

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032816\
 Data File : BG021523.D
 Acq On : 28 Mar 2016 12:58
 Operator : UM/SJ
 Sample : PB89199BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK99

Quant Time: Mar 29 01:40:52 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 29 01:39:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	8094	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	35519	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	23815	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	64611	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	84021	20.00	ng/ul	0.00
86) Perylene-d12	24.90	264	63292	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	1111	6.16	ng/uL	0.00
5) Phenol-d5	7.34	99	21975	28.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	14135	31.18	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	18030	31.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	19395	29.59	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	10084	37.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	10953	35.96	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	17896	32.87	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	4756	6.68	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	79594	40.14	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	86638	35.44	ng/ul	0.00
52) 4-Nitrophenol-d4	15.09	143	11647	36.75	ng/ul	0.00
58) Fluorene-d10	15.66	176	65052	36.39	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	15279	35.79	ng/ul	0.00
71) Anthracene-d10	17.51	188	115607	36.51	ng/ul	0.00
79) Pyrene-d10	19.79	212	144633	36.91	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	120372	41.21	ng/ul	0.00

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

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