

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060115\
 Data File : BG017367.D
 Acq On : 1 Jun 2015 11:57
 Operator : TP/IZ
 Sample : PB83702BL
 Misc : SBLK63
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK63

Quant Time: Jun 02 05:34:34 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG052915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 02 05:29:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	21940	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	100161	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.31	164	70005	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	168347	20.00	ng/ul	0.00
75) Chrysene-d12	21.24	240	182524	20.00	ng/ul	0.00
83) Perylene-d12	23.48	264	174750	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	3051	6.61	ng/uL	0.00
5) Phenol-d5	6.84	99	54625	29.54	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	36750	33.24	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	45969	30.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	47592	30.03	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	24097	31.06	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	27785	30.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	43602	26.86	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	70612	37.89	ng/ul	0.00
43) Dimethylphthalate-d6	13.72	166	162110	32.48	ng/ul	0.00
46) Acenaphthylene-d8	13.99	160	202243	31.29	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	26117	27.43	ng/ul	0.00
57) Fluorene-d10	15.30	176	157328	32.71	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.43	200	26235	22.27	ng/ul	0.00
70) Anthracene-d10	17.16	188	248715	31.53	ng/ul	0.00
76) Pyrene-d10	19.45	212	294672	33.34	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.34	264	256742	32.91	ng/ul	0.00

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

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