syn	RCM	46	20	96			5.58	6.693
<b>Nagano</b>	Chem. Eur. J.	Tot Syn	RCM	51	20	88			5.58	6.580
<b>Nagano</b>	Chem. Eur. J.	Tot Syn	RCM	57	20	81			5.58	6.472
<b>Nakatsuka</b>	Tet. Lett.	Tot Syn	Lactam	1	21	23			1.50	4.261
<b>Narita</b>	Chem. Eur. J.	Tot Syn	Lactone	37	15	90			0.89	5.812
<b>Narita</b>	Chem. Eur. J.	Tot Syn	Lactone	48	15	89			0.90	5.802
<b>Narita</b>	Chem. Eur. J.	Tot Syn	Lactone	68	15	62			0.90	5.331
<b>Newhouse</b>	JACS	Tot Syn	Lactam	11	16	70			0.10	4.535
<b>Nicolaou</b>	ACIE	Tot Syn	Alkylation	3	15	79			1.00	5.693
<b>Nicolaou</b>	ACIE	Tot Syn	DA	24	13	40			7.70	5.693
<b>Nolasco</b>	J. Org. Chem.	Tot Syn	Lactam	3	17	60	1.070	330	3.24	5.845
<b>Nolasco</b>	J. Org. Chem.	Tot Syn	Lactam	25	17	48	0.160	31	5.16	5.756
<b>Nolasco</b>	J. Org. Chem.	Tot Syn	Negishi	10	17	35	0.330	140	2.36	5.005
<b>Numajiri</b>	Chem. As. J.	Tot Syn	Lactam	2	25	56			1.00	5.245
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3a	13	64			0.50	5.118
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3b	9	61			0.30	4.833
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3c	10	50			0.30	4.574
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3d	11	68			0.30	4.975
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3e	12	62			0.50	5.076
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3f	13	69			0.50	5.216
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	3g	14	89			0.50	5.547
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	8a	13	74			0.30	5.085
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	8b	13	55			0.30	4.698
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	8c	13	67			0.30	4.955
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	9a	12	72			0.50	5.271
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	9b	12	52			0.50	4.847
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	9c	12	70			0.50	5.234
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	11a	13	85			0.50	5.487
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	11b	13	73			0.50	5.289
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	11c	13	75			0.50	5.324
<b>Ohba</b>	Chem. Eur. J.	Methodology	Lactone	15	10	45			0.30	4.437
<b>Pattenden</b>	Tetrahedron	Tot Syn	Radical	23,24,25	12	50			1.20	5.176
<b>Peng</b>	BMCL	Med Chem	RCM	17	12	63	0.098	9.758	10.00	6.398
<b>Peng</b>	BMCL	Med Chem	RCM	18a	12	39			10.00	5.773
<b>Peng</b>	BMCL	Med Chem	RCM	18b	12	65			10.00	6.439
<b>Peng</b>	BMCL	Med Chem	RCM	19a	13	12			10.00	4.238
<b>Peng</b>	BMCL	Med Chem	RCM	19b	13	11			10.00	4.124
<b>Peng</b>	BMCL	Med Chem	RCM	20a	14	34			10.00	5.594
<b>Peng</b>	BMCL	Med Chem	RCM	20b	14	21			10.00	4.967
<b>Peng</b>	BMCL	Med Chem	RCM	21a	12	46			10.00	5.988
<b>Peng</b>	BMCL	Med Chem	RCM	21b	12	42			10.00	5.870
<b>Peng</b>	BMCL	Med Chem	RCM	22a	13	34			10.00	5.594
<b>Peng</b>	BMCL	Med Chem	RCM	19	13	32			10.00	5.515
<b>Peng</b>	BMCL	Med Chem	RCM	23a	14	24			10.00	5.141

<b>Peng</b>	BMCL	Med Chem	RCM	23b	14	22			10.00	5.027
<b>Peng</b>	BMCL	Med Chem	RCM	24b	12	38			10.00	5.739
<b>Peng</b>	BMCL	Med Chem	RCM	20	13	56			10.00	6.245
<b>Peng</b>	BMCL	Med Chem	RCM	25b	13	48			10.00	6.044
<b>Peng</b>	BMCL	Med Chem	RCM	26a	14	27			10.00	5.294
<b>Peng</b>	BMCL	Med Chem	RCM	26b	14	31			10.00	5.474
<b>Postema</b>	Org. Lett.	Tot Syn	RCM	18	20	84			3.70	6.341
<b>Purushothaman</b>	Tet. Lett.	Methodology	1,3-cycloaddition	3a	12	70	1.000	30	33.33	7.058
<b>Rajakumar</b>	EJMC	Med Chem	Lactam	1	21	40			0.70	4.651
<b>Rajakumar</b>	EJMC	Med Chem	Lactam	2	21	32			0.70	4.361
<b>Rajakumar</b>	EJMC	Med Chem	Lactam	3	22	28			0.70	4.187
<b>Rajakumar</b>	EJMC	Med Chem	Lactam	4	25	37			0.70	4.550
<b>Rajakumar</b>	EJMC	Med Chem	Lactam	5	32	27			0.70	4.139
<b>Rajakumar</b>	EJMC	Med Chem	Lactam	6	26	32			0.70	4.361
<b>Reddy</b>	Org. Lett.	Tot Syn	RCM	24	15	83			5.00	6.456
<b>Rohanna</b>	Org. Lett.	Methodology	Lactone	5	13	75			1.60	5.829
<b>Rohanna</b>	Org. Lett.	Methodology	Lactone	7	17	90			1.20	5.942
<b>Rohanna</b>	Org. Lett.	Methodology	Lactone	9	16	43			0.60	4.679
<b>Rohanna</b>	Org. Lett.	Methodology	Lactone	11	17	90			0.90	5.817
<b>Rohanna</b>	Org. Lett.	Methodology	Lactone	17	13	81			0.40	5.328
<b>She</b>	Tet. Lett.	Tot Syn	RCM	20,21	14	69	0.164	30	5.47	6.254
<b>Singh</b>	J. Org. Chem.	Tot Syn	Lactone	16	16	77	0.132	204	0.65	5.470
<b>Smith</b>	JACS	Tot Syn	Lactone	19	31	85			0.50	5.487
<b>Smith</b>	J. Org. Chem.	Tot Syn	Heck	80	26	50			10.00	6.097
<b>Smith</b>	Org. Lett.	Tot Syn	HWE	24	21	59	0.020	34	0.59	5.082
<b>Smith</b>	JACS	Tot Syn	RCM	52	11	89			0.08	4.723
<b>Smith</b>	JACS	Tot Syn	RCM	68	11	86			0.06	4.603
<b>Smith</b>	Tetrahedon	Tot Syn	Lactone	101	32	85	0.001	1	0.73	5.649
<b>Smith</b>	Tetrahedon	Tot Syn	Lactone	114	32	81	0.002	4.3	0.52	5.440
<b>Sparling</b>	Tetrahedron	Tot Syn	HWE	3	14	34			9.00	5.549
<b>Sugiyama</b>	Chem. Eur. J.	Tot Syn	HWE	42	36	52	0.096	50	1.92	5.431
<b>Tang</b>	OBC	Tot Syn	NHK	4a	13	70	0.540	377	1.43	5.691
<b>Toumi</b>	EJOC	Tot Syn	Ullmann	36	13	82			7.40	6.611
<b>Toumi</b>	EJOC	Tot Syn	Cyclodehydration	4	13	34			7.40	5.464
<b>Toumi</b>	EJOC	Tot Syn	Pd-amination	4	13	21			10.00	4.967
<b>Trenkle</b>	ACIE	Tot Syn	Ni alkyne-epoxide	2	14	67			140.00	7.624
<b>Trenkle</b>	ACIE	Tot Syn	Ni alkyne-epoxide	2	14	46			70.00	6.833
<b>Trost</b>	JACS	Tot Syn	Ru alkene-alkyne	25	18	99			1.00	5.987
<b>Trost</b>	JACS	Tot Syn	Lactone	73	22	65			1.20	5.518
<b>Trost</b>	ACIE	Tot Syn	Lactone	1	16	26	0.009	4.5	1.96	4.536
<b>Tu</b>	ACIE	Tot Syn	Lactone	11	13	75	0.359	215	1.67	5.848
<b>Tu</b>	ACIE	Tot Syn	Lactone	39	12	72	0.421	305	1.38	5.712
<b>Ueda</b>	J. Org. Chem.	Tot Syn	Lactone	41	14	61	0.039	137.6	0.28	4.810
<b>Urbach</b>	EJOC	Med Chem	RCM	7	13	95			4.70	6.605
<b>Urbach</b>	EJOC	Med Chem	RCM	12b	12	74			5.00	6.307
<b>Urbach</b>	EJOC	Med Chem	RCM	12c	13	62			5.00	6.076
<b>Urbach</b>	EJOC	Med Chem	RCM	17b	12	83			5.00	6.456

Urbach	EJOC	Med Chem	RCM	17c	13	72			5.00	6.271
Urbach	EJOC	Med Chem	RCM	24	12	67			5.00	6.177
Van Berkel	Chem. Comm.	Tot Syn	Huisgen	16	23	23			0.80	3.988
Velazquez	J. Med. Chem.	Med Chem	RCM	11	16	90			9.70	6.849
Wingstrand	Tet. Lett.	Methodology	Alkylation	19	15	56	0.061	1.4	43.57	6.884
Wingstrand	Tet. Lett.	Methodology	Red Amin	20	15	48	0.102	2.27	44.93	6.696
Xing	Org. Lett.	Tot Syn	Lactone	12	12	70	0.170	70	2.43	5.921
Xing	Org. Lett.	Tot Syn	Lactone	15	12	70	0.630	270	2.33	5.903
Yadav	Org. Lett.	Tot Syn	Lactone	25	18	75			0.40	5.227
Yang	New J. Chem.	Methodology	Lactam	3a	30	85			0.50	5.487
Yang	New J. Chem.	Methodology	Lactam	3b	35	50			0.50	4.796
Yang	New J. Chem.	Methodology	Lactam	3c	40	14			0.50	3.137
Yang	New J. Chem.	Methodology	Lactam	3d	45	10			0.50	2.699
Yang	New J. Chem.	Methodology	Lactam	3e	50	6			0.50	2.033
Zhu	OBC	Methodology	Huisgen	1	25	20			10.00	4.903
Zhu	OBC	Methodology	Huisgen	2	34	82			10.00	6.741
Zhu	OBC	Methodology	Huisgen	3	34	85			5.00	6.487
Zhu	OBC	Methodology	Huisgen	4	54	25			5.00	4.893
Zulykama	Tet. Lett.	Methodology	Alkylation	3a	12	61			40.00	6.958
Zulykama	Tet. Lett.	Methodology	Alkylation	3b	12	65			40.00	7.041
Zulykama	Tet. Lett.	Methodology	Alkylation	3c	12	63			40.00	7.000
Zulykama	Tet. Lett.	Methodology	Alkylation	3d	12	61			40.00	6.958
Zulykama	Tet. Lett.	Methodology	Alkylation	3e	12	64			40.00	7.021
Zulykama	Tet. Lett.	Methodology	Alkylation	3f	12	55			40.00	6.823
Zulykama	Tet. Lett.	Methodology	Alkylation	3g	12	57			40.00	6.870
Zulykama	Tet. Lett.	Methodology	Alkylation	3h	12	53			40.00	6.775

## 2010

Name	Journal	Paper type	Reaction	Cmpd no	Ring size	Yield	mmol	Vol (ml)	Conc (mM)	Emac
Ammer	Chem. Eur. J.	Tot syn	Stille	38	29	79	0.010	9.9	1.00	5.692
Ando	Org. Lett.	Methodology	HWE	16	15	84	0.350	35	10.00	6.773
Ando	Org. Lett.	Methodology	HWE	17	13	60			10.00	6.334
Ando	Org. Lett.	Methodology	HWE	18	16	93			10.00	6.905
Ando	Org. Lett.	Methodology	HWE	19	18	77			10.00	6.659
Ando	Org. Lett.	Methodology	HWE	20	12	69			10.00	6.517
Ando	Org. Lett.	Methodology	HWE	21	14	87			10.00	6.819
Ando	Org. Lett.	Methodology	HWE	22	17	90			10.00	6.863
Auzzas	J. Med. Chem.	Med Chem	RCM	32	14	78			1.00	5.676
Auzzas	J. Med. Chem.	Med Chem	RCM	30a	12	58			1.00	5.290
Auzzas	J. Med. Chem.	Med Chem	RCM	30b	13	88			1.00	5.833
Auzzas	J. Med. Chem.	Med Chem	RCM	30c	14	89			1.00	5.848
Auzzas	J. Med. Chem.	Med Chem	RCM	39	14	26			1.00	4.245
Auzzas	J. Med. Chem.	Med Chem	RCM	46	14	55			1.00	5.221
Auzzas	J. Med. Chem.	Med Chem	RCM	54	14	25	0.210	210	1.00	4.194
Auzzas	J. Med. Chem.	Med Chem	RCM	61	14	72	0.150	150	1.00	5.572
Barnickel	J. Org. Chem.	Tot syn	Alkylation	8a	15	64	0.280	36	7.78	6.309

<b>Barnickel</b>	J. Org. Chem.	Tot syn	Alkylation	8b	17	76	0.500	60	8.33	6.563
<b>Barnickel</b>	J. Org. Chem.	Tot syn	Alkylation	4	17	25	0.200	24	8.33	5.115
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6a	12	73	0.010	0.6	16.67	6.812
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6b	13	66			16.67	6.681
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6c	14	80			16.67	6.931
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6d	12	87			16.67	7.040
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6e	14	90			16.67	7.085
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6f	13	35			16.67	5.854
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6g	12	54			16.67	6.419
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6h	13	37			16.67	5.927
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6i	12	37			16.67	5.927
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6j	14	74			16.67	6.830
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6k	14	65			16.67	6.661
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6l	16	62			16.67	6.599
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6m	18	65			16.67	6.661
<b>Bogdan</b>	Chem. Eur. J.	Methodology	Huisgen	6n	22	71			16.67	6.776
<b>Bolduc</b>	JACS	Methodology	RCM	9	18	89			0.20	5.149
<b>Bolduc</b>	JACS	Methodology	RCM	9	18	67			0.40	5.080
<b>Bolduc</b>	JACS	Methodology	RCM	9	18	61			0.80	5.259
<b>Bolduc</b>	JACS	Methodology	RCM	5	16	50	0.089	600	0.15	4.268
<b>Bolduc</b>	JACS	Methodology	RCM	6	16	45	0.089	600	0.15	4.131
<b>Bolduc</b>	JACS	Methodology	RCM	7	15	61	0.089	600	0.15	4.527
<b>Bolduc</b>	JACS	Methodology	RCM	8	17	45	0.089	600	0.15	4.131
<b>Bolduc</b>	JACS	Methodology	RCM	10	15	45	0.089	600	0.15	4.131
<b>Bolduc</b>	JACS	Methodology	Glaser-Hay	11	15	42	0.089	700	0.13	3.974
<b>Bolduc</b>	JACS	Methodology	Glaser-Hay	12	15	40	0.089	700	0.13	3.910
<b>Bolduc</b>	JACS	Methodology	Glaser-Hay	13	18	42	0.089	700	0.13	3.974
<b>Bourcet</b>	EJOC	Tot syn	RCM	57	14	67	0.829	160	5.18	6.193
<b>Brain</b>	Tet	Tot syn	Stille	83	17	52	0.019	12	1.62	5.356
<b>Burns</b>	ACIE	Tot syn	RCM	9	13	82	0.940	320	2.94	6.209
<b>Campbell</b>	Tet. Lett.	Tot syn	Lactam	1a	18	44	0.180	10	18.00	6.186
<b>Carr</b>	Chem Comm	Tot syn	Heck	12a	11	52	0.070	2	35.00	6.692
<b>Chausset</b>	Chem Comm	Tot syn	RCM	1	12	37	0.040	9	4.44	5.352
<b>Clerc</b>	EJOC	Tot syn	RCM	21	12	49	1.620	825	1.96	5.364
<b>Cupido</b>	ACIE	Med Chem	Lactam		15	88			0.05	4.532
<b>Dai</b>	Org. Lett.	Tot syn	Lactam	2	12	15	0.520	200	2.60	3.943
<b>Day</b>	Chem. Eur. J.	Med Chem	RCM	14	14	92	0.280	39	7.18	6.747
<b>Day</b>	Chem. Eur. J.	Med Chem	Huisgen	20a	13	33	0.280	180	1.56	4.747
<b>Doi</b>	OBC	Med Chem	Lactam	6	24	19			3.13	4.332
<b>Dufour</b>	Chem. Eur. J.	Tot syn	Suzuki	24	14	54	0.074	3.72	19.92	6.496
<b>Feng</b>	Org. Lett.	Methodology	C-H arylation	8	12	64			0.005	3.118
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	24	14	82	0.339	68	4.99	6.439
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	23	14	36	0.043	8.7	4.99	5.367
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	25	14	43	0.046	9.2	5.01	5.600
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	26	14	60	0.072	14	5.14	6.046
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	27	14	29	0.086	17	5.05	5.091
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	28	14	72	0.625	125	5.00	6.271

<b>Fuse</b>	Chem. As. J.	Analogues	RCM	29	14	71	0.023	4.6	5.00	6.253
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	30	14	68	0.060	12	4.98	6.194
<b>Fuse</b>	Chem. As. J.	Analogues	RCM	31	14	65	0.043	8.64	5.00	6.138
<b>Fuwa</b>	ACIE	Tot syn	RCM	15	12	85	0.023	7.5	3.03	6.269
<b>Fuwa</b>	Org. Lett.	Tot syn	Lactone	27	16	94	0.008	17.2	0.49	5.608
<b>Fuwa</b>	Tetrahedron	Tot syn	Lactone	28	12	73			1.00	5.590
<b>Gentilucci</b>	J. Med. Chem.	Med Chem	Lactam	14	13	69	0.310	70	4.43	6.163
<b>Gowrisankar</b>	Chem. Eur. J.	Tot syn	RCM	2	20	70	0.023	5	4.60	6.198
<b>Gu</b>	Org. Lett.	Tot syn	Lactam	20	13	58	0.810	300	2.70	5.722
<b>Gung</b>	Chem. Eur. J.	Methodology	Alkyne-aldehyde	14a	13	81	1.500	117.7	12.74	6.831
<b>Han</b>	JACS	Tot syn	Lactone		30	31	0.067	146.35	0.46	4.135
<b>Haug</b>	OBC	Tot syn	Lactone	1	12	65	0.027	51	0.53	5.163
<b>Hickmann</b>	JACS	Tot syn	RCAM	37	14	90	0.011	2	5.70	6.619
<b>Hili</b>	JACS	Methodology	Ugi-like	15	12	84			200	8.074
<b>Hili</b>	JACS	Methodology	Ugi-like	16	18	84	0.200	1	200	8.074
<b>Hili</b>	JACS	Methodology	Ugi-like	17	15	82			200	8.042
<b>Hili</b>	JACS	Methodology	Ugi-like	18	18	84			200	8.074
<b>Hili</b>	JACS	Methodology	Ugi-like	19	18	77			200	7.961
<b>Hili</b>	JACS	Methodology	Ugi-like	20	18	81			200	8.026
<b>Hili</b>	JACS	Methodology	Ugi-like	21	18	88			200	8.134
<b>Hili</b>	JACS	Methodology	Ugi-like	22	18	76			200	7.943
<b>Hoye</b>	ACIE	Tot syn	Lactone	29	16	54	0.032	42.7	0.76	5.077
<b>Hu</b>	Org. Lett.	Tot syn	Lactam	16	17	48	0.300	150	2.00	5.345
<b>Jogireddy</b>	ChemMedChem	Med Chem	Mitsunobu	27	14	74	0.461	45	10.24	6.618
<b>Jogireddy</b>	ChemMedChem	Med Chem	Mitsunobu	29	14	70	0.498	47	10.60	6.560
<b>Jones</b>	Org. Lett.	Tot syn	Julia-Kocienski	19	15	15	0.325	30	10.83	4.563
<b>Lassen</b>	Org. Lett.	Tot syn	Lactam	7	21	25	0.036	40	0.90	4.148
<b>Lassen</b>	Tet. Lett.	Tot syn	Lactam	4	21	48	0.037	37	1.00	5.044
<b>LeClair</b>	Tet. Lett.	Tot syn	NHK	19	14	35	0.009	18.9	0.50	4.329
<b>LeClair</b>	Tet. Lett.	Tot syn	NHK	19	14	61	0.230	22.8	10.09	6.360
<b>Li</b>	J. Org. Chem.	Tot syn	Lactone	87	22	51	0.002	0.3	5.90	5.894
<b>Li</b>	J. Org. Chem.	Tot syn	Heck	87	22	70	0.071	18	3.96	6.133
<b>Li</b>	Chem. Comm.	Tot syn	RCM	1	20	51	0.110	120	0.92	5.085
<b>Liang</b>	Chem. Comm.	Tot syn	Lactam	2	31	46			1.00	4.988
<b>Liu</b>	Chem. Comm.	Tot syn	Lactam	21	31	72			1.00	5.572
<b>Madsen</b>	Tetrahedron	Methodology	Sulfite synth	21	17	51	0.260	22	11.82	6.195
<b>Madsen</b>	Tetrahedron	Methodology	Sulfite synth	23	17	43	0.260	22	11.82	5.973
<b>Madsen</b>	Tetrahedron	Methodology	Sulfite synth	25	15	79	0.260	22	11.82	6.765
<b>Madsen</b>	Tetrahedron	Methodology	Sulfite synth	27	17	74	0.260	22	11.82	6.680
<b>Madsen</b>	Tetrahedron	Methodology	Acylation	29	19	37	0.130	10.74	12.10	5.788
<b>Madsen</b>	Tetrahedron	Methodology	Acylation	31	19	34	0.130	10.74	12.10	5.677
<b>Madsen</b>	Tetrahedron	Methodology	Red Amin	34	15	9	0.100	7.9	12.66	3.965
<b>Madsen</b>	Tetrahedron	Methodology	Red Amin	36	15	46	0.100	7.9	12.66	6.091
<b>Madsen</b>	Tetrahedron	Methodology	Red Amin	40	15	60	0.100	7.9	12.66	6.437
<b>Magauer</b>	Chem. Eur. J.	Tot syn	RCM	56	18	62	0.131	145	0.90	5.333
<b>Magauer</b>	Chem. Eur. J.	Tot syn	RCM	60	16	83	0.098	113	0.87	5.695
<b>Marcaurrelle</b>	JACS	Method/Med Chem	Huisgen	6a	12	78			10.00	6.676

<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	6b	12	75			10.00	6.625
<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	6c	12	59			10.00	6.313
<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	6d	12	56			10.00	6.245
<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	7a	13	57			10.00	6.268
<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	7b	13	58			10.00	6.290
<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	7c	13	76			10.00	6.642
<b>Marcaurelle</b>	JACS	Method/Med Chem	Huisgen	7d	13	78			10.00	6.676
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8a	14	70			10.00	6.535
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8b	14	58			10.00	6.290
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8c	14	50			10.00	6.097
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8d	14	45			10.00	5.960
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8e	14	80			10.00	6.709
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8f	14	93			10.00	6.905
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8g	14	81			10.00	6.725
<b>Marcaurelle</b>	JACS	Method/Med Chem	RCM	8h	14	88			10.00	6.833
<b>Miyatake</b>	Org. Lett.	Tot syn	Ketene	10	18	55	0.380	50	7.60	6.102
<b>Miyatake</b>	Tetrahedron	Tot syn	Ketene	2	18	46	0.150	16	9.38	5.960
<b>Mohapatra</b>	EJOC	Tot syn	NHK	24	13	81	0.540	38	14.21	6.878
<b>Mohapatra</b>	EJOC	Tot syn	NHK	31	13	85	0.640	35	18.29	7.050
<b>Monfette</b>	Chem. Eur. J.	Process	RCM	2	16	82			5.00	6.440
<b>Monfette</b>	Chem. Eur. J.	Process	RCM	2	16	70			5.00	6.234
<b>Monfette</b>	Chem. Eur. J.	Process	RCM	2	16	57			10.00	6.268
<b>Monfette</b>	Chem. Eur. J.	Process	RCM	2	16	43			20.00	6.201
<b>Monfette</b>	Chem. Eur. J.	Process	RCM	2	16	99			5.00	6.686
<b>Monfette</b>	Chem. Eur. J.	Process	RCM	2	16	65			20.00	6.740
<b>Moulin</b>	Tetrahedron	Analogues	RCM	5	24	68	0.007	20	0.35	5.042
<b>Nahrwold</b>	Org. Lett.	Analogues	Lactam	17	17	74	0.200	30	6.67	6.432
<b>Nahrwold</b>	Org. Lett.	Analogues	Huisgen	17	17	32	0.083	40	2.08	4.832
<b>Nair</b>	Tet. Lett.	Med Chem	RCM	17	16	80	0.600	60	10.00	6.709
<b>Napolitano</b>	J. Org. Chem.	Analogues	HWE	15	14	62	0.081	5	16.20	6.587
<b>Navickas</b>	Tetrahedron	Tot syn	Mitsunobu	34a	14	77	0.095	9.5	10.00	6.659
<b>Newhouse</b>	JACS	Tot syn	Lactam	124	15	41	0.012	1.1	10.91	5.876
<b>Newhouse</b>	JACS	Tot syn	Larock Indole	124	15	67	0.039	1.5	26.00	6.893
<b>Nicolaou</b>	JACS	Tot syn	Quinone DA	38a	13	60			5.00	6.033
<b>Nicolaou</b>	JACS	Tot syn	Quinone DA	76	13	40	0.055	7	7.79	5.697
<b>Nshimyumukiza</b>	Tetrahedron	Med Chem	RCM	1a	15	19	0.721	140	5.15	4.548
<b>Nshimyumukiza</b>	Tetrahedron	Med Chem	RCM	1'a	17	49	0.721	140	5.15	5.782
<b>Okura</b>	ACIE	Tot syn/analogues	Lactam	21	15	71	1.430	286	5.00	6.253
<b>Oskarsson</b>	JACS	Analogues	RCM	14	14	75	0.510	1000	0.51	5.333
<b>Ozbek</b>	EJOC	Peptidomimetics	Lactam	14	24	18			4.00	4.368
<b>Park</b>	Org. Lett.	Methodology	Dienyne RCM	entry 6	17	87			4.00	6.421
<b>Paterson</b>	Tetrahedron	Tot syn	Lactone	32	22	87	0.037	52.5	0.70	5.664
<b>Ranyuk</b>	Adv Synth Catal	Methodology	Buch-Hart	3a	21	31	0.250	12	20.83	5.793
<b>Ranyuk</b>	Adv Synth Catal	Methodology	Buch-Hart	3b	21	27	0.250	12	20.83	5.613
<b>Ranyuk</b>	Adv Synth Catal	Methodology	Buch-Hart	5a	20	30	0.250	12	20.83	5.750
<b>Ranyuk</b>	Adv Synth Catal	Methodology	Buch-Hart	5b	21	29	0.250	12	20.83	5.706
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	6a	24	30			2.00	4.732

<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	6b	24	35			2.00	4.933
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	7a	24	22			2.00	4.328
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	8	24	15			0.53	3.253
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	13a	24	58			2.00	5.591
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	13b	24	22			2.00	4.328
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	11a	24	33			2.00	4.857
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	11b	24	68			2.00	5.799
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	15a	24	42			2.00	5.171
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	15b	24	42			2.00	5.171
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	17	24	31			2.00	4.775
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	23a	24	35			2.00	4.933
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	23b	24	30			2.00	4.732
<b>Rzuczek</b>	J. Med. Chem.	Med Chem	Lactam	23c	24	40			2.00	5.107
<b>Schackel</b>	ACIE	Tot syn	Lactone	23	26	26	0.052	6	8.66	5.183
<b>Schmauder</b>	Chem. Eur. J.	Tot syn	Lactam	2a	18	29	0.037	37	1.00	4.387
<b>Shen</b>	EJOC	Methodology	Lactone	13	16	64			100.00	7.419
<b>Shimamura</b>	JACS	Tot syn	SNAr	30	16	95	0.005	5	1.00	5.933
<b>Shimamura</b>	JACS	Tot syn	Larock Indole	37	16	56	0.004	4.4	1.00	5.245
<b>Singh</b>	J. Org. Chem.	Tot syn	Lactone	16	16	77	0.132	229	0.58	5.420
<b>Sinigaglia</b>	Chem. Eur. J.	Tot syn	Enamine attack	15	17	33	0.140	20	7.00	5.401
<b>Smith</b>	Org. Lett.	Analogues	Lactone	9	23	62	0.028	57.68	0.49	5.063
<b>Smith</b>	Org. Lett.	Analogues	Lactone	10	29	48	0.010	20.49	0.49	4.732
<b>Smith</b>	Tetrahedron	Tot syn	RCM	28	17	80	0.350	1000	0.35	5.253
<b>Smith</b>	Tetrahedron	Tot syn	RCM	32/33	17	80	0.039	250	0.16	4.902
<b>Smith</b>	ACIE	Tot syn	RCAM	3	17	63	0.008	6.5	1.23	5.488
<b>Socha</b>	BMC	Med Chem	Lactam	3	16	23			0.10	3.085
<b>Speicher</b>	EJOC	Tot syn	McMurray	30	16	27	1.380	250	5.52	5.036
<b>Speicher</b>	EJOC	Tot syn	Wittig	40	18	44	3.460	500	6.92	5.770
<b>Speicher</b>	EJOC	Tot syn	Wittig	65	20	12	8.030	700	11.47	4.297
<b>Tadross</b>	Org. Lett.	Tot syn	RCM	14	10	77			2.00	5.961
<b>Tanaka</b>	Org. Lett.	Tot syn	Dotz annulation	2	18	58	0.023	11.3	2.00	5.591
<b>Tang</b>	Tetrahedron	Tot syn	NHK	37	13	63	0.620	453	1.37	5.534
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28a	18	44			1.00	4.930
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28b	18	47			1.00	5.016
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	29b	18	33			1.00	4.556
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28c	18	74			1.00	5.608
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28e	18	58			1.00	5.290
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28f	18	48			1.00	5.044
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28g	18	71			1.00	5.554
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	28h	18	43			1.00	4.900
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	30a	19	42			1.00	4.870
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	30b	19	41			1.00	4.838
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	30d	19	84			1.00	5.773
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	31e	19	85			1.00	5.788
<b>Tannert</b>	JACS	Tot syn/analogues	RCM	31f	19	76			1.00	5.642
<b>Terracciano</b>	Tetrahedron	Med Chem	RCM	1	13	15			0.30	3.005
<b>Terracciano</b>	Tetrahedron	Med Chem	RCM	2	17	32			0.30	3.993

<b>Terracciano</b>	Tetrahedron	Med Chem	RCM	3	15	21		0.30	3.444	
<b>Terracciano</b>	Tetrahedron	Med Chem	RCM	6	17	22		0.30	3.504	
<b>Terracciano</b>	Tetrahedron	Med Chem	RCM	7	15	21		0.30	3.444	
<b>Trost</b>	JACS	Tot syn	Pd Diyne coupling	30	22	56	0.003	2.6	1.31	5.361
<b>Trost</b>	JACS	Tot syn	Pd Diyne coupling	40	22	65	0.020	5.1	3.92	6.032
<b>Trost</b>	Chem. Eur. J.	Tot syn	RCM	36	15	44			1.00	4.930
<b>Wang</b>	Tetrahedron	Tot syn	Lactam	2	27	60	0.187	30	6.23	6.129
<b>Wang</b>	ACIE	Tot syn	Suzuki	15	16	66	0.070	85.3	0.82	5.373
<b>Wang</b>	ACIE	Tot syn	SNAr	22	16	62			0.10	4.377
<b>Woo</b>	JACS	Tot syn	Prins	16	14	69	0.159	8	19.88	6.815
<b>Wu</b>	ACIE	Tot syn	Barbier	21	23	40	0.050	25	2.00	5.107
<b>Wu</b>	JACS	Tot syn	Lactam	1	33	54	0.010	8.3	1.20	5.278
<b>Wullschleger</b>	Org. Lett.	Tot syn	RCM	24	14	65	0.087	101	0.86	5.374
<b>Xie</b>	J. Org. Chem.	Tot syn	Lactone	23	18	88	0.090	119.24	0.75	5.711
<b>Yadav</b>	Tetrahedron	Tot syn	RCM	14	10	64	0.462	710	0.65	5.232
<b>Yoshida</b>	Org. Lett.	Tot syn	Lactone	16a	19	65			3.00	5.916
<b>Yu</b>	Org. Lett.	Tot syn	Lactam	25	18	67	0.024	44	0.55	5.215
<b>Yu</b>	Chem. Eur. J.	Methodology	Alkylation	2	12	38	1.000	200	5.00	5.438
<b>Yu</b>	Chem. Eur. J.	Methodology	Alkylation	3	13	78	1.000	200	5.00	6.375
<b>Yu</b>	Chem. Eur. J.	Methodology	Alkylation	4	14	45	1.000	200	5.00	5.659
<b>Yu</b>	Chem. Eur. J.	Methodology	Alkylation	5	15	67	1.000	200	5.00	6.177
<b>Yu</b>	Chem. Eur. J.	Methodology	Alkylation	6	13	61	1.000	200	5.00	6.055
<b>Yu</b>	Chem. Eur. J.	Methodology	Alkylation	9	13	70	1.000	200	5.00	6.234
<b>Yun</b>	ACIE	Tot syn	RCM	21	18	45	0.030	25	1.20	5.039
<b>Zurwerra</b>	Org. Lett.	Tot syn	HWE	20	18	85			3.00	6.265

## 2011

Name	Journal	Paper type	Reaction	Cmpd no	Ring size	Yield	mmol	Vol (ml)	Conc (mM)	Emac
<b>Andersson</b>	J. Med. Chem.	Med chem	RCM	19	14	28	0.019	5.5	3.38	4.871
<b>Andersson</b>	J. Med. Chem.	Med chem	RCM	23	14	59	0.088	15	5.85	6.080
<b>Andersson</b>	J. Med. Chem.	Med chem	RCM	25	14	43	0.061	12	5.05	5.604
<b>Ando</b>	Tet. Lett.	Methodology	HWE	13	15	63	0.200	20	10.00	6.398
<b>Ando</b>	Tet. Lett.	Methodology	HWE	14	16	66	0.200	20	10.00	6.459
<b>Ando</b>	Tet. Lett.	Methodology	HWE	15	13	70	0.200	20	10.00	6.535
<b>Ando</b>	Tet. Lett.	Methodology	HWE	16	14	69	0.200	20	10.00	6.517
<b>Ando</b>	Tet. Lett.	Methodology	HWE	17	17	75	0.200	20	10.00	6.625
<b>Ando</b>	Tet. Lett.	Methodology	HWE	18	18	69	0.200	20	10.00	6.517
<b>Araoz</b>	JACS	Tot syn	RCM	41	22(27?)	75	0.538	108	4.98	6.323
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	3	16	73	0.120	5	24.00	6.970
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	8	14	62			24.00	6.757
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	9	18	74			24.00	6.988
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	10	21	81			24.00	7.106
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	11	23	78			24.00	7.056
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	12	28	98			24.00	7.354
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	13	18	67			24.00	6.858
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	14	27	69			24.00	6.897

<b>Bedard</b>	JACS	Methodology	Glaser-Hay	16	16	98			24.00	7.354
<b>Bedard</b>	JACS	Methodology	Glaser-Hay	15	23	63			24.00	6.778
<b>Benelkebir</b>	BMC	Tot syn	Lactam	31	16	64	0.230	185	1.24	5.513
<b>Berwe</b>	OPRD	NP Process	Lactam	2	15	73			57.00	7.346
<b>Bhansali</b>	J. Med. Chem.	Med chem	Lactam	25		57			1.00	5.268
<b>Bienek</b>	Organomet	Methodology	RCM		15	95			20.00	7.234
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	1	14	80	0.010	0.6	16.67	6.931
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	2	12	42			16.67	6.092
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	3	17	75			16.67	6.847
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	4	15	65			16.67	6.661
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	5	19	21			16.67	5.189
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	6	22	83			16.67	6.979
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	7	22	75			16.67	6.847
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	8	23	80			16.67	6.931
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	9	26	25			16.67	5.416
<b>Bogdan</b>	Org. Lett.	Methodology	Huisgen	10	31	22			16.67	5.249
<b>Bogdan</b>	JACS	Structural	Huisgen	1	14	90			16.67	7.085
<b>Bogdan</b>	JACS	Structural	Huisgen	2	14	80			16.67	6.931
<b>Bogdan</b>	JACS	Structural	Huisgen	3	13	66			16.67	6.681
<b>Bogdan</b>	JACS	Structural	Huisgen	4	12	73			16.67	6.812
<b>Bogdan</b>	JACS	Structural	Huisgen	5	11	20			16.67	5.125
<b>Bogdan</b>	JACS	Structural	Huisgen	7	12	40			16.67	6.028
<b>Bogdan</b>	JACS	Structural	Huisgen	8	12	54			16.67	6.419
<b>Bogdan</b>	JACS	Structural	Huisgen	9	13	35			16.67	5.854
<b>Bogdan</b>	OBC	PK study	Huisgen	1a	12	87	0.010	0.6	16.67	7.040
<b>Bogdan</b>	OBC	PK study	Huisgen	2a	14	80	0.010	0.6	16.67	6.931
<b>Bogdan</b>	OBC	PK study	Huisgen	3a	18	57	0.010	0.6	16.67	6.489
<b>Bogdan</b>	OBC	PK study	Huisgen	4a	22	71	0.010	0.6	16.67	6.776
<b>Bogdan</b>	OBC	PK study	Huisgen	5a	21	63	0.010	0.6	16.67	6.620
<b>Bogdan</b>	OBC	PK study	Huisgen	6a	28	62	0.010	0.6	16.67	6.599
<b>Bogdan</b>	OBC	PK study	Huisgen	7a	25	44	0.010	0.6	16.67	6.152
<b>Bogdan</b>	OBC	PK study	Huisgen	8a	29	59	0.010	0.6	16.67	6.534
<b>Brady</b>	OBC	Method	RCM	37	10	71	0.210	53	3.96	6.152
<b>Brady</b>	OBC	Method	RCM	30	10	71	0.160	53	3.02	6.034
<b>Brandt</b>	ACS MCL	Med Chem	RCM	7	21	83	0.013	13	1.00	5.757
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8a	13	65	0.370	4.3	86.05	7.373
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8b	13	43	0.370	4.3	86.05	6.835
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8c	13	62	0.370	4.3	86.05	7.312
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8d	13	75	0.370	4.3	86.05	7.560
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8e	13	51	0.370	4.3	86.05	7.057
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8f	13	53	0.370	4.3	86.05	7.108
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8g	13	94	0.370	4.3	86.05	7.854
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8h	13	54	0.370	4.3	86.05	7.132
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8i	13	32	0.260	4.25	61.18	6.302
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8j	13	37	0.130	2.12	61.32	6.492
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8k	13	51	0.370	4.3	86.05	7.057
<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8l	13	63	0.094	3	31.33	6.894

<b>Breslin</b>	J. Med. Chem.	Med Chem	Heck	8m	13	72	0.140	3.3	42.42	7.200
<b>Bruno</b>	Org. Lett.	Tot syn	Lactam		36	24	0.096	40	2.41	4.522
<b>Carr</b>	OBC	Tot syn	Germyl-Stille	5	11	9	0.004	3	1.33	2.988
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	3a	14	55	10.000	100	100.00	7.221
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	3b	15	63	10.000	100	100.00	7.398
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	7a	15	60	10.000	100	100.00	7.334
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	7b	15	64	10.000	100	100.00	7.419
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	11	20	46	10.000	100	100.00	6.988
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	14a	14	58	10.000	100	100.00	7.290
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	14b	14	62	10.000	100	100.00	7.377
<b>Chande</b>	J Het Chem	Methodology	Dieckmann	14c	14	50	10.000	100	100.00	7.097
<b>Chavez</b>	J. Med. Chem.	Med chem	RCM	8	20	52	0.110	21	5.24	5.867
<b>Chegondi</b>	J. Org. Chem.	Tot syn	RCM	3	12	82	0.026	5	5.20	6.457
<b>Chen</b>	ACIE	Tot syn	Lactam		20	74	0.247	350	0.71	5.456
<b>Chen</b>	J. Org. Chem.	Tot syn	Stille	34	24	72	0.020	28	0.70	5.417
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	2	12	86			20.00	7.105
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	3	14	95			20.00	7.234
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	4	12	45			20.00	6.261
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	5	14	76			20.00	6.943
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	6	15	70			20.00	6.836
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	7	15	68			20.00	6.799
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	8	17	66			20.00	6.760
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	9	19	66			20.00	6.760
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	10	21	51			20.00	6.424
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	11	16	45			20.00	6.261
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	12	16	67			20.00	6.779
<b>Chouhan</b>	Org. Lett.	Methodology	Huisgen	13	24	85			20.00	7.089
<b>Coghlan</b>	BMC	Med Chem	RCM	16	18	77	0.250	50	5.00	6.358
<b>Coghlan</b>	BMC	Med Chem	RCM	17	18	74	0.200	50	4.00	6.210
<b>Coghlan</b>	BMC	Med Chem	RCM	18	18	72	0.500	120	4.17	6.192
<b>Coghlan</b>	BMC	Med Chem	RCM	19	18	77	0.170	50	3.40	6.191
<b>Coghlan</b>	BMC	Med Chem	RCM	20	18	72			5.00	6.271
<b>Cortes</b>	EJOC	Tot syn	SNAr	27	16	86	0.270	27	10.00	6.803
<b>Crane</b>	ACIE	Tot syn	Prins	20	16	66	0.250	25	10.00	6.459
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	13	13	79			10.00	6.693
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	14	14	85			10.00	6.788
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	15	15	73			10.00	6.590
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	16	16	88			10.00	6.833
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	17	17	83			10.00	6.757
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	18	18	69			10.00	6.517
<b>Dandapani</b>	J. Org. Chem.	Med Chem	RCM	19	15	87			10.00	6.819
<b>Day</b>	ACS Chem. Biol.	Med chem	RCM	12	14	69	0.280	70	4.00	6.119
<b>Day</b>	ACS Chem. Biol.	Med chem	RCM	16	14	87			4.00	6.421
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	42	12	56	0.550	35	15.71	6.441
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	28	12	53	0.560	36	15.56	6.365
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	29	12	92	0.370	25	14.80	7.062
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	30	12	38	1.400	90	15.56	5.931

<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	31	12	19	0.190	12	15.83	5.036
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	32	12	38	0.240	15	16.00	5.943
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	33	12	37	0.400	26	15.38	5.892
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	34	12	24	0.850	54	15.74	5.338
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	35	12	37	0.760	40	19.00	5.983
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	36	12	46	0.490	31	15.81	6.187
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	37	12	35	0.730	47	15.53	5.823
<b>Demillo</b>	J. Med. Chem.	Med chem	Pd-AA	38	12	26	0.290	15	19.33	5.531
<b>Dermenci</b>	Chem. Sci.	Method/analogues	NHK	5	14	59	0.711	720	0.99	5.307
<b>Doi</b>	Chem. As. J.	Tot syn	Lactam	26	25	60	0.033	25	1.32	5.455
<b>ElMarrouni</b>	Org. Lett.	Tot syn	Lactam/ketene	20	12(14?)	75	0.110	1004	0.11	4.665
<b>Fischer</b>	Chem. Eur. J.	Tot syn	Lactone	1f	16	53	0.013	20	0.66	4.992
<b>Franklin</b>	OBC	Analogues	Nitrile oxide	22	15	36	0.082	40	2.05	4.981
<b>Gagnepain</b>	Chem. Eur. J.	Tot syn	RCM	59	24	72	0.359	359	1.00	5.572
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	7a	15	91			2.00	6.178
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	7b	15	70			2.00	5.836
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	10	12	53			2.00	5.474
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	11	13	80			2.00	6.010
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	13	14	92			2.00	6.192
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	15	14	90			2.00	6.164
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	17	15	73			2.00	5.891
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	19	15	75			2.00	5.926
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	21	17	92			2.00	6.192
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	22	18	90			2.00	6.164
<b>Gallenkamp</b>	JACS	Method/Tot syn	RCM	38	12	76	0.086	86	1.00	5.642
<b>Gentilucci</b>	ChemMedChem	Turn mimic	Lactam	11	15	70	0.150	40	3.75	6.109
<b>Gentilucci</b>	ChemMedChem	Turn mimic	Lactam	12	15	55	0.150	40	3.75	5.795
<b>Gentilucci</b>	ChemMedChem	Turn mimic	Lactam	13	15	60	0.150	40	3.75	5.908
<b>Gentilucci</b>	ChemMedChem	Turn mimic	Lactam	14	15	65	0.150	40	3.75	6.013
<b>Gentilucci</b>	ChemMedChem	Turn mimic	Lactam	15	14	65	0.150	40	3.75	6.013
<b>Gersbach</b>	Chem. Eur. J.	Tot syn	RCM	3	12	80			3.00	6.186
<b>Gesinski</b>	JACS	Tot syn	Sakurai	10	16(20?)	76	0.368	7.4	49.73	7.339
<b>Gladding</b>	Tetrahedron	Tot syn	Oxidation/ketal	59	14	66	0.017	8.3	1.99	5.757
<b>Hanson</b>	J. Org. Chem.	Tot syn	RCM	1	24	57			0.50	4.967
<b>He</b>	Chem. As. J.	Tot syn	Lactam	25	15	75	0.110	505	0.22	4.963
<b>Hickmann</b>	JACS	Tot syn	RCAM	42	13	79	0.026	13	0.50	5.392
<b>Hickmann</b>	JACS	Tot syn	RCAM	47	14	94	0.015	1	14.50	7.081
<b>Hickmann</b>	JACS	Tot syn	RCAM	52	14	90	0.011	1	11.40	6.920
<b>Hili</b>	Tetrahedron	Tot syn	Mitsunobu	18	12	74	0.574	11.7	49.06	7.298
<b>Hopkins</b>	Org. Lett.	Tot syn	Lactone	1	30	55	0.088	100	1.50	5.397
<b>Hoveyda</b>	J. Med. Chem.	Med chem	Lactam	2	18	83	78.400	1570	49.94	7.456
<b>Hunter</b>	J. Org. Chem.	Tot syn	Lactam	1	23	81			5.00	6.424
<b>Ibrahim-Ouali</b>	Tet. Lett.	Analogues	RCM	8	17	65			0.008	3.342
<b>Ibrahim-Ouali</b>	Tet. Lett.	Analogues	RCM	16	19	90			0.008	3.766
<b>Inokuchi</b>	ACS MCL	Med chem	Nitrile-oxide	13b	15	33	0.082	40	6.00	5.334
<b>Iwashita</b>	Tetrahedron	Tot syn/analogues	SNAr	24	18	60			10.00	6.334
<b>Jacobsen</b>	J. Org. Chem.	Helix mimic	Huisgen	5	19	83			0.15	4.933

<b>Jakubec</b>	Tet. Lett.	Tot syn	RCAM	18	15	36	0.020	85.6	0.23	4.037
<b>Kang</b>	BMCL	Med Chem	RCM	14	14	46			10.00	5.988
<b>Kim</b>	BMC	Med Chem	Alkylation	22	13	64	0.540	54	10.00	6.419
<b>Kim</b>	BMC	Med Chem	RCM	27	12	23	0.920	184	5.00	4.784
<b>Knapp</b>	Org. Lett.	Methodology	Ketene/lactam	29a	12	94			12.00	6.999
<b>Knapp</b>	Org. Lett.	Methodology	Ketene/lactam	29b	12	92			12.00	6.971
<b>Knapp</b>	Org. Lett.	Methodology	Nitrile oxide	30	12	74			12.00	6.687
<b>Knapp</b>	Org. Lett.	Methodology	Nitrile oxide	32	12	70			12.00	6.614
<b>Knapp</b>	Org. Lett.	Methodology	Nitrile oxide	34	14	72			12.00	6.651
<b>Knapp</b>	Org. Lett.	Methodology	Nitrile oxide	35	14	70			12.00	6.614
<b>Knowles</b>	Chem. Sci.	Tot syn	Aldol	10	13	67	2.140	425	5.04	6.180
<b>Knowles</b>	Chem. Sci.	Tot syn	Suzuki	16	12	50	0.017	3.78	4.37	5.737
<b>Kong</b>	JACS	Tot syn	Barbier	65	15(17?)	56	0.101	15	6.73	6.073
<b>Kostiuk</b>	Chem. Eur. J.	Tot syn	Wittig	43	18	69	0.200	2	100.00	7.517
<b>Kwon</b>	Tetrahedron	Tot syn	RCM	21	13	46	0.044	8.8	4.98	5.685
<b>Larivee</b>	ACIE	Tot syn	Lactone	49	18	68	0.011	23.936	0.45	5.152
<b>Li</b>	J. Nat. Prod.	Tot syn	Lactam	1+2	16	48	0.167	200	0.84	4.965
<b>Lin</b>	Org. Lett.	Tot syn	Lactone	5	19	81	0.230	313	0.73	5.592
<b>Linder</b>	JACS	Tot syn	Lactam	20	24	51	0.040	50	0.80	5.026
<b>Liniger</b>	ACS MCL	Med Chem	Mitsunobu	16	14	49	0.531	306	1.74	5.310
<b>Liniger</b>	ACS MCL	Med Chem	Mitsunobu	26	14	64	1.080	154	7.01	6.264
<b>Liu, J</b>	JACS	Tot syn	Suzuki	13	14	48	0.057	5	11.40	6.101
<b>Liu, K</b>	JACS	Tot syn	Lactone	12	14	64	0.540	170	3.18	5.920
<b>Martinez</b>	J. Org. Chem.	Tot syn	RCM	33	12(15?)	97	0.690	75.8	9.10	6.919
<b>Muranaka</b>	J. Org. Chem.	Method/Tot syn	RCM	43a	13	98			100.00	7.974
<b>Muranaka</b>	J. Org. Chem.	Method/Tot syn	RCM	43b	13	98			100.00	7.974
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10a	16	60	1.000	15	66.67	7.158
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10b	15	63	1.000	15	66.67	7.222
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10c	16	67	1.000	15	66.67	7.302
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10d	17	70	1.000	15	66.67	7.359
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10e	19	59	1.000	15	66.67	7.136
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10f	17	71	1.000	15	66.67	7.378
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10g	17	67	1.000	15	66.67	7.302
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10h	17	68	1.000	15	66.67	7.321
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10i	16	62	1.000	15	66.67	7.201
<b>Muthusamy</b>	Tet. Lett.	Methodology	Carbonyl ylide	10j	18	69	1.000	15	66.67	7.340
<b>Nagasawa</b>	Tetrahedron	Tot syn	Lactone	32	12	30	0.023	28.5	0.81	4.338
<b>Nicolaou</b>	ACIE	Tot syn	RCM	9	13	90	0.140	40	3.50	6.407
<b>Oelke</b>	Chem. Eur. J.	Tot syn	Lactam	71	18	79	0.010	3	3.33	6.216
<b>Park</b>	OBC	Methodology	Dienyne RCM	4j	14	71	0.100	25	4.00	6.156
<b>Paterson</b>	Org. Lett.	Tot syn	Lactone	21a	16	95	0.009	8.4	1.04	5.948
<b>Paterson</b>	Tetrahedron	Tot syn	Lactone	53	18	69	0.060	60	1.00	5.518
<b>Paterson</b>	Tetrahedron	Tot syn	Stille	53	18	64	0.024	14	1.72	5.654
<b>Roberts</b>	J. Nat. Prod.	Tot syn	Suzuki	11	14	42	0.140	7	20.00	6.171
<b>Rodenko</b>	ACS MCL	Med chem	Alkylation	9	48	43	0.500	300	1.67	5.122
<b>Rossle</b>	Org. Lett.	Tot syn	RCM		21	43	0.096	60	1.60	5.105
<b>Rudd</b>	ACS MCL	Med chem	RCM	10j	18	92	100.000	14300	6.99	6.736

<b>Saitman</b>	Org. Lett.	Tot syn	NHK	6	13	82	2.197	449	4.89	6.431
<b>Scheiss</b>	Org. Lett.	Tot syn	RCM	13	16	77	0.360	251	1.43	5.816
<b>Schneider</b>	ACIE	Tot syn	Castro-Stephens	2	12	67	0.062	14	4.43	6.124
<b>Shareef</b>	Chem. Sci.	Methodology	Alkyne/aldehyde	14	12	58	0.085	10	8.50	6.220
<b>Shareef</b>	Chem. Sci.	Methodology	Alkyne/aldehyde	15	11	59	0.103	10	10.30	6.325
<b>Sharpe</b>	J. Org. Chem.	Tot syn	Lactone	20	16	45	0.233	61.8	3.77	5.536
<b>Shen</b>	Tet. Lett.	Tot syn	Ullmann	11	14	52			20.00	6.449
<b>Shibahara</b>	Org. Lett.	Tot syn	Lactam	32	18	25	0.037	5.3	7.04	5.041
<b>Smith</b>	JACS	Tot syn	Lactone	37	23	62	0.028	52.98	0.53	5.100
<b>Smith</b>	Tetrahedron	Tot syn	Lactone	142	31	80	0.002	2.1	0.95	5.688
<b>Smith</b>	Tetrahedron	Tot syn	HWE	51a	21	44	0.020	34	0.59	4.700
<b>Song</b>	J. Org. Chem.	Process	Lactam	8	20	66	98.000	5800	16.90	6.686
<b>Sugiyama</b>	Tetrahedron	Analogues	RCM	23	14	97	0.119	24	4.96	6.656
<b>Sund</b>	BMCL	Med Chem	Lactam	41	15	26			1.00	4.245
<b>Sund</b>	BMCL	Med Chem	Lactam	42	15	28			1.00	4.341
<b>Sund</b>	BMCL	Med Chem	Lactam	43	15	25			1.00	4.194
<b>Sund</b>	BMCL	Med Chem	Lactam	44	15	43			1.00	4.900
<b>Sund</b>	BMCL	Med Chem	Lactam	45	15	7			1.00	2.535
<b>Sund</b>	BMCL	Med Chem	Lactam	46	15	44			1.00	4.930
<b>Tang</b>	BMCL	Med Chem	RCM	10	12	74	0.375	40	9.38	6.580
<b>Tenebaum</b>	ACIE	Tot syn	RCM	25	11	65	0.110	1605	0.07	4.275
<b>Trost</b>	Chem. Eur. J.	Tot syn	RCM	66	25	80	0.018	10	1.80	5.965
<b>Uchiro</b>	Org. Lett.	Tot syn	Ullmann	7	13	42	0.017	3.3	5.06	5.574
<b>Van der Berge</b>	EJOC	Structural	RCM	1a	15	19	0.279	80	3.49	4.379
<b>Van der Berge</b>	EJOC	Structural	RCM	1b	16	41	0.279	80	3.49	5.381
<b>Van der Berge</b>	EJOC	Structural	RCM	1c	17	49	0.279	80	3.49	5.613
<b>Van der Berge</b>	EJOC	Structural	RCM	1d	18	58	0.279	80	3.49	5.833
<b>Velvadapu</b>	J. Org. Chem.	Tot syn	NHK	48	14	50	0.270	158	1.71	5.330
<b>Wang, Y</b>	JACS	Methodology	RCM	6	10	87			2.00	6.120
<b>Wang, Y</b>	JACS	Methodology	RCM	7	11	36			2.00	4.970
<b>Wang, Y</b>	JACS	Methodology	RCM	8	11	43			2.00	5.201
<b>Wang, Y</b>	JACS	Methodology	RCM	9	11	36			2.00	4.970
<b>Wang, Y</b>	JACS	Methodology	RCM	10	12	33			2.00	4.857
<b>Wang, Y</b>	JACS	Methodology	RCM	11	12	59			2.00	5.614
<b>Wang, Y</b>	JACS	Methodology	RCM	12	12	82			2.00	6.042
<b>Wang, Y</b>	JACS	Methodology	RCM	13	12	46			2.00	5.289
<b>Wang, Y</b>	JACS	Methodology	RCM	14	12	79			2.00	5.994
<b>Wang, Y</b>	JACS	Methodology	RCM	15	12	14			2.00	3.739
<b>Wang, Y</b>	JACS	Methodology	RCM	16	14	76			2.00	5.943
<b>Wang, Y</b>	JACS	Methodology	RCM	17	15	43			2.00	5.201
<b>Wang, Y</b>	JACS	Methodology	RCM	18	16	47			2.00	5.317
<b>Wang, Y</b>	JACS	Methodology	RCM	19	14	95			2.00	6.234
<b>Wang, Y</b>	JACS	Methodology	RCM	20	15	76			2.00	5.943
<b>Wang, Y</b>	JACS	Methodology	RCM	21	16	41			2.00	5.139
<b>Wender</b>	JACS	Tot syn	Prins	18	20	65			20.00	6.740
<b>Wender</b>	Tetrahedron	Analogues	Acetalization	5	20	83	0.020	5.5	3.64	6.318
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	16a	18	48	0.052	50	1.04	5.061

<b>Williams</b>	J. Med. Chem.	Med chem	RCM	16b	17	51	0.052	50	1.04	5.140
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	16c	18	45	0.052	50	1.04	4.977
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	16d	17	45	0.052	50	1.04	4.977
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	16e	18	48	0.052	50	1.04	5.061
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	17a	18	43	0.052	50	1.04	4.917
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	17b	18	96	0.052	50	1.04	5.964
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	17d	18	52	0.052	50	1.04	5.165
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	17h	18	56	0.052	50	1.04	5.262
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	21a	18	74	0.052	50	1.04	5.625
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	21g	18	56	0.052	50	1.04	5.262
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	21j	18	73	0.052	50	1.04	5.607
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22a	18	43	0.052	50	1.04	4.917
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22b	18	79	0.052	50	1.04	5.710
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22c	18	56	0.052	50	1.04	5.262
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22d	18	32	0.052	50	1.04	4.532
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22e	18	42	0.052	50	1.04	4.887
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22f	18	48	0.052	50	1.04	5.061
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	22g	18	49	0.052	50	1.04	5.088
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26a	18	73			2.50	5.988
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26b	18	81			2.50	6.123
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26c	17	82			2.50	6.139
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26f	17	73			2.50	5.988
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26h	18	78	2.500	1000	2.50	6.074
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26i	18	74			2.50	6.006
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26j	18	64			2.50	5.816
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26k	18	49			2.50	5.469
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26l	18	53			2.50	5.571
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26m	18	52			2.50	5.546
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26n	18	75			2.50	6.023
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26o	18	42			2.50	5.268
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26p	18	77			2.50	6.057
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	26q	18	63			2.50	5.796
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27a	18	71			2.50	5.952
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27b	18	48			2.50	5.442
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27c	18	53			2.50	5.571
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27d	18	56			2.50	5.643
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27e	18	68			2.50	5.895
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27f	18	59			2.50	5.710
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27h	18	59			2.50	5.710
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27i	18	88			2.50	6.231
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27j	18	50			2.50	5.495
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27k	18	67			2.50	5.876
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27l	18	51			2.50	5.521
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27m	18	62			2.50	5.775
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27n	18	66			2.50	5.857
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27o	18	50			2.50	5.495
<b>Williams</b>	J. Med. Chem.	Med chem	RCM	27p	18	62			2.50	5.775

Williams	J. Med. Chem.	Med chem	RCM	27q	18	44			2.50	5.328
Williams	J. Med. Chem.	Med chem	RCM	27r	18	68			2.50	5.895
Williams	J. Med. Chem.	Med chem	RCM	28a	18	56			2.50	5.643
Williams	J. Med. Chem.	Med chem	RCM	28b	18	49			2.50	5.469
Williams	J. Med. Chem.	Med chem	RCM	28c	18	54			2.50	5.595
Williams	J. Med. Chem.	Med chem	RCM	28d	18	51			2.50	5.521
Williams	J. Med. Chem.	Med chem	RCM	28e	18	44			2.50	5.328
Williams	Tetrahedron	Tot syn	Lactone	66	14	63	0.012	120	0.10	4.402
Wu	Tet. Lett.	Tot syn	Lactam	1	24	56	0.052	77	0.68	5.074
Wu, J	J. Org. Chem.	Tot syn	Barbier	9	23	42			2.00	5.171
Xie, J	JACS	Tot syn	SNAr	10	16	85			5.00	6.487
Xie, J	JACS	Tot syn	Lactam	26	12	76			2.00	5.943
Xue	Org. Lett.	Tot syn	Lactone	8	14	46	1.100	350	3.14	5.486
Yadav	Org. Lett.	Tot syn	Prins	4	14	79	0.280	15	18.67	6.964
Yamamoto	JACS	Tot syn	Lactone	16	26	93			1.60	6.110
Zhan	Chem. Eur. J.	Tot syn	RCM	36	16	84	0.013	12.5	1.00	5.773
Zhan	Chem. Eur. J.	Tot syn	Lactone	63	16	60	0.326	265	1.23	5.424
Zhang, J	Org. Lett.	Tot syn	Lactam	26	14	77	0.180	200	0.90	5.614
Zhang, J	Org. Lett.	Analog syn	Huisgen	9d	17	46			1.00	4.988
Zhang, J	Org. Lett.	Analog syn	Huisgen	17a	13	14			5.00	4.137
Zhang, J	Org. Lett.	Analog syn	Huisgen	17b	14	68			5.00	6.196
Zhang, J	Org. Lett.	Analog syn	Huisgen	17c	15	59			5.00	6.012
Zhang, J	Org. Lett.	Analog syn	Huisgen	17d	16	55			5.00	5.920
Zhang, J	Org. Lett.	Analog syn	Lactam	20a	13	47			0.50	4.715
Zhang, J	Org. Lett.	Analog syn	Lactam	20b	16	70			0.50	5.234
Zhang, Y	J. Org. Chem.	Tot syn	Lactone	49	36	75	0.030	88	0.34	5.158
Zhdanko	Chem. Eur. J.	Analogues	Lactam	2e	18	48	0.074	30	2.46	5.435
Zhu	Org. Lett.	Tot syn	Lactone	22	28	60	0.010	2	5.00	6.033

## Part 2: Full reference list

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