

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050515\
 Data File : BM001214.D
 Acq On : 05 May 2015 18:16
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Quant Time: May 06 06:13:16 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 06 02:27:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	14433	5.00	ng	0.00
6) Naphthalene-d8	10.30	136	57411	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	35478	5.00	ng	0.00
18) Phenanthrene-d10	16.95	188	60443	5.00	ng	0.00
24) Chrysene-d12	21.14	240	77970	5.00	ng	0.00
30) Perylene-d12	23.29	264	64496	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.13	112	310	0.11	ng	0.02
5) Phenol-d6	6.72	99	362	0.10	ng	0.00
7) Nitrobenzene-d5	8.69	82	251	0.08	ng	0.01
13) 2,4,6-Tribromophenol	15.69	330	132	0.09	ng	0.00
14) 2-Fluorobiphenyl	12.81	172	1045	0.09	ng	0.00
26) Terphenyl-d14	19.59	244	1046	0.10	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.37	42	188	0.11	ng	# 85
8) Nitrobenzene	8.73	77	269	0.08	ng	99
9) Naphthalene	10.35	128	1228	0.10	ng	# 87
10) Hexachlorobutadiene	10.64	225	255	0.10	ng	94
11) 2-Methylnaphthalene	11.99	142	838	0.10	ng	95
15) Acenaphthylene	13.90	152	1317	0.09	ng	99
16) Acenaphthene	14.25	154	938	0.09	ng	99
17) Fluorene	15.23	166	1484	0.12	ng	94
19) Hexachlorobenzene	16.24	284	294	0.09	ng	# 83
21) Phenanthrene	16.98	178	2615	0.15	ng	97
22) Anthracene	17.07	178	2059	0.15	ng	98
23) Fluoranthene	19.01	202	2177	0.13	ng	100
25) Pyrene	19.37	202	2256	0.09	ng	100
27) Benzo(a)anthracene	21.12	228	2226	0.10	ng	97
28) Chrysene	21.17	228	2092	0.10	ng	97
29) Indeno(1,2,3-cd)pyrene	25.43	276	1677	0.09	ng	98
31) Benzo(b)fluoranthene	22.64	252	1766	0.09	ng	# 81
32) Benzo(k)fluoranthene	22.68	252	1884	0.10	ng	# 83
33) Benzo(a)pyrene	23.18	252	1550	0.09	ng	# 73
34) Dibenzo(a,h)anthracene	25.44	278	1321	0.09	ng	# 78
35) Benzo(g,h,i)perylene	26.08	276	1406	0.09	ng	# 85

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

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