

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052215\
 Data File : BM001389.D
 Acq On : 22 May 2015 19:05
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC001

Quant Time: May 23 04:22:00 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 23 04:19:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	13479	5.00	ng	0.00
6) Naphthalene-d8	10.18	136	48287	5.00	ng	0.00
12) Acenaphthene-d10	14.08	164	27051	5.00	ng	0.00
18) Phenanthrene-d10	16.82	188	92052	5.00	ng	0.00
25) Chrysene-d12	21.04	240	58045	5.00	ng	0.00
32) Perylene-d12	23.14	264	55021	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	4.99	112	2749	1.04	ng	0.00
5) Phenol-d6	6.60	99	3162	0.97	ng	0.00
7) Nitrobenzene-d5	8.55	82	2797	1.06	ng	0.00
13) 2,4,6-Tribromophenol	15.59	330	854	0.92	ng	0.00
14) 2-Fluorobiphenyl	12.69	172	8641	1.06	ng	0.00
27) Terphenyl-d14	19.49	244	8030	1.04	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.87	88	1162	0.87	ng	100
3) n-Nitrosodimethylamine	3.20	42	1748	1.14	ng	100
8) Nitrobenzene	8.60	77	2925	1.06	ng	100
9) Naphthalene	10.22	128	10603	1.03	ng	100
10) Hexachlorobutadiene	10.52	225	2180	1.08	ng	100
11) 2-Methylnaphthalene	11.85	142	7177	0.98	ng	100
15) Acenaphthylene	13.79	152	11698	1.02	ng	100
16) Acenaphthene	14.14	154	7310	1.02	ng	100
17) Fluorene	15.12	166	13724	2.04	ng	100
19) 4-Bromophenyl-phenylether	16.04	248	1433	0.34	ng	# 100
20) Hexachlorobenzene	16.14	284	3673	0.57	ng	100
21) Pentachlorophenol	16.51	266	603	0.27	ng	100
22) Phenanthrene	16.85	178	18049	0.59	ng	100
23) Anthracene	16.95	178	18353	1.40	ng	100
24) Fluoranthene	18.90	202	16633	0.63	ng	100
26) Pyrene	19.26	202	17808	1.00	ng	100
28) Benzo(a)anthracene	21.02	228	16987	1.01	ng	100
29) Chrysene	21.07	228	15643	0.98	ng	100
30) Bis(2-ethylhexyl)phthalate	20.97	149	8371	1.01	ng	100
31) Indeno(1,2,3-cd)pyrene	25.20	276	17507	1.01	ng	100
33) Benzo(b)fluoranthene	22.51	252	16336	0.96	ng	100
34) Benzo(k)fluoranthene	22.55	252	16216	1.07	ng	100
35) Benzo(a)pyrene	23.04	252	14739	1.01	ng	100
36) Dibenzo(a,h)anthracene	25.21	278	13528	1.01	ng	100
37) Benzo(g,h,i)perylene	25.83	276	14373	1.00	ng	100

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

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