

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100215\
 Data File : BG018993.D
 Acq On : 1 Oct 2015 23:56
 Operator : UM/NP
 Sample : G3865-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E5MY6

Manual Integrations
 APPROVED

MMDadoda
 10/2/2015 5:44:41 PM

Quant Time: Oct 02 13:09:09 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	7.98	152	49579	20.00	ng/ul	0.00
18) Naphthalene-d8	10.78	136	223042	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.60	164	152841	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.35	188	399795	20.00	ng/ul	0.00
78) Chrysene-d12	21.60	240	481742	20.00	ng/ul	-0.01
86) Perylene-d12	24.78	264	471435	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.40	96	5293	5.13	ng/uL	0.00
5) Phenol-d5	7.21	99	58410m	14.35	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.31	67	51296	21.74	ng/ul	0.00
9) 2-Chlorophenol-d4	7.52	132	65111	21.12	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	60808m	18.41	ng/ul	0.02
19) Nitrobenzene-d5	9.15	128	46059m	27.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	48465m	28.52	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.45	165	76702m	21.44	ng/ul	0.02
29) 4-Chloroaniline-d4	10.96	131	92079m	22.12	ng/ul	0.02
44) Dimethylphthalate-d6	14.01	166	357503	29.06	ng/ul	0.00
47) Acenaphthylene-d8	14.30	160	440901	30.25	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	50856m	21.54	ng/ul	0.03
58) Fluorene-d10	15.59	176	330524	30.73	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	31301	16.54	ng/ul	0.00
71) Anthracene-d10	17.44	188	549998	31.25	ng/ul	0.00
79) Pyrene-d10	19.73	212	665233	30.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.56	264	712264	31.04	ng/ul	-0.01

Target Compounds					Qvalue
45) Dimethylphthalate	14.05	163	199197	16.71 ng/ul	99

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

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