

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100215\
 Data File : BG018985.D
 Acq On : 1 Oct 2015 18:53
 Operator : UM/NP
 Sample : G3831-09
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AS2

Manual Integrations
 APPROVED

MMDadoda
 10/2/2015 5:43:50 PM

Quant Time: Oct 02 12:07:45 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	54245	20.00	ng/ul	0.00
18) Naphthalene-d8	10.78	136	230005	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.60	164	146384	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.35	188	378796m	20.00	ng/ul	0.00
78) Chrysene-d12	21.60	240	490805	20.00	ng/ul	0.00
86) Perylene-d12	24.78	264	477748	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.40	96	1054	0.93	ng/uL	0.00
5) Phenol-d5	7.23	99	8564	1.92	ng/ul	0.04
7) Bis-(2-Chloroethyl)ether-d	7.31	67	51934	20.12	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	60809	18.02	ng/ul	0.01
13) 4-Methylphenol-d8	8.74	113	38539m	10.66	ng/ul	0.02
19) Nitrobenzene-d5	9.14	128	43227	24.81	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	46361	26.46	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.45	165	80970m	21.94	ng/ul	0.02
29) 4-Chloroaniline-d4	10.99	131	6366m	1.48	ng/ul	0.05
44) Dimethylphthalate-d6	14.01	166	350407m	29.74	ng/ul	0.00
47) Acenaphthylene-d8	14.29	160	421283	30.18	ng/ul	0.00
52) 4-Nitrophenol-d4	14.96	143	7669m	3.39	ng/ul	0.07
58) Fluorene-d10	15.59	176	314495	30.53	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	32526m	18.14	ng/ul	0.00
71) Anthracene-d10	17.44	188	513815	30.81	ng/ul	0.00
79) Pyrene-d10	19.73	212	638693	28.29	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.56	264	710017	30.54	ng/ul	-0.01

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

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