

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051816\
 Data File : BG021951.D
 Acq On : 19 May 2016 3:18
 Operator : UM/SJ
 Sample : H3096-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9XG6

Quant Time: May 19 04:00:14 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG051816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 19 03:59:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	29951	20.00	ng/ul	0.00
7) Naphthalene-d8	10.98	136	154490	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.78	164	92992	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.52	188	200530	20.00	ng/ul	0.00
29) Chrysene-d12	21.81	240	186577	20.00	ng/ul	0.00
34) Perylene-d12	25.13	264	170305	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.54	96	983	1.51	ng/uL	0.00
3) Phenol-d5	7.31	99	18763	7.09	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.57	67	1334	0.95	ng/ul	0.09
5) 2-Chlorophenol-d4	7.69	132	60785	27.99	ng/ul	0.00
6) 4-Methylphenol-d8	8.86	113	36838	15.64	ng/ul	0.00
8) Nitrobenzene-d5	9.33	128	40822	37.66	ng/ul	0.00
9) 2-Nitrophenol-d4	10.05	143	42504	36.75	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.60	165	63792	31.21	ng/ul	0.00
12) 4-Chloroaniline-d4	11.14	131	1050	0.65	ng/ul	0.02
16) Dimethylphthalate-d6	14.18	166	272327	37.98	ng/ul	0.00
17) Acenaphthylene-d8	14.48	160	358002	38.76	ng/ul	0.00
20) 4-Nitrophenol-d4	15.01	143	2607	2.07	ng/ul	0.03
21) Fluorene-d10	15.77	176	248169	38.46	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.89	200	35434	35.61	ng/ul	0.00
26) Anthracene-d10	17.52	188	200530	21.70	ng/ul	-0.10
30) Pyrene-d10	19.54	212	3977	0.40	ng/ul	-0.36
37) Benzo(a)pyrene-d12	24.91	264	307950	40.72	ng/ul	0.00

Target Compounds Ovalue

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

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