

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041815\
 Data File : BM001042.D
 Acq On : 17 Apr 2015 22:44
 Operator : TP/IZ
 Sample : SSTDICC0.2
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.2

Quant Time: Apr 18 07:48:02 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM041815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 18 05:45:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	16676	5.00	ng	0.00
4) Naphthalene-d8	10.36	136	59174	5.00	ng	0.00
8) Acenaphthene-d10	14.24	164	30693	5.00	ng	0.00
13) Phenanthrene-d10	16.99	188	66039	5.00	ng	0.00
19) Chrysene-d12	21.19	240	55390	5.00	ng	0.00
25) Perylene-d12	23.36	264	50636	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.74	82	552	0.19	ng	0.00
9) 2-Fluorobiphenyl	12.86	172	1995	0.21	ng	0.00
21) Terphenyl-d14	19.63	244	1401	0.21	ng	0.00
Target Compounds						Ovalue
3) n-Nitrosodimethylamine	3.43	42	282	0.18	ng	# 98
6) Naphthalene	10.41	128	2778	0.20	ng	95
7) 2-Methylnaphthalene	12.04	142	1745	0.20	ng	# 90
10) Acenaphthylene	13.95	152	2425	0.18	ng	99
11) Acenaphthene	14.29	154	1902	0.20	ng	95
12) Fluorene	15.30	166	2100	0.20	ng	98
14) Hexachlorobenzene	16.31	284	710	0.19	ng	# 100
15) Pentachlorophenol	16.66	266	128	0.13	ng	# 82
16) Phenanthrene	17.03	178	3394	0.19	ng	99
17) Anthracene	17.11	178	2962	0.19	ng	98
18) Fluoranthene	19.05	202	3408	0.19	ng	100
20) Pyrene	19.42	202	3534	0.20	ng	99
22) Benzo(a)anthracene	21.17	228	3063	0.19	ng	97
23) Chrysene	21.22	228	3398	0.21	ng	96
24) Indeno(1,2,3-cd)pyrene	25.52	276	3343	0.19	ng	99
26) Benzo(b)fluoranthene	22.70	252	2957	0.19	ng	# 90
27) Benzo(k)fluoranthene	22.74	252	3047	0.20	ng	# 90
28) Benzo(a)pyrene	23.25	252	2578	0.19	ng	# 86
29) Dibenzo(a,h)anthracene	25.54	278	2574	0.19	ng	# 91
30) Benzo(g,h,i)perylene	26.18	276	2895	0.20	ng	94

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

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