

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002898.D
 Acq On : 13 Oct 2015 22:12
 Operator : UM/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Quant Time: Oct 14 18:07:58 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	8.65	152	94663	5.00	ng	-0.01
8) Naphthalene-d8	11.50	136	405708	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	210227	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	456068	5.00	ng	0.00
24) Chrysene-d12	22.04	240	416552	5.00	ng	0.00
26) Perylene-d12	24.61	264	379983	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	6.12	112	198906	11.69	ng	-0.02
3) Phenol-d6	7.77	99	262644	10.10	ng	-0.02
9) Nitrobenzene-d5	9.84	82	231418	9.12	ng	-0.02
15) 2,4,6-Tribromophenol	16.71	330	93234	12.14	ng	0.00
18) 2-Fluorobiphenyl	13.87	172	658544	9.57	ng	0.00
25) Terphenyl-d14	20.48	244	621276	9.81	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) 2-Chlorophenol	8.20	128	250359	10.39	ng	# 76
5) Phenol	7.80	94	276484	10.02	ng	# 96
6) 2-Methylphenol	9.09	107	194357	9.94	ng	# 86
7) 3+4-Methylphenols	9.41	107	273441	11.31	ng	# 32
10) 2-Nitrophenol	10.61	139	125862	12.33	ng	# 79
11) 2,4-Dimethylphenol	10.64	122	245172	10.94	ng	# 94
12) 2,4-Dichlorophenol	11.15	162	214694	12.31	ng	# 92
13) 4-Chloro-3-methylphenol	12.74	107	250183	11.34	ng	# 71
16) 2,4,6-Trichlorophenol	13.70	196	165821	12.10	ng	# 74
17) 2,4,5-Trichlorophenol	13.79	196	180647	12.36	ng	# 98
19) 2,4-Dinitrophenol	15.40	184	29653	27.48	ng	# 66
20) 4-Nitrophenol	15.45	139	99605	2.79	ng	# 62
23) Pentachlorophenol	17.62	266	57239	16.54	ng	# 89

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

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