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Alzheimer's disease (AD) is a devastating neurodegenerative disorder with a relentless progression. AD pathogenesis is believed to be triggered by the accumulation of the amyloid- β peptide ($A\beta$), which is due to overproduction of $A\beta$ and/or the failure of clearance mechanisms. Several approaches aimed at inhibiting disease progression have advanced to clinical trials for this unmet medical disorder of neurology.

Among these, strategies targeting the production and clearance of the amyloid- β peptide — a cardinal feature of Alzheimer's disease that is thought to be important in disease pathogenesis — are the most advanced. Approaches aimed at modulating the abnormal aggregation of tau filaments (another key feature of the disease), and those targeting metabolic dysfunction, are also being evaluated preclinical and clinically. We, at Toronto Research Chemicals offer a wide product range for designing different strategies for AD.

Cat No.	Compound Name	Biochemical/ Physiological Action
Acetylcholine Nicotinic Receptors		
V098490	Varenicline Tartrate	$\alpha 4\beta 2$ partial agonist
E582200	Epibatidine	$\alpha 4\beta 2$ and $\alpha 3\beta 4$ agonist
D452485	Dihydro- β -erythroidine Hydrobromide	$\alpha 4$ nicotinic receptors Antagonist
P270280	Pentafluorobenzene	$\alpha 4\beta 2$ Antagonist
Amyloid β		
A604380	(S)-7-Amino-5H,7H-dibenzo[b,d]azepin-6-one	γ -secretase inhibitor, $A\beta$ levels
D422600	Dibenzazepine	γ -secretase inhibitor, $A\beta$ 42 levels
B596070	BMS 299897	γ -secretase inhibitor, $A\beta$ 40 levels
D193290	DAPT	$A\beta$ oligomer Inhibitor
Autophagy		
M275904	3-Methyl Adenine	Autophagy inhibitor via PI-3K
Q550550	6-Chloro-1,2,3,4-tetrahydro-N-[2-(4-methyl-1-piperazinyl)ethyl]-9-acridinamine	Autophagy & lysosomal activity inhibitor
M275905	3-Methyl Adenine-d3	Autophagy inhibitor via PI-3K (L)
Cholinesterase		
A178280	O-Acetylgalanthamine	Cholinesterase inhibitor
A190260	Acotiamide Hydrochloride	BChE inhibitor
C080100	Caffeine	Acetylcholinesterase inhibitor
T773855	Trichlorfon	Dual AChE and BChE inhibitor
Cyclooxygenases (COX)		
B679280	N-[4-(2-Bromoacetyl)phenyl]methanesulfonamide	COX 1 inhibitor
M264695	Desmethyl Naproxen	Selective COX inhibitor
N200500	Nabumetone	Non Selective COX inhibitor
DYRK		
H105250	Harmine	Selective DYRK _{1A} Inhibitor
I655600	INDY	Potent and Selective DYRK _{1A} Inhibitor
P755945	ProINDY	Non Selective DYRK _{1A} Inhibitor
A795155	AZ191	Potent DYRK _{1B} Inhibitor
Glutamate Receptors		

A603550	1-Aminocyclopropane-1-carboxylic Acid	<i>NMDA Agonist</i>
D525365	DL-AP5 Sodium Salt	<i>NMDA Antagonist</i>
R143450	Remacemide Hydrochloride	<i>NMDA Antagonist</i>
P285520	Perampanel	<i>AMPA Antagonist</i>
I232500	IDRA-21	<i>AMPA Modulator</i>
A634275	Ampalex	<i>AMPA Modulator</i>
B129500	BDP 37	<i>AMPA Modulator</i>
M311570	(2S,3R)-3-Methylglutamic Acid Hydrochloride Salt	<i>Kainate Agonist</i>
B684195	Bromoethane	<i>Kainate Antagonist</i>
B684198	Bromoethane-d3	<i>Kainate Antagonist</i>
Metabotropic Glutamate Receptors		
L486690	LY 367385	<i>mGlu₁ Antagonist (Group I)</i>
R637400	Ro 67-4853	<i>mGlu₁ Antagonist (Group I)</i>
V787575	VU 0483605	<i>mGlu₁ Antagonist (Group I)</i>
M745805	MPEP Hydrochloride	<i>mGlu₅ Antagonist (Group I)</i>
R637400	Ro 67-4853	<i>mGlu₅ Antagonist (Group I)</i>
A171990	Acetyl L-Carnitine Hydrochloride	<i>mGlu₂ Antagonist (Group II)</i>
A171991	Acetyl L-Carnitine-13C3 Chloride	<i>mGlu₂ Antagonist (Group II)</i>
A171992	Acetyl L-Carnitine-d3 Hydrochloride	<i>mGlu₂ Antagonist (Group II)</i>
L486680	LY 354740 Hydrate	<i>mGlu_{2/3} Antagonist (Group II)</i>
L486720	LY 379268	<i>mGlu₃ Antagonist (Group II)</i>
M425325	ML-289	<i>mGlu₃ Antagonist (Group II)</i>
C379860	N-(3-Chlorophenyl)-2-pyridinecarboxamide	<i>mGlu₄ Antagonist (Group III)</i>
T012600	TCN 238	<i>mGlu₄ Antagonist (Group III)</i>
V787555	VU 0364439	<i>mGlu₄ Antagonist (Group III)</i>
V787575	VU 0483605	<i>mGlu₄ Antagonist (Group III)</i>
X743850	XAP 044	<i>mGlu₇ Antagonist (Group III)</i>
L473475	LSN 2463359	<i>mGlu₈ Antagonist (Group III)</i>
GSK3 (Glycogen Synthase Kinase 3)		
A626820	3-[[[6-(3-Aminophenyl)-7H-pyrrolo[2,3-d]pyrimidin-4-yl]oxy]phenol	Potent GSK-3 β inhibitor
A808085	AZD1080	Selective GSK-3 inhibitor
A808095	A808095	GSK-3 α/β inhibitor
C367705	2-Chloro-4(1H)-pyrimidinone	Inhibits GSK-3 in vivo
γ-Secretase		
B682305	3-Bromo-2-chloropyridine	γ -Secretase Modulator
D422600	Dibenzazepine	Potent γ -Secretase Inhibitor
B596070	BMS 299897	Non selective A β 40 and A β 42 inhibitor
D193290	DAPT	γ -Secretase Inhibitor

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