

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121715\
 Data File : BG020226.D
 Acq On : 16 Dec 2015 17:27
 Operator : UM/SJ
 Sample : PB7250BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK50

Quant Time: Dec 17 00:44:10 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 17 00:37:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	16815	20.00	ng/ul	0.00
18) Naphthalene-d8	10.91	136	78432	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.73	164	62625	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.46	188	161192	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	186881	20.00	ng/ul	0.00
86) Perylene-d12	25.00	264	170322	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.24	99	170	0.11	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.42	67	214	0.24	ng/ul	0.00
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.80	113	122	0.10	ng/ul	0.00
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	9.99	143	107	0.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	11.05	131	119	0.08	ng/ul	0.00
44) Dimethylphthalate-d6	14.12	166	1503	0.30	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	1436	0.24	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.71	176	1740	0.38	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.46	188	161192	21.70	ng/ul	-0.10
79) Pyrene-d10	19.84	212	1954	0.22	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.78	264	1584	0.19	ng/ul	0.00

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

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