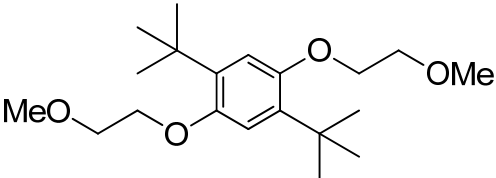


<b>Description</b>	1,4-di-tert-butyl-2,5-bis(2-methoxyethoxy)benzene
<b>CAS</b>	1350770-63-6
<b>Formula</b>	C <sub>20</sub> H <sub>34</sub> O <sub>4</sub>
<b>FW</b>	338.48
<b>LOT #</b>	KP01015
<b>Purity</b>	99.9%
<b>Batch Size</b>	154 g
<b>Manufactured</b>	3/4/2011
<b>Structure</b>	

Analysis	Instrument/Method	Results	Analysis By:
<b>HPLC</b>	Agilent 1260 Infinity Agilent Eclipse Plus C18, 3.5 um, 4.6x100, UV 225, water/ACN gradient	99.91% <sup>1</sup>	K. Pupek
	Agilent Eclipse Plus C18, 3.5 um, 4.6x100, UV 275, water/ACN gradient	99.94% <sup>1</sup>	K. Pupek
	Agilent Eclipse Plus C18, 3.5 um, 4.6x100, RID, water/ACN isocratic	100.00% <sup>1</sup>	K. Pupek
	Phenomenex Kinetex PFP, 2.6 um, 4.6x100, UV 225, water/ACN gradient	99.84% <sup>1</sup>	K. Pupek
<b>GC/MSD</b>	Agilent 7890A/5975C Triple-Axis Agilent HP-5MS, 0.25 um, 30m x 0.250 mm, 45 to 300 deg, 30 deg/min	99.96% <sup>1</sup> M <sup>+</sup> = 338	T. Dzwiniel
<b>Melting Point</b>	Buchi M-565 Automatic, range method	69.7-70.5 C	T. Michaelos
<b>KF Moisture Titration</b>	KEM MCU-610 Coulometric, WaterMark 1612/1613	15 ppm	K. Pupek
<b>FTIR</b>	Bruker Vertex 70 Attenuated Total Reflection	Consistent with Structure	S. Gallagher
<b>NMR</b>	Bruker 500 MHz <sup>1</sup> H observed in CDCl <sub>3</sub> solution	Consistent with Structure	K. Pupek

<sup>1</sup> By area integration.

Properties	Value
<b>Redox Potential (vs Li+/Li)</b>	4.00
<b>Solubility in Electrolyte (1.2M LiPF6 EC/EMC (3/7 V/V))</b>	0.5 M