

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060615\
 Data File : BM001594.D
 Acq On : 06 Jun 2015 14:39
 Operator : TP/IZ
 Sample : G2491-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SS043MW100-150602

Quant Time: Jun 08 05:39:39 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 05:27:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2	5.00	ng	0.00
6) Naphthalene-d8	10.47	136	6	5.00	ng	0.00
12) Acenaphthene-d10	14.34	164	11	5.00	ng	-0.01
18) Phenanthrene-d10	17.09	188	12	5.00	ng	0.00
25) Chrysene-d12	21.28	240	104	5.00	ng	0.00
32) Perylene-d12	23.51	264	138	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.22	112	1	2.26	ng	-0.05
5) Phenol-d6	6.86	99	3	5.53	ng	0.02
7) Nitrobenzene-d5	8.86	82	6	14.68	ng	0.01
13) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
14) 2-Fluorobiphenyl	12.96	172	6	1.77	ng	0.00
27) Terphenyl-d14	19.72	244	16	1.19	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.17	88	1	5.70	ng	# 1
3) n-Nitrosodimethylamine	3.49	42	4	13.10	ng	# 1
8) Nitrobenzene	8.95	77	4	9.13	ng	# 43
9) Naphthalene	10.47	128	10	7.72	ng	# 1
11) 2-Methylnaphthalene	12.09	142	1	1.23	ng	# 58
15) Acenaphthylene	14.04	152	7	1.47	ng	# 1
16) Acenaphthene	14.40	154	5	1.67	ng	# 43
17) Fluorene	15.36	166	2	1.04	ng	# 65
19) 4-Bromophenyl-phenylether	16.27	248	8	8.38	ng	# 18
21) Pentachlorophenol	16.75	266	2	10.91	ng	# 66
22) Phenanthrene	17.12	178	41	9.79	ng	# 66
23) Anthracene	17.12	178	49	27.79	ng	# 34
24) Fluoranthene	19.15	202	38	11.67	ng	# 70
26) Pyrene	19.52	202	49	1.46	ng	# 67
28) Benzo(a)anthracene	21.25	228	53	1.81	ng	# 1
29) Chrysene	21.31	228	84	3.00	ng	# 7
30) Bis(2-ethylhexyl)phthalate	21.19	149	128	6.53	ng	# 54
31) Indeno(1,2,3-cd)pyrene	25.76	276	256	8.52	ng	# 61
33) Benzo(b)fluoranthene	22.83	252	82	1.99	ng	# 1
34) Benzo(k)fluoranthene	22.88	252	94	2.50	ng	# 1
35) Benzo(a)pyrene	23.42	252	94	2.61	ng	# 1
36) Dibenzo(a,h)anthracene	25.79	278	216	6.12	ng	# 1
37) Benzo(g,h,i)perylene	26.46	276	185	4.98	ng	# 13

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

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