

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100615\
 Data File : BG019052.D
 Acq On : 6 Oct 2015 19:19
 Operator : UM/NP
 Sample : PB85947BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK47

Quant Time: Oct 07 01:13:10 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 07 01:07:23 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	16145	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	74489	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	51529	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	149342	20.00	ng/ul	0.00
78) Chrysene-d12	21.69	240	199576	20.00	ng/ul	0.00
86) Perylene-d12	24.92	264	192504	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	1877	5.25	ng/uL	0.00
5) Phenol-d5	7.20	99	43702	31.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.37	67	30013	33.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.58	132	33594	32.17	ng/ul	0.00
13) 4-Methylphenol-d8	8.75	113	36786	33.63	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	18335	32.56	ng/ul	0.00
22) 2-Nitrophenol-d4	9.93	143	20296	35.03	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	36559	31.53	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	50134	35.45	ng/ul	0.00
44) Dimethylphthalate-d6	14.08	166	156561	38.53	ng/ul	0.00
47) Acenaphthylene-d8	14.36	160	161851	32.21	ng/ul	0.00
52) 4-Nitrophenol-d4	14.86	143	33268	40.53	ng/ul	0.00
58) Fluorene-d10	15.66	176	128383	34.90	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.77	200	23428	28.92	ng/ul	0.00
71) Anthracene-d10	17.51	188	239624	34.07	ng/ul	0.00
79) Pyrene-d10	19.80	212	306767	33.19	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.71	264	334111	33.92	ng/ul	0.00

Target Compounds Ovalue

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

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