

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002899.D
 Acq On : 13 Oct 2015 22:48
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM101315

Quant Time: Oct 14 17:42:48 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 17:36:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	79629	5.00	ng	0.00
8) Naphthalene-d8	11.50	136	329015	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	175958	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	410542	5.00	ng	0.00
24) Chrysene-d12	22.04	240	398291	5.00	ng	0.00
26) Perylene-d12	24.61	264	364890	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	6.15	112	13107	0.85	ng	0.02
3) Phenol-d6	7.80	99	19115	0.87	ng	0.02
9) Nitrobenzene-d5	9.86	82	17117	0.94	ng	0.00
15) 2,4,6-Tribromophenol	16.71	330	5531	0.86	ng	0.00
18) 2-Fluorobiphenyl	13.88	172	53344	0.93	ng	0.01
25) Terphenyl-d14	20.48	244	54409	0.90	ng	0.00

Target Compounds						Qvalue
4) 2-Chlorophenol	8.21	128	18291	0.90	ng	# 70
5) Phenol	7.82	94	19566	0.91	ng	81
6) 2-Methylphenol	9.11	107	14596	0.89	ng	# 85
7) 3+4-Methylphenols	9.46	107	18035	0.89	ng	# 36
10) 2-Nitrophenol	10.64	139	7146	0.86	ng	# 59
11) 2,4-Dimethylphenol	10.66	122	16421	0.90	ng	100
12) 2,4-Dichlorophenol	11.21	162	13006	0.92	ng	94
13) 4-Chloro-3-methylphenol	12.76	107	15613	0.88	ng	# 86
16) 2,4,6-Trichlorophenol	13.72	196	9745	0.85	ng	# 81
17) 2,4,5-Trichlorophenol	13.83	196	10729	0.88	ng	# 97
19) 2,4-Dinitrophenol	15.52	184	260	0.82	ng	# 1
20) 4-Nitrophenol	15.52	139	3131	0.88	ng	# 76
22) 4,6-Dinitro-2-methylphenol	16.40	198	2043	0.87	ng	81
23) Pentachlorophenol	17.63	266	1891	0.96	ng	90

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

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