

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051315\
 Data File : BM001271.D
 Acq On : 13 May 2015 03:17
 Operator : TP/IZ
 Sample : G2183-01MS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-1(7-8)MS

Quant Time: May 13 05:50:10 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 13 05:24:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	17664	5.00	ng	0.00
6) Naphthalene-d8	10.22	136	68869	5.00	ng	0.00
12) Acenaphthene-d10	14.12	164	39293	5.00	ng	0.00
18) Phenanthrene-d10	16.85	188	101863	5.00	ng	0.00
24) Chrysene-d12	21.07	240	70811	5.00	ng	-0.01
30) Perylene-d12	23.18	264	58084	5.00	ng	-0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.04	112	39852	11.58	ng	0.00
5) Phenol-d6	6.63	99	54420	14.02	ng	0.00
7) Nitrobenzene-d5	8.60	82	36062	8.23	ng	0.00
13) 2,4,6-Tribromophenol	15.63	330	16201	11.99	ng	0.00
14) 2-Fluorobiphenyl	12.73	172	102474	8.40	ng	0.00
26) Terphenyl-d14	19.51	244	83446	8.77	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.90	88	10135	7.17	ng	# 58
3) n-Nitrosodimethylamine	3.23	42	16261	8.13	ng	# 98
8) Nitrobenzene	8.64	77	39565	8.40	ng	97
9) Naphthalene	10.27	128	116267	7.93	ng	98
10) Hexachlorobutadiene	10.56	225	21969	7.67	ng	99
11) 2-Methylnaphthalene	11.90	142	80871	8.48	ng	94
15) Acenaphthylene	13.83	152	147744	9.13	ng	99
16) Acenaphthene	14.18	154	84307	8.14	ng	98
17) Fluorene	15.17	166	68098	7.26	ng	# 18
19) Hexachlorobenzene	16.18	284	51533	7.86	ng	# 94
20) Pentachlorophenol	16.52	266	41247	20.90	ng	88
21) Phenanthrene	16.88	178	112259	7.25	ng	# 98
22) Anthracene	16.98	178	122879	8.14	ng	# 97
23) Fluoranthene	18.93	202	193723	7.97	ng	99
25) Pyrene	19.30	202	201421	9.20	ng	99
27) Benzo(a)anthracene	21.06	228	172810	8.99	ng	95
28) Chrysene	21.11	228	165180	8.48	ng	96
29) Indeno(1,2,3-cd)pyrene	25.27	276	143225	7.23	ng	99
31) Benzo(b)fluoranthene	22.56	252	166070	9.75	ng	98
32) Benzo(k)fluoranthene	22.60	252	135811	8.39	ng	98
33) Benzo(a)pyrene	23.09	252	140604	9.54	ng	97
34) Dibenzo(a,h)anthracene	25.28	278	115074	8.30	ng	97
35) Benzo(g,h,i)perylene	25.90	276	119426	7.85	ng	97

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

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