

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100515\
 Data File : BG019041.D
 Acq On : 6 Oct 2015 00:01
 Operator : UM/NP
 Sample : G3900-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AA3

Quant Time: Oct 06 05:37:29 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 06 04:27:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	37546	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	175709	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	116943	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	265605	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	291360	20.00	ng/ul	0.00
86) Perylene-d12	24.91	264	282378	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	3919	4.72	ng/uL	0.00
5) Phenol-d5	7.22	99	118954	36.76	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.37	67	72705	34.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	88236	36.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.75	113	94783	37.26	ng/ul	0.01
19) Nitrobenzene-d5	9.21	128	46400	34.93	ng/ul	0.00
22) 2-Nitrophenol-d4	9.93	143	54239	39.69	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	103188	37.73	ng/ul	0.01
29) 4-Chloroaniline-d4	10.99	131	91661	27.48	ng/ul	0.00
44) Dimethylphthalate-d6	14.08	166	321898	34.91	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	396296	34.75	ng/ul	0.00
52) 4-Nitrophenol-d4	14.87	143	65122	34.96	ng/ul	0.02
58) Fluorene-d10	15.66	176	292380	35.02	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.77	200	49790	34.56	ng/ul	0.00
71) Anthracene-d10	17.52	188	438260	35.04	ng/ul	0.00
79) Pyrene-d10	19.79	212	478200	35.44	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.69	264	525948	36.40	ng/ul	0.02

Target Compounds

						Ovalue
6) Phenol	7.24	94	3995	1.21	ng/ul	97
45) Dimethylphthalate	14.12	163	189978	20.18	ng/ul	99

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

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