

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090415\
 Data File : BE090734.D
 Acq On : 4 Sep 2015 17:49
 Operator : UM/IZ
 Sample : G3562-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 090215-D534-F-U-B

Quant Time: Sep 04 18:28:55 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2	5.00	ng	0.02
6) Naphthalene-d8	10.31	136	13	5.00	ng	0.00
12) Acenaphthene-d10	14.15	164	7	5.00	ng	-0.02
18) Phenanthrene-d10	16.86	188	26	5.00	ng	-0.05
25) Chrysene-d12	21.00	240	20	5.00	ng	-0.08
32) Perylene-d12	23.25	264	20	5.00	ng	-0.12

System Monitoring Compounds						
4) 2-Fluorophenol	5.20	112	4	10.25	ng	0.00
5) Phenol-d6	6.77	99	7	12.57	ng	0.02
7) Nitrobenzene-d5	8.71	82	11	12.80	ng	0.00
13) 2,4,6-Tribromophenol	15.57	330	28	133.76	ng	-0.10
14) 2-Fluorobiphenyl	12.77	172	4	2.06	ng	-0.02
27) Terphenyl-d14	19.47	244	5	2.26	ng	-0.06

Target Compounds						Qvalue
2) 1,4-Dioxane	3.18	88	18	71.69	ng	# 69
3) n-Nitrosodimethylamine	3.47	42	69	198.24	ng	# 1
8) Nitrobenzene	8.75	77	19	17.71	ng	# 38
9) Naphthalene	10.39	128	30	10.58	ng	# 1
10) Hexachlorobutadiene	10.65	225	9	20.29	ng	# 43
11) 2-Methylnaphthalene	11.97	142	17	11.09	ng	# 53
15) Acenaphthylene	13.86	152	23	2.06	ng	# 1
16) Acenaphthene	14.20	154	11	5.88	ng	# 8
17) Fluorene	15.19	166	57	24.45	ng	# 6
19) 4-Bromophenyl-phenylether	16.13	248	24	27.08	ng	# 1
20) Hexachlorobenzene	16.15	284	31	7.39	ng	# 37
21) Pentachlorophenol	16.53	266	6	16.13	ng	# 58
22) Phenanthrene	16.93	178	17	2.64	ng	# 1
23) Anthracene	17.01	178	66	12.27	ng	# 69
24) Fluoranthene	18.91	202	7	1.03	ng	# 30
26) Pyrene	19.25	202	11	2.66	ng	# 20
28) Benzo(a)anthracene	21.00	228	18	4.10	ng	# 1
29) Chrysene	21.04	228	10	2.12	ng	# 1
30) Bis(2-ethylhexyl)phthalate	20.92	149	71	23.68	ng	# 54
31) Indeno(1,2,3-cd)pyrene	25.62	276	13	2.75	ng	# 50
33) Benzo(b)fluoranthene	22.55	252	13	3.00	ng	# 1
34) Benzo(k)fluoranthene	22.60	252	11	2.18	ng	# 1
35) Benzo(a)pyrene	23.15	252	22	5.67	ng	# 1
36) Dibenzo(a,h)anthracene	25.62	278	1	0.34	ng	# 1
37) Benzo(g,h,i)perylene	26.32	276	10	2.24	ng	# 1

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

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