

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000764.D
 Acq On : 31 Mar 2015 21:50
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC002

Quant Time: Apr 01 04:06:27 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM033115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 03:53:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	36469	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	126260	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	62019	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	128075	5.00	ng	0.00
19) Chrysene-d12	21.21	240	109129	5.00	ng	0.00
25) Perylene-d12	23.40	264	97244	5.00	ng	0.00

System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	13416	2.10	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	40370	2.70	ng	0.00
21) Terphenyl-d14	19.66	244	27711	1.78	ng	0.00

Target Compounds				Qvalue		
3) n-Nitrosodimethylamine	3.44	42	7490	1.45	ng	# 99
6) Naphthalene	10.45	128	61167	2.42	ng	100
7) 2-Methylnaphthalene	12.07	142	39217	2.46	ng	97
10) Acenaphthylene	13.98	152	61004	0.78	ng	100
11) Acenaphthene	14.33	154	38130	2.77	ng	99
12) Fluorene	15.32	166	45269	2.67	ng	99
14) Hexachlorobenzene	16.33	284	15559	3.61	ng	# 100
15) Pentachlorophenol	16.68	266	3543	0.08	ng	97
16) Phenanthrene	17.05	178	71105	2.55	ng	100
17) Anthracene	17.15	178	61302	2.48	ng	99
18) Fluoranthene	19.08	202	70937	2.66	ng	100
20) Pyrene	19.44	202	74880	2.68	ng	100
22) Benzo(a)anthracene	21.19	228	64618	2.99	ng	# 88
23) Chrysene	21.24	228	67492	2.96	ng	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	70417	4.33	ng	100
26) Benzo(b)fluoranthene	22.74	252	66870	3.59	ng	99
27) Benzo(k)fluoranthene	22.79	252	60101	2.74	ng	99
28) Benzo(a)pyrene	23.30	252	55576	3.36	ng	95
29) Dibenzo(a,h)anthracene	25.60	278	55498	4.46	ng	99
30) Benzo(g,h,i)perylene	26.26	276	58940	3.72	ng	99

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

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