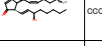
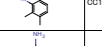
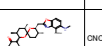
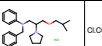
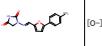
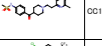


Structure	SMILES	BPAD	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	N-(2S,3R,4E)-1,3-dihydroxystearic acid	C18 Ceramide	537.9	1	ICCB Known Bioactives Library	24696-26-2	SL-115	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	2S,3R,4E-2-octylamino-1,3-diol	C8 Ceramide	411.7	1	ICCB Known Bioactives Library	170959-06-4	SL-145	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	(S)-1-hydroxy-3-(octadecyloxy)propan-2-yl decanoate	(S)-1-Methandamide	381.6	1	ICCB Known Bioactives Library	157182-49-5	FA-021	1mM
	[R]([C@H]1[C@@H]([C@H]1)C(=O)C(C)(C)(O)C(O)C)C(=O)COC2=C(C)C1=C(O)C(O)C(C)C1=O	1,3-bis(sn-3'-phosphatidyl)-sn-glycerol-3-phosphate	1,3-bis(sn-3'-phosphatidyl)-sn-glycerol-3-phosphate	418.6	1	ICCB Known Bioactives Library	32222-06-3	DM-200	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-(decyloxy)-3-hydroxypropan-2-yl decanoate	1,2-Didecanoyl glycerol (100)	400.6	1	ICCB Known Bioactives Library	82950-64-9	DG-100	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	2S-1-hydroxy-3-(octadecyloxy)propan-2-yl octanoate	1,2-Octanoyl-sn-glycerol	344.5	1	ICCB Known Bioactives Library	75685-80-2	DG-112	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-hydroxy-3-(2R)-octadec-9-enyloxy)propan-2-yl (2Z)-octadec-9-enoate	1,2-Dioleoyl glycerol (18:1)	621.0	1	ICCB Known Bioactives Library	3728-74-7	DG-115	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	(±)-11,12-Epoxyoctadecanoic acid	(±)-11,12-Epoxyoctadecanoic acid	320.5	1	ICCB Known Bioactives Library	123231-40-8	EE-11	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	12R-HETE	12R-HETE	320.5	1	ICCB Known Bioactives Library	82327-46-0	H-112	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	12S-HETE	12S-HETE	320.5	1	ICCB Known Bioactives Library	54397-83-0	H-012	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	12S-HPETE	12S-HPETE	320.5	1	ICCB Known Bioactives Library	71774-10-2	HP-012	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	12-methoxydecanoic acid	12-Methoxydecanoic acid	230.3	1	ICCB Known Bioactives Library	92169-28-3	G-234	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	13-HODE	13-HODE	296.4	1	ICCB Known Bioactives Library	28623-29-7	HO-013	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	13-HPODE	13-HPODE	312.4	1	ICCB Known Bioactives Library	33964-75-9	HP-013	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	13,14-Dihydro-PGE1	13,14-Dihydro-PGE1	356.5	1	ICCB Known Bioactives Library	19313-29-1	PG-048	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	13-cis-retinoic acid	13-cis-retinoic acid	300.4	1	ICCB Known Bioactives Library	4759-48-2	GR-102	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	13-Keto-octadeca-9,11-dienoic acid	13-Keto-octadeca-9,11-dienoic acid	294.4	1	ICCB Known Bioactives Library	54739-30-9	KO-013	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	(±)-14,15-Epoxyoctadecanoic acid	(±)-14,15-Epoxyoctadecanoic acid	320.5	1	ICCB Known Bioactives Library	81276-03-1	EE-14	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	15S-HETE	15S-HETE	320.5	1	ICCB Known Bioactives Library	54845-95-3	H-015	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	15S-HPETE	15S-HPETE	320.5	1	ICCB Known Bioactives Library	70981-96-3	HP-015	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	15-deoxy-Prostaglandin J2	15-deoxy-Prostaglandin J2	316.4	1	ICCB Known Bioactives Library	87893-55-8	PG-050	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	15-Ketosteatraenoic acid	15-Ketosteatraenoic acid	318.5	1	ICCB Known Bioactives Library	81416-72-0	K-015	0.1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	16,16-Dimethyl-prostaglandin E2	16,16-Dimethyl-prostaglandin E2	380.5	1	ICCB Known Bioactives Library	38746-25-3	PG-021	1mM
	[R]([C@H]1[C@@H]([C@H]1)C(=O)C(C)(C)C(O)C)C(=O)C	17-Octadecynoic acid	17-Octadecynoic acid	280.4	1	ICCB Known Bioactives Library	34450-19-5	E1-103	1mM
	[R]([C@H]1[C@@H]([C@H]1)C(=O)C(C)(C)C(O)C)C(=O)C	17-Phenyl-tri-nor-prostaglandin E2	17-Phenyl-tri-nor-prostaglandin E2	386.5	1	ICCB Known Bioactives Library	38315-43-4	PG-042	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Acyl-PAF	1-Acyl-PAF	537.7	1	ICCB Known Bioactives Library	84062-61-3	L-134	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Hexadecyl-2-arachidonyl-glycerol	1-Hexadecyl-2-arachidonyl-glycerol	603.0	1	ICCB Known Bioactives Library	126721-87-7	DG-140	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Hexadecyl-2-methylglycero-3-PC	1-Hexadecyl-2-methylglycero-3-PC	495.7	1	ICCB Known Bioactives Library	78958-44-3	L-106	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Hexadecyl-2-O-acetyl-glycerol	1-Hexadecyl-2-O-acetyl-glycerol	398.6	1	ICCB Known Bioactives Library	77133-35-8	E1-150	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Hexadecyl-2-O-methyl-glycerol	1-Hexadecyl-2-O-methyl-glycerol	330.5	1	ICCB Known Bioactives Library	84327-41-7	E1-149	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Octadecyl-2-methylglycero-3-PC	1-Octadecyl-2-methylglycero-3-PC	523.7	1	ICCB Known Bioactives Library	77286-66-9	L-108	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Oleoyl-2-acetyl-glycerol	1-Oleoyl-2-acetyl-glycerol	398.6	1	ICCB Known Bioactives Library	88390-77-4	DG-125	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Stearoyl-2-oleoyl-glycerol	1-Stearoyl-2-oleoyl-glycerol	621.0	1	ICCB Known Bioactives Library	34487-26-8	DG-135	1mM
	CCCCCCCC(O)C(O)C(O)CCCCCCCC	1-Stearoyl-2-arachidonyl-glycerol	1-Stearoyl-2-arachidonyl-glycerol	645.0	1	ICCB Known Bioactives Library	65914-64-3	DG-130	1mM





Structure	SMILES	BPAD	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	<chem>[O][C@@H]1[C@@H](O)[C@@H](O)[C@@H](O)C(C)C(O)C1(O)C(C)C(C)C(C)C</chem>	3R,6R-6-[(1R,3aS,4E,7aR)-4-[2-[(1Z,5Z)-5-hydroxy-2-methylenecyclohexylidene]ethylidene]-7a-methyl-octahydro-1H-inden-1-yl]-2-methylheptane-2,3-diol	24,25-Dihydroxyvitamin D3	416.6	1	ICCB Known Bioactives Library	40013-87-4	DM-300	1uM
	<chem>[O][C@@H]1[C@@H](O)[C@@H](O)C(C)C(O)C1(O)C(C)C(C)C(C)C</chem>	1S,3Z)-3-[2-[(1R,3aS,4E,7aR)-1-[(2R)-6-hydroxy-6-methylheptan-2-yl]-7a-methyl-octahydro-1H-inden-4-ylidene]ethylidene]-4-methylenecyclohexan-1-ol	25-Dihydroxyvitamin D3	400.6	1	ICCB Known Bioactives Library	19356-17-3	DM-100	1uM
	<chem>CCCCCCCCCCCCCCCC(=O)COC</chem>	1,3-dihydroxypropan-2-yl (3Z,8Z,11Z,14Z)-icoso-5,8,11,14-tetraenoate	2-Arachidonylglycerol	378.5	1	ICCB Known Bioactives Library	53847-30-6	FA-028	1uM
	<chem>CCCCCCCCCCCC(F)C(=O)O</chem>	2-fluorohexadecanoic acid	2-Fluoropalmitic acid	274.4	1	ICCB Known Bioactives Library	89270-22-4	G-231	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	2-hydroxytetradecanoic acid	2-Hydroxytridecanoic acid	244.4	1	ICCB Known Bioactives Library	2507-55-3	G-230	1uM
	<chem>CC(=O)Nc1ccc(O)cc1</chem>	2E,4E,6E,8E)-N-(4-hydroxyphenyl)-3,7-dimethyl-9-(2,6,6-trimethylcyclohex-1-en-1-yl)nona-2,4,6,8-tetraenamide	4-Hydroxyphenylacetamide	391.5	1	ICCB Known Bioactives Library	6546-68-6	GR-103	1uM
	<chem>CCCCCCCCCCCCCCCCC(=O)O</chem>	3-(decyl)propanoic acid	4-Octadecanoic acid	230.3	1	ICCB Known Bioactives Library	7420-16-8	G-232	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	5S,8E,12Z,14Z)-5-hydroxyicoso-6,8,11,14-tetraenoic acid	5S-HETE	320.5	1	ICCB Known Bioactives Library	70608-72-9	H-005	0.1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	5S,8E,12Z,14Z)-5-hydroperoxyicoso-6,8,11,14-tetraenoic acid	5S-HPETE	320.5	1	ICCB Known Bioactives Library	71774-08-6	HP-005	0.1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	4-[3-[(2Z,5Z)-tetradeca-2,5,9-trien-1-yl]oxiran-2-yl]butanoic acid	5,6-Epoxyicosatrienoic acid	320.5	1	ICCB Known Bioactives Library	184489-43-5	EE-56	0.1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	icoso-5,8,11,14-tetraenoic acid	5,8,11,14-Eicosatetraenoic acid	296.4	1	ICCB Known Bioactives Library	1191-85-1	ET-004	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	icoso-5,8,11-trienoic acid	5,8,11-Eicosatrynoic acid	300.4	1	ICCB Known Bioactives Library	13488-22-7	ET-003	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	6E,8Z,11Z,14Z)-5-oxoicoso-6,8,11,14-tetraenoic acid	5-Ketoicosatetraenoic acid	318.5	1	ICCB Known Bioactives Library	106154-18-1	K-005	0.1uM
	<chem>[O]C=Oc1c[nH]c2c3c[nH]c4c3c[nH]c41</chem>	9H-indol-3-ylcarbazole-6-carbaldehyde	6-Formylindole [3,2-b]carbazole	284.3	1	ICCB Known Bioactives Library	172922-91-7	GR-206	1uM
	<chem>CCCC(C)C(O)C(=O)O</chem>	7-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxyoct-1-en-1-yl]cyclopentyl]-6-oxoheptanoic acid	6-Keto-prostaglandin F _{2α}	370.5	1	ICCB Known Bioactives Library	58962-34-8	PG-017	1uM
	<chem>CCCCCCCCCCCC(C)(C)C(=O)O</chem>	3Z,8Z)-7,7-dimethylcoso-5,9-dienoic acid	7,7-Dimethylcosadecanoic acid	338.6	1	ICCB Known Bioactives Library	89560-01-0	EI-112	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	5Z)-7-[3-[(2Z,5Z)-undeca-2,5-dien-1-yl]oxiran-2-yl]hept-5-enoic acid	8,9-Epoxyicosatrienoic acid	320.5	1	ICCB Known Bioactives Library	81246-65-7	EE-89	0.1uM
	<chem>CCCC(C)C(O)C(=O)O</chem>	5Z)-7-[(1S,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxyoct-1-en-1-yl]cyclopentyl]hept-5-enoic acid	8-epi-Prostaglandin F _{2α}	354.5	1	ICCB Known Bioactives Library	27415-26-5	PG-048	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	9R,10E,12Z)-9-hydroxyoctadeca-10,12-dienoic acid	9S-HODE	296.4	1	ICCB Known Bioactives Library	72543-67-6	HO-409	0.1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	9S,10E,12Z)-9-hydroperoxyoctadeca-10,12-dienoic acid	9S-HPODE	312.4	1	ICCB Known Bioactives Library	29774-12-7	HP-019	0.1uM
	<chem>CCCCCCCCCCCCCNC=O</chem>	9Z)-octadec-9-enamide	9,10-Octadecanamide	281.5	1	ICCB Known Bioactives Library	301-02-0	FA-025	1uM
	<chem>CCCC(C)C(O)C(=O)O</chem>	5Z)-7-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxyoct-1-en-1-yl]cyclopentyl]hept-5-enoic acid	9a,11b-Prostaglandin F ₂	354.5	1	ICCB Known Bioactives Library	38432-87-0	PG-031	1uM
	<chem>CCCC(C)C(O)C(=O)O</chem>	5Z)-7-[(1R,2R,3R,5R)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxyoct-1-en-1-yl]cyclopentyl]hept-5-enoic acid	9a,11a-Prostaglandin F ₂	354.5	1	ICCB Known Bioactives Library	4510-16-1	PG-030	1uM
	<chem>CC(=O)Nc1ccc(O)cc1</chem>	2E,4E,6E,8E)-3,7-dimethyl-9-(2,6,6-trimethylcyclohex-1-en-1-yl)nona-2,4,6,8-tetraenoic acid	9-cis Retinoic acid	300.4	1	ICCB Known Bioactives Library	5300-03-6	GR-101	1uM
	<chem>CCOC(=O)C(O)C(=O)OC</chem>	methyl 2-(2-methoxy-2-oxoacetamido)acetate	Dimethylsalsalicylate	175.1	1	ICCB Known Bioactives Library	89464-63-1	EI-347	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	7Z,10Z,13Z,16Z)-docosa-7,10,13,16-tetraenoic acid	Adrenic acid (22:4, n-6)	322.5	1	ICCB Known Bioactives Library	28974-58-0	FA-011	1uM
	<chem>CC(=O)Nc1ccc(O)cc1</chem>	2R)-3-[EZE]-3,7-dimethyl-2,6-dien-1-yl]sulfonyl-2-acetamidopropanoic acid	N-Acetyl-S-garanyl-L-cysteine	290.4	1	ICCB Known Bioactives Library	135784-50-8	G-221	1uM
	<chem>CC(=O)Nc1ccc(O)cc1</chem>	2R)-2-acetamido-3-[(2E,6E,10E)-3,7,11,15-tetramethylhexadeca-2,6,10,14-tetraen-1-yl]sulfonylpropanoic acid	N-acetyl-S-garany-L-Cysteine	435.7	1	ICCB Known Bioactives Library	135332-94-8	G-220	1uM
	<chem>CC1=CC=NC=C1</chem>	1-(2,4-dichlorophenyl)-5-(4-iodophenyl)-4-methyl-N-(piperidin-1-yl)-1H-pyrazole-3-carboxamide	AM-251	552.2	1	ICCB Known Bioactives Library	18323-66-6	GR-108	1uM
	<chem>CC1=CC=NC=C1</chem>	4-(5S,8,8-tetramethyl-5,8,7,8-tetrahydrophthalene-2-amido)benzoic acid	AM-580	351.4	1	ICCB Known Bioactives Library	102121-60-8	GR-104	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	8Z,12Z)-N-(2-hydroxyethyl)octadeca-9,12-dienamide	Anandamide (18:2, n-6)	323.5	1	ICCB Known Bioactives Library	68171-52-8	FA-026	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	8Z,11Z,14Z)-N-(2-hydroxyethyl)icoso-6,11,14-trienamide	Anandamide (20:3, n-6)	349.6	1	ICCB Known Bioactives Library	150314-34-4	FA-019	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	5Z,8Z,11Z,14Z)-N-(2-hydroxyethyl)icoso-5,8,11,14-tetraenamide	Anandamide (20:4, n-6)	347.5	1	ICCB Known Bioactives Library	84421-68-8	FA-017	1uM
	<chem>CCCCCCCCCCCC(O)C(=O)O</chem>	7Z,10Z,13Z,16Z)-N-(2-hydroxyethyl)docosa-7,10,13,16-tetraenamide	Anandamide (22:4, n-6)	375.6	1	ICCB Known Bioactives Library	150314-35-5	FA-020	1uM

Structure	SMILES	BPAD	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	<chem>CCCCCCCCCCCCCCCC(=O)N</chem>		Arachidonamide	302.5	1	ICGB Known Bioactives Library	85146-53-6	FA-022	1mM
	<chem>CCCCCCCC=CCCC=CCCC=CCCC(=O)O</chem>		Arachidonic acid (20:4, n-6)	304.5	1	ICGB Known Bioactives Library	506-32-1	FA-003	1mM
	<chem>CCCCCCCCCCCCCCCC(COP(=O)([O-])[O-])CCCCCCCCCCCCCCCC</chem>		Arachidonyl-PAF	768.1	1	ICGB Known Bioactives Library	86288-11-1	L-102	1mM
	<chem>COCC1=CC2=CC(=O)C3=CC(=O)C(C)O3C1=CC(=O)O</chem>		BML-190	426.9	1	ICGB Known Bioactives Library	2854-32-2	CR-104	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		C2 Ceramide	341.5	1	ICGB Known Bioactives Library	3102-57-6	SL-100	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		C2 Dihydroceramide	343.5	1	ICGB Known Bioactives Library	13031-64-6	SL-101	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		C8 Ceramide	425.7	1	ICGB Known Bioactives Library	74713-59-0	SL-112	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		C8 Dihydroceramide	427.7	1	ICGB Known Bioactives Library	145714-33-0	SL-113	1mM
	<chem>CCCC(C)CCC(O)C(=O)C</chem>		Carbonyl	350.5	1	ICGB Known Bioactives Library	69552-46-1	PG-035	1mM
	<chem>CC1=CC(C2=CC(=O)O2)C(=O)C1</chem>		Cilgitezone	333.4	1	ICGB Known Bioactives Library	74772-77-3	GR-205	1mM
	<chem>CCCC(O)CC(C)OCC1=CC=CC=C1C1</chem>		Cilfibrate	242.7	1	ICGB Known Bioactives Library	637-07-0	EI-160	1mM
	<chem>O=C(O)C=C(C)C(=O)O</chem>		Cloprostenol	424.9	1	ICGB Known Bioactives Library	55028-72-3	RA-106	1mM
	<chem>CCCCCCCCCCCCCCCC(COP(=O)([O-])[O-])CCCCCCCCCCCCCCCC</chem>		C-PAF	538.7	2	ICGB Known Bioactives Library	51575-58-5	L-120	1mM
	<chem>CCCC(C)CCC(O)C(=O)C</chem>		D12-Protaglandin I ₂	334.4	2	ICGB Known Bioactives Library	87893-54-7	PG-047	1mM
	<chem>CCCCCCCC=CCCC=CCCC=CCCC(=O)O</chem>		Dhomo-gamma-linolenic acid	306.5	2	ICGB Known Bioactives Library	1783-84-2	FA-009	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		Dihydrospingosine	301.5	2	ICGB Known Bioactives Library	3102-56-5	SL-125	1mM
	<chem>O=C1=CC2=CC(=O)C3=CC(=O)C(C)O3C2=CC=C1</chem>		Didolymethane	246.3	2	ICGB Known Bioactives Library	1988-05-4	GR-207	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		DL-Dihydrospingosine	301.5	2	ICGB Known Bioactives Library	73938-69-9	SL-205	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		DL-PPMP	390.6	2	ICGB Known Bioactives Library	72527-80-4	SL-210	1mM
	<chem>CCCCCCCCCCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		DL-PPMP	474.7	2	ICGB Known Bioactives Library	149022-18-4	SL-215	1mM
	<chem>CCCCCCCC=CCCC=CCCC=CCCC=CCCC(=O)O</chem>		Docosahexaenoic acid (22:6 n-3)	328.5	2	ICGB Known Bioactives Library	6217-54-5	FA-002	1mM
	<chem>CCCCCCCC=CCCC=CCCC=CCCC(=O)O</chem>		Docosapentaenoic acid	330.5	2	ICGB Known Bioactives Library	2234-74-4	FA-012	1mM
	<chem>CCCCCCCC=CCCC=CCCC(=O)O</chem>		Docosatrienoic acid (22:3 n-3)	334.5	2	ICGB Known Bioactives Library	28845-86-5	FA-010	1mM
	<chem>CCCCCCCC=CCCC=CCCC(=O)O</chem>		Eicosa-5,8-dienoic acid (20:2 n-12)	308.5	2	ICGB Known Bioactives Library	2555-07-1	FA-024	1mM
	<chem>CCCCCCCC=CCCC=CCCC(=O)O</chem>		Eicosadienoic acid (20:2 n-6)	308.5	2	ICGB Known Bioactives Library	2091-39-6	FA-007	1mM
	<chem>CCCCCCCC=CCCC=CCCC=CCCC(=O)O</chem>		Eicosapentaenoic acid (20:5 n-3)	302.5	2	ICGB Known Bioactives Library	10417-94-4	FA-001	1mM
	<chem>CCCCCCCC=CCCC=CCCC(=O)O</chem>		Eicosatrienoic acid (20:3 n-3)	306.5	2	ICGB Known Bioactives Library	2091-27-2	FA-008	1mM
	<chem>CCCCCCCCCCCCCCCC(COP(=O)([O-])[O-])CCCCCCCCCCCCCCCC</chem>		Enantiomeric PAF C16	523.7	2	ICGB Known Bioactives Library	117985-53-6	L-110	1mM
	<chem>CCC(O)CCCC(O)C(=O)O</chem>		Farnesylthioctic acid	296.5	2	ICGB Known Bioactives Library	13784-48-4	G-222	1mM
	<chem>O=C(O)C=C(C)C(=O)O</chem>		Fluprostenol	458.5	2	ICGB Known Bioactives Library	40666-16-0	RA-105	1mM
	<chem>CCCCCCCC=CCCC=CCCC(=O)O</chem>		Gamma-linolenic acid (18:3 n-6)	278.4	2	ICGB Known Bioactives Library	506-20-3	FA-006	1mM
	<chem>CCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		Leukotain A (8:10-EODE)	296.4	2	ICGB Known Bioactives Library	6814-92-4	LE-009	0.1mM
	<chem>CCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		Leukotain B (12:13-EODE)	296.4	2	ICGB Known Bioactives Library	848310-68-7	LE-012	0.1mM
	<chem>CCCCCCCC(O)CCCCCCCCCCCCCCCC</chem>		Leukotriene B4	338.5	2	ICGB Known Bioactives Library	71160-24-2	LB-004	0.1mM

Structure	SMILES	BPAD	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	CCCCC=CC/C=C/C/C=C/C([C@@H](S[C@@H](NC(=O)CC[C@@H](NCO)O)C(=O)NCCC)O)C(=O)O	55.6R.7E.9E.11Z.14Z)-6-[[2R]-2-[(4S)-4-amino-4-carboxybutanamido]-2-[(carboxymethyl)carbamoyl]sulfonyl]-5-hydroxyocta-7,8,11,14-tetraenoic acid	Leukotriene C4	625.9	2	ICDB Known Bioactives Library	72025-60-6	LC-004	0.1mM
	CCCCC=CC/C=C/C/C=C/C([C@@H](S[C@@H](NC(=O)O)C(=O)O)C(=O)O)O	55.6R.7E.9E.11Z.14Z)-6-[[2R]-2-amino-2-[(carboxymethyl)carbamoyl]sulfonyl]-5-hydroxyocta-7,8,11,14-tetraenoic acid	Leukotriene D4	496.7	2	ICDB Known Bioactives Library	73836-78-9	LD-004	0.1mM
	CCCCC=CC/C=C/C/C=C/C([C@@H](S[C@@H](NCO)O)C(=O)O)O	55.6R.7E.9E.11Z.14Z)-6-[[2R]-2-amino-2-carboxyethylsulfonyl]-5-hydroxyocta-7,8,11,14-tetraenoic acid	Leukotriene E4	438.6	2	ICDB Known Bioactives Library	75715-69-8	LE-004	0.1mM
	CCCCC=CC/C=C/C/C=CCCCCCCC(N)=O	9Z,12Z)-octadeca-9,12-dienamide	Linoleamide	279.5	2	ICDB Known Bioactives Library	3072-13-7	FA-023	1mM
	CCCCC=CC/C=C/C/C=CCCCCCCC(=O)O	9Z,12Z)-octadeca-9,12-dienoic acid	Linoleic acid	280.4	2	ICDB Known Bioactives Library	60-33-3	FA-014	1mM
	CC=CC/C=C/C/C=C/C/C=CCCCCCCC(=O)O	9Z,12Z,15Z)-octadeca-9,12,15-trienoic acid	Linolenic acid (18:3 n-3)	278.4	2	ICDB Known Bioactives Library	463-40-1	FA-005	1mM
	CCCCC[C@@H](O)C(C)=CC/C=C/C/C=C/C([C@@H](O)C(=O)O)O	55.6R.7E.9E.11Z.13E.15S)-5,8,15-trihydroxyocta-7,8,11,13-tetraenoic acid	Lipoin A4	352.5	2	ICDB Known Bioactives Library	89663-86-5	LXA-004	0.1mM
	CCCCC=CC(CCCC)C(C)N(C(=O)N(C)C)C(=O)O	2S)-2-hexadecanamide-3-(phosphonooxy)propanoic acid	L-NASPA	423.5	2	ICDB Known Bioactives Library	155915-46-1	LP-101	1mM
	CCCC1=C(O)CCCCOCC2=NNHNN2C(=O)O=C1	1-(2-hydroxy-3-propyl-4-[(4S)-1H-1,2,3,4-tetrazol-5-yl]butoxy)ethan-1-one	LY-171883	318.4	2	ICDB Known Bioactives Library	88107-10-2	RA-101	1mM
	CCCCC=CC(CCCC)C(C)OP(O)(=O)C1=CC=CC=C1	2-[[2R]-3-(hexadecyloxy)-2-hydroxypropyl phosphonate]oxyethyltrimethylazanium	Lyso-PAF C16	431.6	2	ICDB Known Bioactives Library	52691-62-0	L-101	1mM
	[Na+]CCCC=CC(CCCC)C(C)OP(O)(=O)C(=O)O	sodium [2R]-3-(hydrogen phosphonate)-2-hydroxypropyl [9Z]-octadec-9-enoate	Lyso-phosphatidic acid Na	438.5	2	ICDB Known Bioactives Library	325465-93-9	LP-100	1mM
	CCCCC=CC(CCCC)C(C)OP(O)(=O)C(=O)O	N-[(1S,2R)-1-hydroxy-1-phenylpropan-2-yl]tetradecanamide	D-erythro-MAPP	361.6	2	ICDB Known Bioactives Library	66085-59-4	SL-221	1mM
	CCCCC=CC(CCCC)C(C)OP(O)(=O)C(=O)O	N-[(1R,2S)-1-hydroxy-1-phenylpropan-2-yl]tetradecanamide	L-erythro-MAPP	361.6	2	ICDB Known Bioactives Library	143482-38-0	SL-222	1mM
	CCCC=CC/C=C/C/C=C/C/C=CCCCCCCC(=O)O	5Z,8Z,11Z)-icosa-5,8,11-trienoic acid	Mead acid (20:3 n-6)	306.5	2	ICDB Known Bioactives Library	20590-32-3	FA-004	1mM
	CCCC=CC/C=C/C/C=C/C/C=CCCCCCCC(N)=O	5Z,8Z,11Z)-N-(2-hydroxyethyl)icosa-5,8,11-trienamide	Mead ethanolamide	348.6	2	ICDB Known Bioactives Library	169232-04-6	FA-027	1mM
	CCCC(C)CCCC(C)C/C=C/C/C=C/C/C(=O)O	2E,4E)-11-methoxy-3,7,11-trimethyldeca-2,4-dienoic acid	Methoprene acid	268.4	2	ICDB Known Bioactives Library	53082-52-7	GR-106	1mM
	CCCCC(C)C(C)C=C(C)OP(O)(=O)C(=O)O	7-[[1R,2R,3R]-3-hydroxy-2-[[1E]-4-hydroxy-4-methylbut-1-en-1-yl]-5-oxocyclopentyl]heptanoic acid	Misoprostol, free acid	368.5	2	ICDB Known Bioactives Library	59122-46-2	PG-052	1mM
	CCCCC(C)C(C)C=C(C)OP(O)(=O)C(=O)O	2S,3R,4E)-2-(dimethylamino)octadec-4-ene-1,3-diol	N,N-Dimethylsphingosine	327.5	2	ICDB Known Bioactives Library	119567-63-4	SL-105	1mM
	CCCCC=CC/C=C/C/C=C/C([C@@H](S[C@@H](NC(=O)C)O)C(=O)O)O	55.6R.7E.9E.11Z.14Z)-6-[[2-carboxy-2-acetamidethylsulfonyl]-5-hydroxyocta-7,8,11,14-tetraenoic acid	N-Acetyl-leukotriene E4	481.6	2	ICDB Known Bioactives Library	80115-95-3	LEA-004	0.1mM
	CCCCC=CC/C=C/C/C=C/C/C=CCCCCCCC(=O)O	2-[[5Z,8Z,11Z,14Z]-icosa-5,8,11,14-tetraenamide]acetic acid	N-arachidonylglycine	381.5	2	ICDB Known Bioactives Library	179113-91-8	FA-029	1mM
	CCCCC=CC/C=C/C/C=CCCCCCCC(=O)O	2-[[9Z,12Z)-octadeca-9,12-dienamide]acetic acid	N-isooleoylglycine	337.5	2	ICDB Known Bioactives Library	2764-03-6	FA-030	1mM
	CCCCC=CC(CCCC)C(C)OP(O)(=O)C1=CC=CC=C1	2-[[2R]-2-(acetoxy)-3-(hexadecyloxy)propyl phosphonate]oxyethyltrimethylazanium	PAF C16	523.7	2	ICDB Known Bioactives Library	74389-68-7	L-100	1mM
	CCCCC=CC(CCCCC)C(C)OP(O)(=O)C1=CC=CC=C1	2-[[2R]-2-(acetoxy)-3-(octadecyloxy)propyl phosphonate]oxyethyltrimethylazanium	PAF C18	551.7	2	ICDB Known Bioactives Library	79549-29-1	L-105	1mM
	CCCCC=CC(CCCCC)C(C)OP(O)(=O)C1=CC=CC=C1	2-[[2R]-2-(acetoxy)-3-[[9Z]-octadec-9-en-1-yl]oxy]propyl phosphonate]oxyethyltrimethylazanium	PAF C18:1	548.7	2	ICDB Known Bioactives Library	86288-90-1	L-107	1mM
	CCCCC=CC(CCCC)C(C)OP(O)(=O)O	N-(2-hydroxyethyl)hexadecanamide	Palmitolethanolamide	299.5	2	ICDB Known Bioactives Library	544-31-0	FA-018	1mM
	CCCCC(C)C(C)C=C(C)OP(O)(=O)C(=O)O	2,3-bis(hexadecanoyloxy)propane-1,3-diphosphonic acid	Dipalmitoylphosphatidic acid	648.9	2	ICDB Known Bioactives Library	71065-67-7	LP-103	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	7-[[1R,2S]-2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-5-oxocyclopent-3-en-1-yl]heptanoic acid	Prostaglandin A1	336.5	2	ICDB Known Bioactives Library	14152-28-4	PG-001	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	5Z)-7-[[1R,2S]-2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-5-oxocyclopent-3-en-1-yl]hept-5-enoic acid	Prostaglandin A2	334.4	2	ICDB Known Bioactives Library	13345-50-1	PG-002	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	7-(2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-5-oxocyclopent-1-en-1-yl]heptanoic acid	Prostaglandin B1	336.5	2	ICDB Known Bioactives Library	13345-51-2	PG-003	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	5Z)-7-(2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-5-oxocyclopent-1-en-1-yl]hept-5-enoic acid	Prostaglandin B2	334.4	2	ICDB Known Bioactives Library	13387-65-6	PG-004	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	5Z)-7-[[1R,2R,5S]-5-hydroxy-2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-3-oxocyclopentyl]hept-5-enoic acid	Prostaglandin D2	352.5	2	ICDB Known Bioactives Library	41598-07-6	PG-005	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	7-[[1R,2R,3R]-3-hydroxy-2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-5-oxocyclopentyl]heptanoic acid	Prostaglandin E1	354.5	2	ICDB Known Bioactives Library	745-65-3	PG-006	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	5Z)-7-[[1R,2R,3R]-3-hydroxy-2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]-5-oxocyclopentyl]hept-5-enoic acid	Prostaglandin E2	352.5	2	ICDB Known Bioactives Library	363-24-6	PG-007	1mM
	CCCCC[C@@H](O)C=C(C)OP(O)(=O)C(=O)O	7-[[1R,2R,3R,5S]-2,5-dihydroxy-2-[[1E,3S]-3-hydroxyoct-1-en-1-yl]cyclopentyl]heptanoic acid	Prostaglandin F2a	358.5	2	ICDB Known Bioactives Library	745-62-0	PG-010	1mM

Structure	SMILES	BPAD	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	CCCCO[C@@H](O)C[C@H](O)C[C@H](O)C[C@H](O)C(=O)O	SZ-7-[[1R,3R,5S]-5-hydroxy-2-[(1E,3S)-3-hydroxyoct-1-en-1-yl]cyclopentyl]hept-5-enoic acid	Prostaglandin F _{2α}	354.5	2	ICCB Known Bioactives Library	551-11-1	PG-008	1nM
	CCCCO[C@@H](O)C[C@H](O)C[C@H](O)C[C@H](O)C(=O)O	5-[[2Z,3R,4R,5R,6S]-5-hydroxy-4-[(1E,3S)-3-hydroxyoct-1-en-1-yl]heptahydro-2H-cyclopenta[b]furan-2-ylidene]pentanoic acid	Prostaglandin L ₆	352.5	2	ICCB Known Bioactives Library	61849-14-7	PG-011	1nM
	CCCCO[C@@H](O)C[C@H](O)C[C@H](O)C(=O)O	SZ-7-[[1S,5R]-5-[(1E,3S)-3-hydroxyoct-1-en-1-yl]-4-oxocyclopent-2-en-1-yl]hept-5-enoic acid	Prostaglandin J ₂	354.4	2	ICCB Known Bioactives Library	60203-57-8	PG-023	1nM
	CC(C)C=CC(=C)C=CC(=O)O	2E,4E,6E,8E)-3,7-dimethyl-9-(2,6,8-trimethylcyclohex-1-en-1-yl)nona-2,4,6,8-tetraenoic acid	All trans retinoic acid	300.4	2	ICCB Known Bioactives Library	302-79-4	GR-100	1nM
	CCCCOC1C(=O)C(C)C(=O)N(C)C1	1-[2-(quinidin-2-ylmethoxy)phenyl]hexan-1-ol	REV-5901	335.4	2	ICCB Known Bioactives Library	101910-24-1	EI-209	1nM
	CCCC(C)SC(C)C	methyl [(2R)-2-amino-3-[[[2E,6E]-3,7,11-trimethyldec-2,6,10-trien-1-yl]sulfonyl]propanoate	S-farnesyl-L-cysteine	339.5	2	ICCB Known Bioactives Library		G-201	1nM
	CCCC(C)C(N)C(C)C(C)O	2S,3R,4E)-2-aminooctadec-4-ene-1,3-diol	Sphingosine	299.5	2	ICCB Known Bioactives Library	123-78-4	EI-155	1nM
	CC1=CC=C(C=C1)C(=O)O	SZ-7-[[2-(phenylcarbamoyl)amino]methyl]-7-oxabicyclo[2.2.1]heptan-2-yl]hept-5-enoic acid	SO-39548	367.5	2	ICCB Known Bioactives Library	88298-61-7	RA-103	1nM
			Blank		2	ICCB Known Bioactives Library			
	CC1=CC=C(C=C1)C(=O)O	4-[(1E)-2-(5S,8R)-tetramethyl-5,8,7,8-tetrahydroaphthalen-2-yl]prop-1-en-1-yl]benzoic acid	TNPB	348.5	2	ICCB Known Bioactives Library	71441-28-6	GR-105	1nM
	CCCCO[C@@H](O)C[C@H](O)C[C@H](O)C(=O)O	SZ-7-[[1R,4S,5S,6R]-6-[(1E,3S)-3-hydroxyoct-1-en-1-yl]-2-oxabicyclo[2.2.1]heptan-5-yl]hept-5-enoic acid	U-46619	350.5	2	ICCB Known Bioactives Library	58985-40-1	PG-023	1nM
	CCCCO[C@@H](O)C[C@H](O)C[C@H](O)C(=O)O	6-[[1E,5Z]-3-hydroxyundec-1,5-dien-1-yl]pyridin-2-yl]hexane-1,5-diol	U-75302	361.5	2	ICCB Known Bioactives Library	119477-85-9	RA-102	0.1nM
	CS(=O)(=O)O	11R)-2-methyl-11-(morpholin-4-ylmethyl)-3-naphthalene-1-carboxylate	WN 55212-2 mesylate	529.8	2	ICCB Known Bioactives Library	131543-23-2	OR-105	1nM
	CC1=CC=C(C=C1)C(=O)O	2-[[6-chloro-6-[[2,3-dimethylphenyl]amino]pyridin-2-yl]sulfonyl]acetic acid	WY-14643	323.8	2	ICCB Known Bioactives Library	50882-23-4	GR-200	1nM
	NC1=CC=C(N)C=C1	pyridin-4-amine	4-Aminopyridine	94.1	2	ICCB Known Bioactives Library	504-24-5	AC-119	5ng/mL
	CCCCO[C@@H](O)C[C@H](O)C[C@H](O)C(=O)O	5-hydroxydecanic acid	5-Hydroxydecanate	188.5	2	ICCB Known Bioactives Library	624-00-0	KC-141	5ng/mL
	[H][C@H]1[C@@H]2[C@H]3[C@H]4[C@@H](O)C=C(O)C=C4[C@H]5[C@@H](O)C=C(C)C=C5[C@H]6[C@@H](O)C=C(C)C=C6[C@H]7[C@@H](O)C=C(C)C=C7[C@H]8[C@@H](O)C=C(C)C=C8[C@H]9[C@@H](O)C=C(C)C=C9[C@H]10[C@@H](O)C=C(C)C=C10	2R,3R,4S,5R,6S,7S,8R,9R,10R,11R,14R,16R,18R)-6-(acetyl-11-ethyl-5,7,14-trihydroxy-8,16,18-trimethoxy-13-(methoxymethyl)-11-azabicyclo[7.2.2.1 ^{0,9'} .0 ^{1,10'} .0 ^{2,11'}]-nonadeca-4-yl]benzoate	Aconitine	645.7	2	ICCB Known Bioactives Library	302-27-2	AC-128	5ng/mL
	ClC1=CC=C(C=C1)C(=O)O	N-(2-(2-chlorophenylethyl)amino)-2-hydroxypropan-1-yl]methanesulfonamide hydrochloride	AM 8016 HCl	483.8	2	ICCB Known Bioactives Library	178884-81-0	AC-115	5ng/mL
	ClC1=CC=C(C=C1)C(=O)O	adamantan-1-amine hydrochloride	Amantadine HCl	187.7	2	ICCB Known Bioactives Library	665-66-7	AC-133	5ng/mL
	ClN1C=CC(=O)N1	3,5-diamino-N-carbamimidyl-6-chloropyrazine-2-carboxamide hydrochloride	Amiloride HCl	286.1	2	ICCB Known Bioactives Library	2016-88-8	CA-200	5ng/mL
	CC1=CC=C(C=C1)C(=O)O	2-(4-(2-butyl-1-benzofuran-3-carboxyl)-2,6-diodophenoxy)ethyl]diethylamine hydrochloride	Anisodrone HCl	681.8	2	ICCB Known Bioactives Library	19774-62-4	AC-105	5ng/mL
	OC1=CC=C(C=C1)C(=O)O	5-(methylamino)-2-[[2S,3S,8S,9R]-3,8,11-trimethyl-8-(1-oxo-1-(1H-pyridin-2-yl)propan-2-yl)-1,7-dioxaspiro[5.5]undecan-2-yl]methyl-1,3-benzoxazole-4-carboxylic acid	A-23187	523.8	2	ICCB Known Bioactives Library	52665-69-7	CA-100	5ng/mL
	COc1ccc(C)c(O)c1	methyl 2,6-dimethyl-5-nitro-4-[[2-(trifluoromethyl)phenyl]-1,4-dihydropyridin-3-yl]carboxylate	(R)-(-)-BAY K-8644	358.3	2	ICCB Known Bioactives Library	88791-67-4	CA-410	5ng/mL
	ClN1C=CC(=O)N1	3,5-diamino-N-[(E)-N-benzylcarbamimidyl]-6-chloropyrazine-2-carboxamide hydrochloride	Benzamide HCl	358.2	2	ICCB Known Bioactives Library	2898-76-2	NA-103	5ng/mL
	ClC1=CC=C(C=C1)C(=O)O	N-benzyl-N-(2-(2-methylpropanoyl)-pyrrolidin-1-yl)pyridin-4-amine hydrochloride	Bepirid HCl	403.0	3	ICCB Known Bioactives Library	74764-40-2	AC-104	5ng/mL
	[H][C@H]1[C@H]2[C@H]3[C@H]4[C@@H](O)C=C(O)C=C4[C@H]5[C@@H](O)C=C(C)C=C5[C@H]6[C@@H](O)C=C(C)C=C6[C@H]7[C@@H](O)C=C(C)C=C7[C@H]8[C@@H](O)C=C(C)C=C8[C@H]9[C@@H](O)C=C(C)C=C9[C@H]10[C@@H](O)C=C(C)C=C10	2R,3S,5Z,8R)-5-(1-hydroxyethylidene)-8,8-dimethyl-7,16-dioxapentacyclo[8.6.1.0 ^{2,7} .0 ^{3,10} .0 ^{4,11} .0 ^{5,12}]tetradeca-1(7),11,12,15(18)-tetraene-4,8-dione	Cyclopiazonic acid	338.4	3	ICCB Known Bioactives Library	18172-33-3	CA-415	5ng/mL
	[O-][N-](=O)(=O)C1=CC=C(C=C1)C(=O)O	1-[[E]-[5-(4-nitrophenyl)furan-2-yl]methylidene]imidazolidin-2,4-dione	Dantrolene	314.3	3	ICCB Known Bioactives Library	7261-97-4	AC-100	5ng/mL
	ClC1=CC=C(C=C1)C(=O)O	7-chloro-3-methyl-2H-1,2,4-benzothiadiazin-1,1-dione	Diazoxide	230.7	3	ICCB Known Bioactives Library	364-98-7	KC-115	5ng/mL
	ClN1C=CC(=O)N1	3,5-diamino-N-[(E)-amino[[2,4-dichlorophenyl]amino]methylidene]-6-chloropyrazine-2-carboxamide hydrochloride	Dichlorobenzamide HCl	425.1	3	ICCB Known Bioactives Library	1166-01-4	CA-204	5ng/mL
	ClC1=CC=C(C=C1)C(=O)O	2S,3S)-5-[[2-(dimethylamino)ethyl]-2-(4-methoxyphenyl)-4-oxo-2,3,4,5-tetrahydro-1,5-benzothiazepin-3-yl]acetate hydrochloride	Diltiazem HCl	451.0	3	ICCB Known Bioactives Library	33288-22-5	CA-205	5ng/mL
	CC1=CC=C(C=C1)C(=O)O	N-(4-{1-[2-(8-methylpyridin-2-yl)ethyl]piperidin-4-carboxyl}phenyl)methanesulfonamide	E-6031	401.5	3	ICCB Known Bioactives Library	113558-89-7	KC-158	5ng/mL
	NC1=CC=C(C=C1)C(=O)O	5-amino-1-[[2,6-dichloro-4-(trifluoromethyl)phenyl]-4-trifluoromethanesulfonyl-1H-pyrazole-3-carbonitrile	Fipronil	437.1	3	ICCB Known Bioactives Library	120068-37-3	DM-102	5ng/mL
	CC(=O)O	N-(6-piperidin-2-ylmethyl)-2,5-bis[2,2,2-trifluoroethyl]benzamide acetate	Flecainide acetate	473.4	3	ICCB Known Bioactives Library	54143-56-5	AC-130	5ng/mL
	CC1=CC=C(C=C1)C(=O)O	2-[E)-(trifluoromethyl)phenyl]amino]benzoic acid	Fluoramic acid	281.2	3	ICCB Known Bioactives Library	530-78-9	AC-123	5ng/mL

Structure	SMILES	BIPAC	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	ClC1C(F)CC(C)C1CN1C(C)CC(NC)C(=O)O1C1=CC=C(C)C1	1-[5-(4-fluorophenylmethyl)-4-(2E)-3-phenylprop-2-en-1-yl]piperazine dihydrochloride	Flunitrazepam HCl	477.4	3	ICCB Known Bioactives Library	30484-77-6	CA-225	5mg/mL
	FC1=CC=C(C)C=C1C(=O)OCC(O)C(=O)N(C)C(=O)C1=CC=C(C)C1=CC=C(C)C1	8-[4-(4-bis(4-fluorophenyl)butyl)-1-phenyl-1,3,8-triazaspiro[4.5]decan-4-one	Flupirtine	475.6	3	ICCB Known Bioactives Library	1841-19-6	AD-116	5mg/mL
	CCC(=O)C1=CC=C(C)C=C1C(=O)N(C)C(=O)C(=O)C1=CC=C(C)C1	methyl 4-(2-benzyloxy-5-methyl-1H-pyridin-3-yl)carboxylate	FPL-64178	347.4	3	ICCB Known Bioactives Library	120934-96-5	CA-245	5mg/mL
	CCCCC(OH)C(=O)C(=O)C(=O)C(=O)C(=O)C(=O)C(=O)C(=O)C	3S)-5-hydroxy-1-(4-hydroxy-3-methoxyphenyl)decan-3-one	6-Gingerol	294.4	3	ICCB Known Bioactives Library	23513-14-6	CA-422	5mg/mL
	CC1=NC(=O)C(N)C(=O)C1=CC=C(C)C(=O)C(=O)N(C)C(=O)C(=O)C	N-(2-{4-[(cyclohexylcarbonylamino)sulfonylphenylethyl]-5-methylpyridin-2-yl}carbamoyl)pyridin-2-carboxamide	Olipide	445.5	3	ICCB Known Bioactives Library	28084-61-9	AD-120	5mg/mL
	CCC1=CC=C(C)C=C1C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C(=O)C(=O)C	5-chloro-N-(2-{4-[(cyclohexylcarbonylamino)sulfonylphenylethyl]-2-methylbenzamide	Oltipride	494.0	3	ICCB Known Bioactives Library	10230-21-6	KC-120	5mg/mL
	[H](O)C1[C@@H](O)C[C@H](O)C[C@@H](O)C[C@H](O)C[C@H](O)C1O	(1S,3R,4R,5S,8R,10R,16R)-5,5,14-tetramethyltetrahydro[11.2.1.0.0'.0'']hexadecane-3,4,6,8,14,16-hexol	Grayanotoxin II	370.5	3	ICCB Known Bioactives Library	4678-45-9	NA-135	5mg/mL
	C1=CC=C(C=C1)C(=O)O	2-(E)-6,7-dichloro-2-cyclopentyl-2-methyl-1-oxo-2,3-dihydro-1H-inden-5-yl]oxy]acetic acid	R(-)-9AA-94	357.2	3	ICCB Known Bioactives Library	54193-31-8	CM-101	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C(=O)C	2R,3R)-5-[2-(dimethylamino)ethyl]-2-(4-methoxyphenyl)-4-oxo-2,3,4,5-tetrahydro-1,5-benzothiazepin-3-yl acetate hydrochloride	L-cis-Diltiazem HCl	451.0	3	ICCB Known Bioactives Library	42399-54-2	CA-207	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C(=O)C	2-(diethylamino)-N-(2,6-dimethylphenyl)acetamide hydrate hydrochloride	Lidocaine HCl	288.8	3	ICCB Known Bioactives Library	6108-05-0	AD-128	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C(=O)C	4-[4-(4-chlorophenyl)-4-hydroxypiperidin-1-yl]-N,N-dimethyl-2,2-diphenylbutanamide hydrochloride	Looperamide HCl	513.5	3	ICCB Known Bioactives Library	34552-83-5	ALX-550-253	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C(=O)C	5-(2-(3,4-dimethoxyphenylethyl)(methylamino)-2-(propan-2-yl)-2-(3,4,5-trimethoxyphenyl)pentanenitrile hydrochloride	Methoxypropimide HCl	521.1	3	ICCB Known Bioactives Library	18662-47-8	AD-103	5mg/mL
	NC1=NC(=O)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C	2,6-diamino-4-(piperidin-1-yl)pyrimidin-1-ium-1-olate	Minoxidil	209.2	3	ICCB Known Bioactives Library	38304-91-5	KC-125	5mg/mL
	NC1=NC(=O)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C	2,6-diamino-4-(piperidin-1-yl)-1-sulfonatoxy)pyrimidin-1-ium	Minoxidil sulfate	289.3	3	ICCB Known Bioactives Library	83701-22-6	KC-130	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C	3-[2-(benzyl(methylamino)ethyl) 5-methyl 2,6-dimethyl-4-(3-nitrophenyl)-1,4-dihydropyridin-3-yl]dicarboxylate hydrochloride	Nicardipine HCl	516.0	3	ICCB Known Bioactives Library	54527-84-3	AD-108	5mg/mL
	CC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C	5-(acetyloxy)-2,6-dimethyl-4-(2-nitrophenyl)-1,4-dihydropyridin-3-yl acetate	Nifedipine	348.5	3	ICCB Known Bioactives Library	21829-29-4	CA-210	5mg/mL
	CC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C	2-[(3-(trifluoromethyl)phenyl)amino]pyridine-3-carboxylic acid	Nitfunic acid	282.2	3	ICCB Known Bioactives Library	4394-00-7	AD-134	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C(=O)C	3-[2-(4,4-dihydroquinolindin-1-yl)propyl] 5-methyl 2,6-dimethyl-4-(3-nitrophenyl)-1,4-dihydropyridin-3-yl dicarboxylate hydrochloride	Nigalipine	648.2	3	ICCB Known Bioactives Library	113317-61-6	CA-218	5mg/mL
	CCCC(=O)C1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	3-(2-methoxyethyl) 5-propan-2-yl 2,6-dimethyl-4-(3-nitrophenyl)-1,4-dihydropyridin-3-yl dicarboxylate	Nimodipine	418.4	3	ICCB Known Bioactives Library	66085-59-4	ALX-550-277	5mg/mL
	CC(=O)C1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	3-ethyl 5-methyl 2,6-dimethyl-4-(3-nitrophenyl)-1,4-dihydropyridin-3-yl dicarboxylate	Nitrendipine	380.4	3	ICCB Known Bioactives Library	39562-70-4	ALX-550-278	5mg/mL
	CC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	2-phenylamino]benzoic acid	N-Phenylanthranic acid	212.2	3	ICCB Known Bioactives Library	91-40-7	AD-132	5mg/mL
	CC(=O)C1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	5-nitro-2-[(3-phenylpropyl)amino]benzoic acid	NPPB	300.3	3	ICCB Known Bioactives Library	107254-88-4	CM-100	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	1-[2-hydroxy-5-(trifluoromethyl)phenyl]-5-(trifluoromethyl)-2,3-dihydro-1H-1,3-benzodiazol-2-one	NS-1619	382.2	3	ICCB Known Bioactives Library	153581-01-0	KC-127	5mg/mL
	[H](O)C1[C@@H](O)C[C@H](O)C[C@@H](O)C[C@H](O)C1O	(1S,2R,5S,7R,11S,14S)-11-hydroxy-7-(2-hydroxypropyl)-2-(1,2-dimethyl-6-oxo-2,3-azabicyclo[12.10.0.0']{2.0'.0''.0'''}tetracosane-9,16,24,17,18,21-pentane-8-one	Pavline	435.6	3	ICCB Known Bioactives Library	57186-25-1	KC-155	5mg/mL
	CC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	3S,4R)-3-hydroxy-2,2-dimethyl-4-[(3-oxocyclopent-1-en-1-yl)oxy]-2,4-dihydro-2H-1-benzopyran-6-carbonitrile	PCD-400	299.3	3	ICCB Known Bioactives Library	121055-10-5	KC-131	5mg/mL
	[H](O)C1[C@@H](O)C[C@H](O)C[C@@H](O)C[C@H](O)C1O	(1S,2R,5S,8S,8R,9S,10R,12S,15R,16S,25R,27S,28R)-21-chloro-15,16,33,33-tetramethyl-24-methyldene-10-(prop-1-en-2-yl)-7,11,32-trioxo-18-azadecacyclo[25.4.2.0'.0''.0''.0''.0''.0''.0''.0''.0''.0''.0''.0''.0''']tetracosane-9,16,24,17,18,21-pentane-8-one	Pentam A	634.2	3	ICCB Known Bioactives Library	12627-35-9	KC-157	5mg/mL
	NC1=NC(=O)C(=O)N(C)C(=O)C(=O)C(=O)C	3,5-diamino-6-chloro-N-(N-phenylcarbamimidyl)pyrazine-2-carboxamide	Phenamil	305.7	3	ICCB Known Bioactives Library	2038-35-9	AD-131	5mg/mL
	CC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	3-[(4S)-dihydro-1H-indazol-2-ylmethyl]4-methylphenylamino]phenol	Phentolamine	281.4	3	ICCB Known Bioactives Library	73-05-2	AD-121	5mg/mL
	O=C1NC(=O)N1C(=O)C(=O)C(=O)C	5,5-diphenylmizalidine-2,4-dione	Phenytoin	252.3	3	ICCB Known Bioactives Library	57-41-0	AD-125	5mg/mL
	FC1=CC=C(C)C=C1C(=O)OCC(O)C(=O)N(C)C(=O)C1=CC=C(C)C1=CC=C(C)C1	1-[1-(4,4-bis(4-fluorophenyl)butyl)piperidin-4-yl]-2,3-dihydro-1H-1,3-benzodiazol-2-one	Pinoclide	461.5	3	ICCB Known Bioactives Library	2062-78-4	AD-109	5mg/mL
	CC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	E)-2-cyano-1-(3,3-dimethylbutan-2-yl)-3-(pyridin-4-yl)guanidine	Pivacil	245.3	3	ICCB Known Bioactives Library	85371-64-8	AD-118	5mg/mL
	CC(=O)OCC(=O)N(C)C(=O)C(=O)C(=O)C	4-amino-N-(2-(diethylamino)ethyl)benzamide	Procainamide	235.3	3	ICCB Known Bioactives Library	614-39-1	AD-127	5mg/mL
	CC(=O)OCC(=O)N(C)C(=O)C(=O)C(=O)C	1-[2-(2-hydroxy-3-(propylamino)propoxy)phenyl]-3-phenylpropan-1-one	Propafenone	341.4	3	ICCB Known Bioactives Library	54063-53-5	AD-124	5mg/mL
	ClC1=CC=C(C=C1)C(=O)N(C)C(=O)C(=O)C(=O)C	[(1R,2R)-5-ethenyl-1-azabicyclo[2.2.2]octan-2-yl](8-methoxyquinolin-4-yl)methanol hydrate hydrochloride	Quinidine HCl	378.9	3	ICCB Known Bioactives Library	6151-40-2	AD-129	5mg/mL

Structure	SMILES	BPAD	Name	MW	Plate number	Plate description	CAS	Catalog number	Concentration
	<chem>CC1=CC=C(C=C1)C=CC(=O)N#CC(=O)C=CC(=O)C=CC(=O)C1=CC=C(C=C1)O</chem>	(E)-3-phenylprop-2-en-1-yl (Z)-2-cyano-3-(3,4-dihydroxyphenyl)prop-2-enoate	CDG	321.5	6	ICCB Known Bioactives Library	132465-11-3	EI-211	5mg/mL
	<chem>[H][C@H](O)[C@H](O)C(=O)N[C@@H]1OC[C@H]1C=CC=CC=C</chem>	(R)-3-[(4E,7E)-nona-4,7-dienyl]oxirane-2-carboxamide	Cerulenin	223.3	6	ICCB Known Bioactives Library	17397-89-6	G-237	5mg/mL
	<chem>CCOC1=CC=C(C=C1)N(CCC1=CC=C(C=C1)Cl)CC2=CC=C(C=C2)</chem>	benzyl(2-chloroethyl)(1-phenoxyprop-2-yl)amine	Phenylethylamine	303.8	6	ICCB Known Bioactives Library	63-82-3	CA-305	5mg/mL
	<chem>CN(C)COC1=NC2=CC=CC=C2N1C3=CC=CC=C3Br</chem>	1-(3-bromo-9H-carbazol-9-yl)-3-(dimethylamino)propan-2-ol	Wakostatin	426.1	6	ICCB Known Bioactives Library	253449-04-6	T-126	5mg/mL

Compound description	Solvent	Plate port number
Bioactive lipid: Activates PKC zeta	DMSO	2840
Bioactive lipid: Ceramide analog. Apoptosis inducer	DMSO	2840
CNS receptor ligand: Cannabinoid CB1 receptor agonist	DMSO	2840
Nuclear receptor ligand: Vitamin D receptor agonist	DMSO	2840
Bioactive lipid: Activates PKC	DMSO	2840
Bioactive lipid: Activates PKC	DMSO	2840
Bioactive lipid: Activates PKC	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Fatty acid hydroperoxide	DMSO	2840
Bioactive lipid: Myristic acid analog	DMSO	2840
Bioactive lipid: Bioactive linoleic acid metabolite	DMSO	2840
Bioactive lipid: Fatty acid hydroperoxide	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Nuclear receptor ligand: Retinoid receptor ligand	DMSO	2840
Bioactive lipid: Bioactive linoleic acid metabolite	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Fatty acid hydroperoxide	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Prostaglandin EP receptor agonist	DMSO	2840
Lipid biosynthesis: Inhibits fatty acid omega oxidation	DMSO	2840
Bioactive lipid: Prostaglandin EPI receptor agonist	DMSO	2840
Bioactive lipid: PAF agonist	DMSO	2840
Bioactive lipid: DAG analog	DMSO	2840
Bioactive lipid: PAF receptor agonist	DMSO	2840
Bioactive lipid: Blocks DAG activation of PKC	DMSO	2840
Bioactive lipid: Blocks DAG activation of PKC	DMSO	2840
Bioactive lipid: Inhibits PI-specific PLC	DMSO	2840
Bioactive lipid: PKC activator	DMSO	2840
Bioactive lipid: PKC activator	DMSO	2840
Bioactive lipid: PKC activator	DMSO	2840

Compound description	Solvent	Plate port number
Nuclear receptor ligands: Vitamin D receptor ligand	DMSO	2840
Nuclear receptor ligands: Vitamin D receptor ligand	DMSO	2840
Endocannabinoids: Cannabinoid CB1 receptor agonist	DMSO	2840
Bioactive lipids: Protein palmitoylation inhibitor	DMSO	2840
Bioactive lipids: Protein myristoylation inhibitor	DMSO	2840
Nuclear receptor ligands: Retinoid receptor agonist / apoptosis inducer	DMSO	2840
Bioactive lipids: Myristic acid analog	DMSO	2840
Bioactive lipids: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipids: Fatty acid hydroperoxide	DMSO	2840
Bioactive lipids: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipids: Cyclooxygenase & lipoxygenase inhibitor	DMSO	2840
Bioactive lipids: Lipoxygenase inhibitor	DMSO	2840
Bioactive lipids: 5-KETE receptor (RS27) agonist	DMSO	2840
Bioactive lipids: AHR agonist	DMSO	2840
Bioactive lipids: Bioactive prostaglandin	DMSO	2840
Bioactive lipids: PLA2 inhibitor	DMSO	2840
Bioactive lipids: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipids: Thromboxane TP receptor agonist	DMSO	2840
Bioactive lipids: Bioactive linoleic acid metabolite	DMSO	2840
Bioactive lipids: Fatty acid hydroperoxide	DMSO	2840
Endocannabinoids: Endogenous sleep inducing lipid	DMSO	2840
Bioactive lipids: Bioactive prostaglandin	DMSO	2840
Bioactive lipids: Bioactive prostaglandin	DMSO	2840
Nuclear receptor ligands: Retinoid RXR agonist	DMSO	2840
Inhibitor: Phyl-4-hydroxylase inhibitor	DMSO	2840
Bioactive lipids: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipids: Negative control for AGGC and AFC	DMSO	2840
Bioactive lipids: ICMT inhibitor	DMSO	2840
CNS receptor ligands: Cannabinoid CB1 receptor antagonist	DMSO	2840
Nuclear receptor ligands: Retinoid RAR agonist	DMSO	2840
Endocannabinoids: Cannabinoid receptor agonist	DMSO	2840
Endocannabinoids: Cannabinoid receptor agonist	DMSO	2840
Endocannabinoids: Cannabinoid receptor agonist	DMSO	2840
Endocannabinoids: Cannabinoid receptor agonist	DMSO	2840

Compound description	Solvent	Plate port number
Endocannabinoids: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: PAF precursor	DMSO	2840
Bioactive lipid: Cannabinoid CB1 inverse agonist	DMSO	2840
Bioactive lipid: Apoptosis inducer	DMSO	2840
Bioactive lipid: Negative control for C2 ceramide	DMSO	2840
Bioactive lipid: Stimulates Cer-activated PK	DMSO	2840
Bioactive lipid: Negative control for C8 ceramide	DMSO	2840
Bioactive lipid: PPAR delta agonist	DMSO	2840
Nuclear receptor ligand: PPAR gamma agonist	DMSO	2840
Nuclear receptor ligand: PPAR alpha agonist	DMSO	2840
Bioactive lipid: Prostaglandin FP receptor agonist	DMSO	2840
Bioactive lipid: PAF receptor agonist	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Apoptosis inducer	DMSO	2840
Nuclear receptor ligand: AHR agonist	DMSO	2840
Lipid biosynthesis: Sphingosine kinase inhibitor	DMSO	2840
Lipid biosynthesis: Glucosylceramide synthase inhibitor	DMSO	2840
Lipid biosynthesis: Glucosylceramide synthase inhibitor	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Negative control for PAF	DMSO	2840
Bioactive lipid: Carbonylmethylation inhibitor	DMSO	2840
Bioactive lipid: Prostaglandin FP receptor agonist	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Bioactive linoleic acid metabolite	DMSO	2840
Bioactive lipid: Bioactive linoleic acid metabolite	DMSO	2840
Bioactive lipid: Leukotriene B4 receptor agonist	DMSO	2840

Compound description	Solvent	Plate port number
Bioactive lipid: CysLT receptor agonist	DMSO	2840
Bioactive lipid: CysLT receptor agonist	DMSO	2840
Bioactive lipid: CysLT1 receptor agonist	DMSO	2840
Endocannabinoids: Bioactive linoleic acid metabolite	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Bioactive lipid: LPA agonist / antagonist	DMSO	2840
Bioactive lipid: Leukotriene D4 receptor antagonist	DMSO	2840
Bioactive lipid: Inactive PAF metabolite	DMSO	2840
Bioactive lipid: LPA receptor antagonist	DMSO	2840
Lipid biosynthesis: Ceramidase inhibitor	DMSO	2840
Lipid biosynthesis: Negative control for D-erythro-MAPP	DMSO	2840
Bioactive lipid: Polyunsaturated fatty acid	DMSO	2840
Endocannabinoids: Cannabinoid receptor agonist	DMSO	2840
Nuclear receptor ligand: Retinoid RXR agonist	DMSO	2840
Bioactive lipid: Prostaglandin EP receptor agonist	DMSO	2840
Lipid biosynthesis: Sphingosine kinase inhibitor	DMSO	2840
Bioactive lipid: Bioactive arachidonic acid metabolite	DMSO	2840
Endocannabinoids: FAAH inhibitor	DMSO	2840
Endocannabinoids: FAAH inhibitor	DMSO	2840
Bioactive lipid: PAF receptor agonist	DMSO	2840
Bioactive lipid: PAF receptor agonist	DMSO	2840
Bioactive lipid: PAF receptor agonist	DMSO	2840
Endocannabinoids: Cannabinoid CB2 receptor agonist	DMSO	2840
Bioactive lipid: Activates MAP kinase cascade	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Bioactive prostaglandin	DMSO	2840
Bioactive lipid: Prostaglandin DP receptor agonist	DMSO	2840
Bioactive lipid: Prostaglandin EP receptor agonist	DMSO	2840
Bioactive lipid: Prostaglandin EP receptor agonist	DMSO	2840
Bioactive lipid: Prostaglandin FP receptor agonist	DMSO	2840

Compound description	Solvent	Plate port number
Bioactive lipids: Prostaglandin FP receptor agonist	DMSO	2840
Bioactive lipids: Prostaglandin IP receptor agonist	DMSO	2840
Bioactive lipids: Bioactive prostaglandin	DMSO	2840
Nuclear receptor ligands: Retinoid RAR agonist	DMSO	2840
Lipid biosynthesis: 5-Lipoxygenase inhibitor	DMSO	2840
Bioactive lipids: MOR ATPase activator	DMSO	2840
Bioactive lipids: PKC inhibitor	DMSO	2840
Bioactive lipids: Thromboxane A2 antagonist	DMSO	2840
		2840
Nuclear receptor ligands: Retinoid RAR agonist	DMSO	2840
Bioactive lipids: Thromboxane TP receptor agonist	DMSO	2840
Bioactive lipids: Leukotriene B4 receptor antagonist	DMSO	2840
CNS receptor ligands: Cannabinoid CB1/CB2 receptor agonist	DMSO	2840
Nuclear receptor ligands: PPAR alpha agonist	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Misc. channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium ionophore	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Intracellular calcium	DMSO	2840
Ion channel ligands: Intracellular calcium	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Misc. channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840

Compound description	Solvent	Plate port number
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Intracellular calcium	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Misc. channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Misc. channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Misc. channels	DMSO	2840
Ion channel ligands: Misc. channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840

Compound description	Solvent	Plate port number
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Intracellular calcium	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Intracellular calcium	DMSO	2840
Ion channel ligands: Intracellular calcium	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Sodium channels	DMSO	2840
Ion channel ligands: Calcium channels	DMSO	2840
Ion channel ligands: Potassium channels	DMSO	2840
Protease inhibitors: Serine protease inhibitor	DMSO	2840
Protease inhibitors: Neutral endopeptidase inhibitor	DMSO	2840
Protease inhibitors: Prolyl endopeptidase inhibitor	DMSO	2840
Protease inhibitors: Cathepsin L inhibitor	DMSO	2840
Protease inhibitors: Granzyme B inhibitor	DMSO	2840
Protease inhibitors: Gamma secretase inhibitor	DMSO	2840
Protease inhibitors: Aminopeptidase inhibitor	DMSO	2840
Protease inhibitors: Tripeptidyl peptidase II inhibitor	DMSO	2840
Protease inhibitors: Calpain inhibitor	DMSO	2840
Protease inhibitors: Calpain inhibitor	DMSO	2840
Inhibitors: Bcl-2/Bcl-XL ligand induces apoptosis	DMSO	2840
SK2 inh.	DMSO	2840
Inhibitors: inhibits mitochondrial permeability pore opening	DMSO	2840
Inhibitors: induces mitochondrial permeability pore opening	DMSO	2840
		2840
Inhibitors: iNOS inhibitor	DMSO	2840
Lipid biosynthesis: inhibits ceramide synthase	DMSO	2840
Inhibitors: inhibits NF-kappaB transcriptional activation	DMSO	2840

Compound description	Solvent	Plate port number
Inhibitors: AP-1 inhibitor	DMSO	2840
Inhibitors: induces intracellular acidification	DMSO	2840
Inhibitors: inhibits TNFalpha production	DMSO	2840
Inhibitors: topoisomerase I inhibitor	DMSO	2840
Inhibitors: topoisomerase I inhibitor	DMSO	2840
Inhibitors: IkappaB kinase inhibitor	DMSO	2840
Protease inhibitors: Proteasome inhibitor	DMSO	2840
Inhibitors: NFkappaB inhibitor	DMSO	2840
Nuclear receptor ligands: PPAR alpha agonist	DMSO	2840
Inhibitors: FXBP12 inhibitor	DMSO	2840
Inhibitors: binds to FXBP inhibits calcineurin	DMSO	2840
Inhibitors: superoxide/free-radical inhibitor	DMSO	2840
Inhibitors: glutathione peroxidase mimetic	DMSO	2840
Inhibitors: iNOS/NOS inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE1/2) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE1) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE1) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE3) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE3) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE4) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE4) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE5) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE5) inhibitor	DMSO	2840
Kinase inhibitors: kinase inhibitor	DMSO	2840
Inhibitors: farnesylation inhibitor	DMSO	2840
Inhibitors: transcription inhibitor	DMSO	2840
Ion channel ligands: Vanilloid receptor agonist	DMSO	2840
Kinase inhibitors: Tyrosine kinase inhibitor. Broad spectrum	DMSO	2840
Kinase inhibitors: EGF-R tyrosine kinase inhibitor	DMSO	2840
Kinase inhibitors: PDGF receptor kinase inhibitor	DMSO	2840
Kinase inhibitors: JAK2 inhibitor	DMSO	2840
Kinase inhibitors: NGF receptor inhibitor	DMSO	2840
Inhibitors: monovalent cation ionophore	DMSO	2840
Kinase inhibitors: c-kit, FGF and PDGF kinase inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE3) inhibitor	DMSO	2840

Compound description	Solvent	Plate port number
Inhibitors: p53 reactivator	DMSO	2840
Inhibitors: inhibits clathrin coated pit mediated endocytosis	DMSO	2840
Inhibitors: PARP inhibitor	DMSO	2840
Inhibitors: ADP-ribose polymerase, apoptosis inhibitor	DMSO	2840
Inhibitors: HSP-90 inhibitor	DMSO	2840
Inhibitors: MAP kinase activator	DMSO	2840
Inhibitors: DNA polymerase inhibitor	DMSO	2840
Lipid biosynthesis: phospholipase A2 inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE1) inhibitor	DMSO	2840
Inhibitors: phosphodiesterase (PDE3) inhibitor	DMSO	2840
Inhibitors: farnesyltransferase inhibitor	DMSO	2840
Nuclear receptor ligands: PPAR gamma antagonist	DMSO	2840
Inhibitors: vacuolar ATPase inhibitor	DMSO	2840
Inhibitors: cell permeable Ca ⁺⁺ chelator	DMSO	2840
Inhibitors: Myosin II inhibitor	DMSO	2840
Inhibitors: ANT inhibitor	DMSO	2840
CNS receptor ligands: Adenosine receptor agonist	DMSO	2840
Inhibitors: ARF GEF inhibitor	DMSO	2840
Inhibitors: NO synthase inhibitor	DMSO	2840
Activators: PKA activator	DMSO	2840
Activators: PKG activator	DMSO	2840
Inhibitors: Na ⁺ K ⁺ Cl ⁻ cotransport inhibitor	DMSO	2840
CNS receptor ligands: adenosine receptor agonist	DMSO	2840
Lipid biosynthesis: 5 lipoygenase inhibitor	DMSO	2840
Protease inhibitors: Cathepsin B inhibitor	DMSO	2840
Kinase inhibitors: GSK3 beta inhibitor	DMSO	2840
Protease inhibitors: Calpain inhibitor	DMSO	2840
Kinase inhibitors: PKC inhibitor	DMSO	2840
Inhibitors: PPI, PP2A inhibitor	DMSO	2840
Inhibitors: Topoisomerase I inhibitor	DMSO	2840
Inhibitors: PP2A inhibitor	DMSO	2840
Ion channel ligands: vanilloid receptor agonist	DMSO	2840
Ion channel ligands: vanilloid receptor antagonist	DMSO	2840
Inhibitors: glucosidase inhibitor	DMSO	2840

Compound description	Solvent	Plate port number
Inhibitors: cGMP phosphodiesterase inhibitor	DMSO	2840
Inhibitors: inhibitor of mitochondrial Na ⁺ /Ca ²⁺ exchange	DMSO	2840
Kinase inhibitors: PKC inhibitor	DMSO	2840
Inhibitors: Antioxidant/ NFkappa B inhibitor	DMSO	2840
CNS receptor ligands: adrenoceptor agonist (beta)	DMSO	2840
Inhibitors: PTP1B inhibitor	DMSO	2840
CNS receptor ligands: adrenoceptor agonist (alpha)	DMSO	2840
CNS receptor ligands: adrenoceptor agonist (alpha)	DMSO	2840
CNS receptor ligands: Dopamine antagonist	DMSO	2840
Inhibitors: NFkappaB inhibitor	DMSO	2840
Inhibitors: protein synthesis inhibitor	DMSO	2840
Inhibitors: calcineurin inhibitor	DMSO	2840
Inhibitors: calcineurin inhibitor	DMSO	2840
Inhibitors: F actin capper	DMSO	2840
Inhibitors: F actin capper	DMSO	2840
Lipid biosynthesis: PC-PLC inhibitor	DMSO	2840
Kinase inhibitors: p58ckc inhibitor	DMSO	2840
Activators: GC stimulator / HIF-1 alpha inhibitor	DMSO	2840
Inhibitors: lowers GTP levels	DMSO	2840
Inhibitors: mannosidase inhibitor	DMSO	2840
Inhibitors: glucosidase inhibitor	DMSO	2840
Nuclear receptor ligands: corticosteroid	DMSO	2840
Activators: PKA activator	DMSO	2840
Activators: PKA activator	DMSO	2840
Nuclear receptor ligands: PPAR gamma antagonist	DMSO	2840
CNS receptor ligands: Dopamine agonist	DMSO	2840
Inhibitors: Iron chelator	DMSO	2840
Inhibitors: flavoprotein inhibitor	DMSO	2840
Ion channel ligands: ER Ca ²⁺ ATPase inhibitor	DMSO	2840
Inhibitors: topoisomerase II inhibitor; induces apoptosis	DMSO	2840
Kinase inhibitors: CK2 inhibitor	DMSO	2840
Inhibitors: SOD mimetic	DMSO	2840
Inhibitors: calmodulin inhibitor	DMSO	2840
Protease inhibitors: calpain/cathepsin inhibitor	DMSO	2840

Compound description	Solvent	Plate port number
Inhibitors: Phosphodiesterase (PDE2) inhibitor/adenosine deaminase inhibitor	DMSO	2840
Nuclear receptor ligands: PXR/SXR agonist	DMSO	2840
Nuclear receptor ligands: estrogen	DMSO	2840
Inhibitors: topoisomerase II inhibitor	DMSO	2840
Inhibitors: mitochondrial uncoupler	DMSO	2840
Activators: Adenylate cyclase activator	DMSO	2840
Inhibitors: HSP90 inhibitor	DMSO	2840
Kinase inhibitors: Tyrosine kinase inhibitor	DMSO	2840
Kinase inhibitors: PKC inhibitor	DMSO	2840
Protease inhibitors: broad spectrum MMP inhibitor	DMSO	2840
Kinase inhibitors: PKG inhibitor	DMSO	2840
Lipid biosynthesis: Cox 2 inhibitor	DMSO	2840
Kinase inhibitors: kinase inhibitor	DMSO	2840
Kinase inhibitors: PKA inhibitor	DMSO	2840
Kinase inhibitors: kinase inhibitor	DMSO	2840
Kinase inhibitors: inhibitor of Rho-dependent kinases	DMSO	2840
Kinase inhibitors: kinase inhibitor	DMSO	2840
Kinase inhibitors: PKC inhibitor	DMSO	2840
GHS receptor ligands: Histamine receptor agonist	DMSO	2840
Kinase inhibitors: Insulin receptor TK inhibitor	DMSO	2840
Inhibitors: DNA minor groove binder	DMSO	2840
Inhibitors: acetylcholinesterase inhibitor	DMSO	2840
Nuclear receptor ligands: LXR agonist	DMSO	2840
Inhibitors: topo II inhibitor that does not cause DNA breaks	DMSO	2840
Lipid biosynthesis: cyclooxygenase inhibitor	DMSO	2840
Ion channel ligands: Ca ²⁺ ionophore	DMSO	2840
Inhibitors: PDE inhibitor (broad spec), adenosineR agonist	DMSO	2840
Inhibitors: CDC25 phosphatase inhibitor	DMSO	2840
Inhibitors: PDI inhibitor	DMSO	2840
Kinase inhibitors: Kinase inhibitor (Broad spectrum)	DMSO	2840
Kinase inhibitors: CaM kinase II inhibitor	DMSO	2840
Kinase inhibitors: PKA inhibitor	DMSO	2840
Inhibitors: Ras farnesyltransferase inhibitor	DMSO	2840
Activators: NO donor	DMSO	2840

Compound description	Solvent	Plate port number
Inhibitors: Actin inhibitor	DMSO	2840
Kinase inhibitors: Tyrosine kinase inhibitor EGF-R	DMSO	2840
Kinase inhibitors: CDC-2 kinase inhibitor	DMSO	2840
Protease inhibitors: protease inhibitor	DMSO	2840
Inhibitors: Inhibits SK kinase activation	DMSO	2840
Inhibitors: Calcineurin inhibitor	DMSO	2840
Inhibitors: NO synthesis inhibitor	DMSO	2840
Lipid biosynthesis: PI-3-Kinase inhibitor	DMSO	2840
Inhibitors: Inhibits NO-activation of guanylate cyclase	DMSO	2840
Kinase inhibitors: ERK-2 inhibitor	DMSO	2840
Lipid biosynthesis: Phospholipase A2 inhibitor	DMSO	2840
Inhibitors: ras farnesylation inhibitor	DMSO	2840
Activators: activates heterotrimeric GTPases	DMSO	2840
Inhibitors: antioxidant, cytoprotectant	DMSO	2840
Protease inhibitors: Inhibitor of 20S-proteasome chymotrypsin activity	DMSO	2840
Kinase inhibitors: BTK inhibitor	DMSO	2840
CNS receptor ligands: D1FR inhibitor	DMSO	2840
CNS receptor ligands: nicotinic cholinergic agonist	DMSO	2840
Lipid biosynthesis: Inhibitor HMG-CoA reductase	DMSO	2840
Inhibitors: cross links DNA	DMSO	2840
Kinase inhibitors: kinase inhibitor	DMSO	2840
Kinase inhibitors: kinase inhibitor	DMSO	2840
Inhibitors: E2F inhibitor	DMSO	2840
Kinase inhibitors: cRAF1 kinase inhibitor	DMSO	2840
Ion channel ligands: Na ⁺ ionophore	DMSO	2840
Inhibitors: FKBP ligand	DMSO	2840
Inhibitors: tubulin inhibitor	DMSO	2840
Inhibitors: phospholipase A2 inhibitor	DMSO	2840
Inhibitors: PPI/PP2A inhibitor	DMSO	2840
Inhibitors: Fo ATP synthase inhibitor	DMSO	2840
Kinase inhibitors: CDK inhibitor	DMSO	2840
Inhibitors: Na ⁺ /K ⁺ ATPase inhibitor	DMSO	2840
Bioactive lipids: PAF antagonist	DMSO	2840
Kinase inhibitors: MEK inhibitor	DMSO	2840

Compound description	Solvent	Plate port number
Kinase inhibitors: VEGF-R (Flk-1) tyrosine kinase	DMSO	2840
Protease inhibitors: protease inhibitor	DMSO	2840
Inhibitors: PAMP inhibitor	DMSO	2840
Activators: PKC activator	DMSO	2840
Inhibitors: p53 inhibitor	DMSO	2840
Nuclear receptor ligands: Corticosteroid receptor (CAR) agonist	DMSO	2840
Lipid biosynthesis: COX1 inhibitor	DMSO	2840
Inhibitors: Hedgehog pathway inhibitor	DMSO	2840
Protease inhibitors: src family tyrosine kinase inhibitor	DMSO	2840
CNS receptor ligands: adrenoceptor agonist	DMSO	2840
Inhibitors: DNA intercalator	DMSO	2840
CNS receptor ligands: adrenoceptor antagonist (beta)	DMSO	2840
Inhibitors: protein synthesis inhibitor	DMSO	2840
Kinase inhibitors: ALK4, ALK5, ALK7 inhibitor	DMSO	2840
Kinase inhibitors: JNK inhibitor	DMSO	2840
Kinase inhibitors: kinase inhibitor (plus other)	DMSO	2840
Inhibitors: FRAP inhibitor	DMSO	2840
Inhibitors: integrin inhibitor	DMSO	2840
Lipid biosynthesis: DAG lipase inhibitor	DMSO	2840
Inhibitors: VHR phosphatase inhibitor	DMSO	2840
Kinase inhibitors: CDK inhibitor	DMSO	2840
Kinase inhibitors: PKC delta inhibitor	DMSO	2840
Inhibitors: CD45 phosphatase inhibitor	DMSO	2840
Kinase inhibitors: Suppressor of MAPKAP kinase-2	DMSO	2840
Kinase inhibitors: MAP kinase inhibitor	DMSO	2840
Kinase inhibitors: GSK-3beta inhibitor	DMSO	2840
Kinase inhibitors: GSK-3beta inhibitor	DMSO	2840
CNS receptor ligands: serotonin receptor agonist	DMSO	2840
Activators: SIRT1 activator	DMSO	2840
Inhibitors: sir2p inhibitor	DMSO	2840
Kinase inhibitors: Src family tyrosine kinase inhibitor	DMSO	2840
Kinase inhibitors: Syk inhibitor	DMSO	2840
Inhibitors: adenylyate cyclase inhibitor	DMSO	2840
Inhibitors: aldose reductase inhibitor	DMSO	2840

Compound description	Solvent	Plate port number
Kinase inhibitors: kinase inhibitor - non-selective	DMSO	2840
Inhibitors: protein glycosylation inhibitor	DMSO	2840
Nuclear receptor ligands: estrogen antagonist	DMSO	2840
Inhibitors: microtubule stabilizer	DMSO	2840
Kinase inhibitors: EGF-R tyrosine kinase inhibitor	DMSO	2840
Inhibitors: TNFalpha synthesis inhibitor	DMSO	2840
Inhibitors: MDC inhibitor	DMSO	2840
Kinase inhibitors: PKC inhibitor	DMSO	2840
Kinase inhibitors: HER-1,2 tyrosine kinase inhibitor	DMSO	2840
Ion channel ligands: voltage-dependent Na channel inhibitor	DMSO	2840
Inhibitors: cell permeable heavy metal chelator	DMSO	2840
Inhibitors: histone deacetylase inhibitor	DMSO	2840
Inhibitors: calmodulin inhibitor-possibly only at high concentrations!	DMSO	2840
Inhibitors: glycosylation inhibitor	DMSO	2840
Inhibitors: Serine protease inhibitor	DMSO	2840
Kinase inhibitors: tyrosine kinase inhibitor	DMSO	2840
Inhibitors: Calcineurin inhibitor	DMSO	2840
Kinase inhibitors: PDGF-R tyrosine kinase inhibitor	DMSO	2840
Lipid biosynthesis: PLC inhibitor	DMSO	2840
Kinase inhibitors: MEK inhibitor	DMSO	2840
Ion channel ligands: K ⁺ ionophore	DMSO	2840
Inhibitors: tubulin inhibitor	DMSO	2840
Inhibitors: calmodulin antagonist	DMSO	2840
Lipid biosynthesis: PI-3K/kinase, other kinases inhibitor	DMSO	2840
Kinase inhibitors: ROCK inhibitor	DMSO	2840
OR5 receptor ligands: Adrenoreceptor antagonist (alpha)	DMSO	2840
Protease inhibitors: Proteasome inhibitor	DMSO	2840
Kinase inhibitors: Raf inhibitor	DMSO	2840
Protease inhibitors: Caspase inhibitor (broad spectrum)	DMSO	2840
Lipid biosynthesis: Cox-2 inhibitor	DMSO	2840
Lipid biosynthesis: 5-lipoxygenase inhibitor	DMSO	2840
Inhibitors: Inosine-5'-monophosphate dehydrogenase inhibitor	DMSO	2840
Inhibitors: Apoptosis inducer, p53 dependent	DMSO	2840
Inhibitors: IP3 receptor blocker	DMSO	2840

Compound description	Solvent	Plate part number
Lipid biosynthesis: 12-Lipoxygenase inhibitor	DMSO	2840
Lipid biosynthesis: Fatty acid biosynthesis inhibitor	DMSO	2840
Inhibitors: Calmodulin antagonist	DMSO	2840
Inhibitors: N-WASP inhibitor	DMSO	2840