

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051816\
 Data File : BG021936.D
 Acq On : 18 May 2016 17:15
 Operator : UM/SJ
 Sample : PB90629BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK29

Quant Time: May 19 03:09:34 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG051816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 19 03:03:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	24834	20.00	ng/ul	0.00
7) Naphthalene-d8	10.98	136	131182	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.79	164	84280	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.53	188	195722	20.00	ng/ul	0.00
29) Chrysene-d12	21.81	240	171693	20.00	ng/ul	0.00
34) Perylene-d12	25.14	264	153519	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.54	96	2787	5.15	ng/uL	0.00
3) Phenol-d5	7.31	99	53949	24.60	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.48	67	32312	27.67	ng/ul	0.00
5) 2-Chlorophenol-d4	7.69	132	49241	27.35	ng/ul	0.00
6) 4-Methylphenol-d8	8.87	113	48882	25.02	ng/ul	0.00
8) Nitrobenzene-d5	9.33	128	25610	27.83	ng/ul	0.00
9) 2-Nitrophenol-d4	10.06	143	27463	27.97	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.60	165	42280	24.36	ng/ul	0.00
12) 4-Chloroaniline-d4	11.12	131	65749	48.10	ng/ul	0.00
16) Dimethylphthalate-d6	14.27	166	1271	0.20	ng/ul	0.08
17) Acenaphthylene-d8	14.48	160	235201	28.10	ng/ul	0.00
20) 4-Nitrophenol-d4	14.99	143	21378	18.70	ng/ul	0.00
21) Fluorene-d10	15.78	176	167709	28.67	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.89	200	20024	20.62	ng/ul	0.00
26) Anthracene-d10	17.53	188	195722	21.70	ng/ul	-0.09
30) Pyrene-d10	19.91	212	286590	31.64	ng/ul	0.00
37) Benzo(a)pyrene-d12	24.91	264	213290	31.28	ng/ul	0.00

Target Compounds Ovalue

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

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