

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112015\
 Data File : BG019732.D
 Acq On : 20 Nov 2015 23:22
 Operator : UM/NP
 Sample : PB86714BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Nov 21 01:08:51 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG111615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 21 01:06:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	28315	20.00	ng/ul	0.00
18) Naphthalene-d8	10.95	136	104834	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.76	164	83910	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.50	188	229281	20.00	ng/ul	0.00
78) Chrysene-d12	21.77	240	309446	20.00	ng/ul	0.00
86) Perylene-d12	25.06	264	295990	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.41	96	3392	5.12	ng/uL	0.00
5) Phenol-d5	7.27	99	63172	21.78	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.43	67	40324	21.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	40661	21.86	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	48507	20.20	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	21080	24.49	ng/ul	0.00
22) 2-Nitrophenol-d4	10.02	143	24293	25.08	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	47798	22.82	ng/ul	0.00
29) 4-Chloroaniline-d4	11.08	131	59283	29.81	ng/ul	0.00
44) Dimethylphthalate-d6	14.15	166	184163	24.27	ng/ul	0.00
47) Acenaphthylene-d8	14.45	160	190376	23.14	ng/ul	0.00
52) 4-Nitrophenol-d4	14.95	143	26286	21.98	ng/ul	0.00
58) Fluorene-d10	15.75	176	155704	22.93	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.86	200	26275	17.74	ng/ul	0.00
71) Anthracene-d10	17.60	188	286967	25.78	ng/ul	0.00
79) Pyrene-d10	19.88	212	373431	26.48	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.84	264	411917	26.66	ng/ul	0.00

Target Compounds Ovalue

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

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