

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000762.D
 Acq On : 31 Mar 2015 20:39
 Operator : TP/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.8

Quant Time: Apr 01 04:05:45 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	33909	5.00	ng	0.00
4) Naphthalene-d8	10.40	136	118580	5.00	ng	0.00
8) Acenaphthene-d10	14.28	164	58715	5.00	ng	0.00
13) Phenanthrene-d10	17.01	188	122880	5.00	ng	0.00
19) Chrysene-d12	21.22	240	110719	5.00	ng	0.01
25) Perylene-d12	23.40	264	101421	5.00	ng	0.00

System Monitoring Compounds						
5) Nitrobenzene-d5	8.77	82	4485	0.75	ng	0.00
9) 2-Fluorobiphenyl	12.89	172	14637	1.03	ng	0.00
21) Terphenyl-d14	19.66	244	10519	0.67	ng	0.00

Target Compounds				Qvalue		
3) n-Nitrosodimethylamine	3.45	42	2496	0.52	ng	99
6) Naphthalene	10.45	128	22137	0.93	ng	100
7) 2-Methylnaphthalene	12.07	142	13899	0.93	ng	98
10) Acenaphthylene	13.98	152	19977	0.27	ng	100
11) Acenaphthene	14.33	154	14095	1.08	ng	99
12) Fluorene	15.32	166	16157	1.01	ng	100
14) Hexachlorobenzene	16.33	284	5659	1.37	ng	# 100
15) Pentachlorophenol	16.68	266	842	0.02	ng	99
16) Phenanthrene	17.05	178	26849	1.00	ng	100
17) Anthracene	17.15	178	22373	0.94	ng	99
18) Fluoranthene	19.08	202	26738	1.05	ng	100
20) Pyrene	19.45	202	27872	0.98	ng	100
22) Benzo(a)anthracene	21.20	228	24244	1.11	ng	98
23) Chrysene	21.25	228	27061	1.17	ng	# 89
24) Indeno(1,2,3-cd)pyrene	25.59	276	27459	1.67	ng	100
26) Benzo(b)fluoranthene	22.74	252	24687	1.27	ng	99
27) Benzo(k)fluoranthene	22.79	252	25023	1.09	ng	99
28) Benzo(a)pyrene	23.31	252	21357	1.24	ng	99
29) Dibenzo(a,h)anthracene	25.60	278	21432	1.65	ng	99
30) Benzo(g,h,i)perylene	26.26	276	23149	1.40	ng	99

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

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