

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100215\
 Data File : BG018992.D
 Acq On : 1 Oct 2015 23:18
 Operator : UM/NP
 Sample : G3865-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E5MY8

Manual Integrations
 APPROVED

MMDadoda
 10/2/2015 5:44:36 PM

Quant Time: Oct 02 13:04:09 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 02 11:51:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.97	152	48477	20.00	ng/ul	0.00
18) Naphthalene-d8	10.78	136	229565	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.60	164	159630	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.35	188	409742	20.00	ng/ul	0.00
78) Chrysene-d12	21.60	240	500794	20.00	ng/ul	-0.01
86) Perylene-d12	24.77	264	501195	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.40	96	3502	3.47	ng/uL	0.00
5) Phenol-d5	7.22	99	36690m	9.22	ng/ul	0.04
7) Bis-(2-Chloroethyl)ether-d	7.31	67	31147	13.50	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	50872m	16.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.75	113	28861m	8.94	ng/ul	0.04
19) Nitrobenzene-d5	9.15	128	28603m	16.45	ng/ul	0.00
22) 2-Nitrophenol-d4	9.88	143	31075m	17.77	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.45	165	56225m	15.27	ng/ul	0.02
29) 4-Chloroaniline-d4	10.96	131	62056m	14.48	ng/ul	0.02
44) Dimethylphthalate-d6	14.01	166	216213m	16.83	ng/ul	0.00
47) Acenaphthylene-d8	14.30	160	296414	19.47	ng/ul	0.00
52) 4-Nitrophenol-d4	14.93	143	27445m	11.13	ng/ul	0.04
58) Fluorene-d10	15.59	176	221895	19.75	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	15172m	7.82	ng/ul	0.00
71) Anthracene-d10	17.44	188	373593	20.71	ng/ul	0.00
79) Pyrene-d10	19.73	212	455664	19.78	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.55	264	488779	20.04	ng/ul	-0.02

Target Compounds					Qvalue
45) Dimethylphthalate	14.05	163	136128	10.93 ng/ul#	98

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

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