





















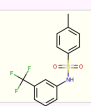




df_frag_final_supplier_oe.csv																	
Molecule	eMolecules SMILES	pKas in [3,11]	XlogP	MolWt	Availability (mg)	Price	group	N_Rot	N_UV_chrom	Selection	Bin index	Priority	Final list	Supplier Name	Catalog Number	canonical isomeric SMILES	eMolecules ID
	Oc1cc2c3CCCNc1=O)c3oc2cc1	[9.119]	0.72	217.221	184.0	533.0	fragment-like	0	27	picked	4	1	True	ChemDiv	5816-0042	c1cc2c(c1O)c3c(c2)c1=O)NCCC3	6679830
	FC(c1cccc(c1)Nc1nncn2c1ccc2)(F)F	[4.05]	3.27	289.255	101.0	355.0	fragment-like	3	36	picked	17	2	True	ChemDiv	3232-0333	c1ccc2c(c1)c(ncn2)Nc3cccc(c3)C(F)(F)F	1327907
	Clc1ccc(o1)c1=O)Nc1cccc1N1CCCCC1	[5.346]	3.79	304.771	424.3	148.0	fragment-like	4	18	picked	19	1	True	Enamine Screening Compounds	Z119335440	c1ccc(c(c1)NC(=O)c2ccc(o2)C1)N3CCCCC3	18908671
	Clc1ccc(cc1)CNc1nncn2c1ccc2	[5.564]	3.52	269.72900000000004	415.5	148.0	fragment-like	3	36	picked	18	2	True	Enamine Screening Compounds	Z126957826	c1cc2c(c1)c(ncn2)Nc3ccc(cc3)Cl	30719859
	O=C(c1ccsc1)Nc1nnc(s1)C1cccc1	[7.12]	3.61	301.387	379.0	148.0	fragment-like	5	14	picked	18	1	True	Enamine Screening Compounds	Z27474679	c1ccc(cc1)Cc2nnc(s2)NC(=O)c3ccsc3	1228629
	c1ccc(cc1)CNc1nncn2c1ccc2	[5.564]	2.91	235.284	208.7	223.0	fragment-like	3	36	picked	15	2	True	Enamine Screening Compounds	Z57161635	c1ccc(cc1)CNc2c3cccc3ncn2	1327878
	Brc1ncc(c1)C(=O)Nc1cccc2c1ncc2	[4.001, 10.328]	2.92	328.163	406.0	148.0	fragment-like	3	37	picked	15	1	True	Enamine Screening Compounds	Z28487401	c1cc2ccncc2c1)NC(=O)c3cc(cnc3)Br	18893169
	OC(=O)C1c1c=O)[nH]c2c(c1c1cccc1)c(c2)C	[4.109]	2.80	293.317	232.1	223.0	fragment-like	3	59	picked	14	2	True	Enamine Screening Compounds	Z57157353	Cc1ccc2c1)c(c1c=O)[nH]2)CC(=O)O)c3cccc3	1367649
	OC1cccc(c1)Nc1nncn2c1ccc2.Cl	[4.05]	3.14	287.744	119.7	148.0	fragment-like	3	36	picked	16	1	True	Enamine Screening Compounds	Z220564816	OC1cccc(c1)Nc2c3cccc3nncn2.Cl	1865544
	Nc1ccc2c(c1)Ncnc2c1ccc1	[6.348]	2.33	209.247	50213.0	148.0	fragment-like	1	40	picked	12	1	True	Enamine Screening Compounds	Z57290870	c1ccc(cc1)n2cnc3c2cc(c3)N	3165344
	Oc1ccc(cc1)n1nnc2c1ccc2	[5.82, 8.709]	2.22	210.231	21650.2	148.0	fragment-like	1	40	picked	11	1	True	Enamine Screening Compounds	Z1318268952	c1ccc2c(c1)nnc2c3ccc(cc3)O	37095168
	O=C(Nc1nc2c(s1)cccc2)CNc1=O)c1cccc1	[8.672]	2.72	311.358	149.1	148.0	fragment-like	6	28	picked	14	1	True	Enamine Screening Compounds	Z69130143	c1ccc(cc1)C(=O)NCC(=O)Nc2nc3cccc3s2	23354217
	O=C1Nc2c(C)/C1=C/c1ccc(cc1)C(C)C)ccc2	[3.158]	4.23	263.334	283.7	400.0	fragment-like	2	24	picked	22	1	True	Life Chemicals	F0807-0563	CC(C)c1ccc(cc1)/C=C/C2/c3cccc3NC2=O	45809595
	COc1cc2c(ncn2c1OC)Nc1cccc(c1)C	[4.267]	2.60	295.336	1864.0	414.0	fragment-like	4	36	picked	13	2	True	Maybridge	GK03474	Cc1cccc(c1)Nc2c3cc(c(c3nnc2)OC)OC	5828805
	FC(c1ccc(cc1)c1nnc2n1nc(cc2)NC1CC1)(F)F	[4.394]	4.13	319.284	3175.0	414.0	fragment-like	4	32	picked	21	2	True	Maybridge	RH00113	c1cc(ccc1c2nnc3n2nc(cc3)NC4CC4)(F)(F)F	5784088
	Nc1nncn2c1cn2c1ccc1	[3.869]	1.50	211.223	3430.0	414.0	fragment-like	1	31	picked	8	1	True	Maybridge	RJC00689	c1ccc(cc1)n2c3c(cn2)c(cn3)N	719540
	Clc1cccc(c1)Nc1nncn2c1ccc2.Cl	[4.05]	3.90	292.163	7366.0	414.0	fragment-like	2	36	picked	20	2	True	Maybridge	DP00818	c1ccc2c1)c(c1nnc2)Nc3cccc(c3)Cl.Cl	1859493
	Nc1nncn2c1nc2c(n1CC1CC1)ccc2	[4.082]	2.44	255.275	300.0	168.0	fragment-like	3	28	picked	12	2	True	Vitas M Labs	STL482461	c1ccc2c(c1)nc(n2CC3CC3)c4c(nnn4)N	8332960
	O=C(c1(c)C)cccc1C)Nc1cccc1	[4.714, 9.847]	2.93	267.111	385.0	275.0	fragment-like	3	20	picked	15	3	True	Vitas M Labs	STK098832	c1cc(c(c1)C)C(=O)Nc2cnc2)Cl	1284691

	<chem>OCCNc1ncnc2c1c(c1ccccc1)c(o2)c1ccccc1</chem>	[4.294]	3.08	331.36800000000005	159.0	168.0	fragment-like	5	71	picked	16	2	True	Vitas M Labs	STK012643	<chem>c1ccc(cc1)c2c3c(ncnc3oc2c4ccccc4)NCCO</chem>	1415762
	<chem>c1ccc(cc1)C5c1nnc(o1)c1ccncc1</chem>	[4.902]	3.31	269.322	170.0	168.0	fragment-like	4	30	picked	17	1	True	Vitas M Labs	STK032731	<chem>c1ccc(cc1)C5c2nnc(o2)c3ccncc3</chem>	1444229
	<chem>NC(=O)c1ccc2c(c1)nc([nH]2)c1ccccc1</chem>	[6.342]	2.19	237.257	2000.0	168.0	fragment-like	2	42	picked	11	2	True	Vitas M Labs	STL497402	<chem>c1ccc(cc1)c2[nH]c3ccc(cc3n2)C(=O)N</chem>	37053191
	<chem>Cc1ccc(cc1)[S](=O)(=O)Nc1cccc(c1)C(F)(F)F</chem>	[7.903]	3.75	315.311	527.0	168.0	fragment-like	4	24	picked	19	2	True	Vitas M Labs	STK053880	<chem>Cc1ccc(cc1)[S](=O)(=O)Nc2cccc(c2)C(F)(F)F</chem>	1377874
	<chem>O=C15/C(=C/c2ccc(o2)c2ccc(c(c2)C)F)/C(=O)N1</chem>	[8.052]	3.91	323.7270000000001	125.0	168.0	fragment-like	2	31	picked	20	1	True	Vitas M Labs	STL282834	<chem>c1cc(c(c1c2ccc(o2)/C=C/3/C(=O)NC(=O)S3)C)F</chem>	25775231
	<chem>COc1cc(ccc1OC)c1csc(c1C(=O)OC)C(K)N</chem>	[10.184]	4.08	321.391	10117.0	168.0	fragment-like	6	28	picked	21	1	True	Vitas M Labs	STK348738	<chem>CC(C)OC(=O)c1c1c(csc1N)c2ccc(c(c2)OC)OC</chem>	837243