

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102115\
 Data File : BG019284.D
 Acq On : 21 Oct 2015 17:10
 Operator : UM/NP
 Sample : PB86173BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK73

Quant Time: Oct 21 18:15:29 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 21 12:55:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	14275	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	66206	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.62	164	49097	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.37	188	115766	20.00	ng/ul	0.00
78) Chrysene-d12	21.63	240	125535	20.00	ng/ul	0.00
86) Perylene-d12	24.83	264	127355	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.42	96	1319	5.15	ng/uL	0.00
5) Phenol-d5	7.16	99	36939	29.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.31	67	22640	28.24	ng/ul	0.00
9) 2-Chlorophenol-d4	7.52	132	28514	31.40	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	30950	29.32	ng/ul	0.00
19) Nitrobenzene-d5	9.15	128	15357	30.11	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	18341	33.88	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.42	165	34098	31.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.93	131	43463	31.48	ng/ul	0.00
44) Dimethylphthalate-d6	14.03	166	135642	31.59	ng/ul	0.00
47) Acenaphthylene-d8	14.31	160	150206	32.36	ng/ul	0.00
52) 4-Nitrophenol-d4	14.84	143	23155	30.00	ng/ul	0.00
58) Fluorene-d10	15.61	176	114464	31.19	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.74	200	21390	29.95	ng/ul	0.00
71) Anthracene-d10	17.47	188	191913	35.09	ng/ul	0.00
79) Pyrene-d10	19.76	212	214674	37.70	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.61	264	232948	36.21	ng/ul	0.00

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

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