

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002891.D
 Acq On : 13 Oct 2015 17:57
 Operator : UM/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

MMdadoda
 10/14/2015 6:25:51 PM

Quant Time: Oct 14 17:29:41 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 17:20:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	85503	5.00	ng	0.00
8) Naphthalene-d8	11.50	136	351739	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	189278	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	438365	5.00	ng	0.00
24) Chrysene-d12	22.04	240	401662	5.00	ng	0.00
26) Perylene-d12	24.61	264	380150	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	6.20	112	1786m	0.12	ng	0.06
3) Phenol-d6	7.83	99	2865	0.12	ng	0.05
9) Nitrobenzene-d5	9.90	82	1940m	0.09	ng	0.04
15) 2,4,6-Tribromophenol	16.74	330	658	0.10	ng	0.02
18) 2-Fluorobiphenyl	13.88	172	6975	0.11	ng	0.01
25) Terphenyl-d14	20.48	244	6875	0.11	ng	0.00
Target Compounds						
4) 2-Chlorophenol	8.24	128	2234	0.10	ng	Qvalue # 78
5) Phenol	7.86	94	2396m	0.10	ng	
6) 2-Methylphenol	9.15	107	2153	0.12	ng	# 66
7) 3+4-Methylphenols	9.51	107	2138	0.10	ng	# 45
10) 2-Nitrophenol	10.68	139	859	0.10	ng	# 33
11) 2,4-Dimethylphenol	10.68	122	2005	0.10	ng	89
12) 2,4-Dichlorophenol	11.26	162	1164	0.08	ng	# 83
13) 4-Chloro-3-methylphenol	12.81	107	1830	0.10	ng	# 71
16) 2,4,6-Trichlorophenol	13.76	196	1193	0.10	ng	# 64
17) 2,4,5-Trichlorophenol	13.88	196	1154	0.09	ng	# 95
22) 4,6-Dinitro-2-methylphenol	16.49	198	101	0.03	ng	# 1

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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