

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032216\
 Data File : BG021465.D
 Acq On : 22 Mar 2016 8:34
 Operator : UM/SJ
 Sample : PB89036BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK36

Quant Time: Mar 23 01:32:04 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 23 01:30:27 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	13533	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.88	136	65449	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	43405	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	94441	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	110047	20.00	ng/ul	0.00
86) Perylene-d12	24.89	264	105967	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	1630	5.40	ng/uL	0.00
5) Phenol-d5	7.33	99	36462	28.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.39	67	24176	31.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	30028	30.90	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	33337	30.42	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	16697	34.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	19025	33.90	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	31870	31.76	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	47685	36.35	ng/ul	0.00
44) Dimethylphthalate-d6	14.08	166	124236	34.37	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	147543	33.12	ng/ul	0.00
52) 4-Nitrophenol-d4	15.07	143	13654	23.64	ng/ul	0.00
58) Fluorene-d10	15.67	176	104512	32.07	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	17672	28.32	ng/ul	0.00
71) Anthracene-d10	17.52	188	157531	34.04	ng/ul	0.00
79) Pyrene-d10	19.80	212	171857	33.49	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	187199	38.28	ng/ul	0.00

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

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