

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052516\
 Data File : BG022063.D
 Acq On : 25 May 2016 18:00
 Operator : UM/SJ
 Sample : PB90804BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK04

Quant Time: May 26 03:12:13 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG051816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 26 03:11:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	37165	20.00	ng/ul	0.00
7) Naphthalene-d8	10.98	136	212111	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.78	164	139693	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.53	188	307181	20.00	ng/ul	0.00
29) Chrysene-d12	21.81	240	237345	20.00	ng/ul	0.00
34) Perylene-d12	25.14	264	251491	20.00	ng/ul	-0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.53	96	1643	2.03	ng/uL	0.00
3) Phenol-d5	7.32	99	73766	22.48	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.47	67	41339	23.65	ng/ul	0.00
5) 2-Chlorophenol-d4	7.69	132	64509	23.94	ng/ul	0.00
6) 4-Methylphenol-d8	8.87	113	68225	23.34	ng/ul	0.00
8) Nitrobenzene-d5	9.33	128	34331	23.07	ng/ul	0.00
9) 2-Nitrophenol-d4	10.05	143	39795	25.06	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.60	165	66209	23.59	ng/ul	0.00
12) 4-Chloroaniline-d4	11.11	131	98120	44.39	ng/ul	0.00
16) Dimethylphthalate-d6	14.18	166	260027	24.14	ng/ul	0.00
17) Acenaphthylene-d8	14.48	160	323688	23.33	ng/ul	0.00
20) 4-Nitrophenol-d4	14.99	143	33325	17.59	ng/ul	0.00
21) Fluorene-d10	15.77	176	227581	23.48	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.89	200	35699	23.42	ng/ul	0.00
26) Anthracene-d10	17.63	188	334649	23.65	ng/ul	0.00
30) Pyrene-d10	19.54	212	4087	0.33	ng/ul	-0.37
37) Benzo(a)pyrene-d12	24.91	264	242156	21.68	ng/ul	0.00

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

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