

Data Path : Z:\HPCHEM1\BNA G\DATA\BG111715\
 Data File : BG019690.D
 Acq On : 18 Nov 2015 11:59
 Operator : UM/NP
 Sample : G4374-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Nov 19 03:09:26 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG111615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 19 03:06:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	24988	20.00	ng/ul	0.00
18) Naphthalene-d8	10.98	136	108397	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.78	164	91315	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.52	188	259317	20.00	ng/ul	0.00
78) Chrysene-d12	21.79	240	328116	20.00	ng/ul	0.00
86) Perylene-d12	25.11	264	310654	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	422	0.72	ng/uL	0.00
5) Phenol-d5	7.30	99	14539	5.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.46	67	44564	26.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.67	132	34992	21.32	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	29205	13.78	ng/ul	0.00
19) Nitrobenzene-d5	9.32	128	25614	28.78	ng/ul	0.00
22) 2-Nitrophenol-d4	10.05	143	29056	29.01	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.59	165	58159	26.85	ng/ul	0.00
29) 4-Chloroaniline-d4	11.12	131	1828	0.89	ng/ul	0.01
44) Dimethylphthalate-d6	14.18	166	235420	28.51	ng/ul	0.00
47) Acenaphthylene-d8	14.48	160	258549	28.88	ng/ul	0.00
52) 4-Nitrophenol-d4	14.97	143	6847	5.26	ng/ul	0.00
58) Fluorene-d10	15.77	176	216996	29.36	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.88	200	39924	23.83	ng/ul	0.00
71) Anthracene-d10	17.62	188	356239	28.29	ng/ul	0.00
79) Pyrene-d10	19.90	212	435716	29.14	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.88	264	448963	27.68	ng/ul	0.00

Target Compounds Ovalue

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

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