

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051816\
 Data File : BG021945.D
 Acq On : 18 May 2016 23:16
 Operator : UM/SJ
 Sample : H3092-22
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9XD4

Quant Time: May 19 03:10:29 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	8.16	152	25335	20.00	ng/ul	0.00
7) Naphthalene-d8	10.98	136	141759	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.78	164	87103	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.53	188	185509	20.00	ng/ul	0.00
29) Chrysene-d12	21.81	240	157176	20.00	ng/ul	-0.01
34) Perylene-d12	25.13	264	147584	20.00	ng/ul	-0.02

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.54	96	2081	3.77	ng/uL	0.00
3) Phenol-d5	7.31	99	33244	14.86	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.48	67	20913	17.55	ng/ul	0.00
5) 2-Chlorophenol-d4	7.69	132	33370	18.17	ng/ul	0.00
6) 4-Methylphenol-d8	8.86	113	19598	9.83	ng/ul	0.00
8) Nitrobenzene-d5	9.33	128	16360	16.45	ng/ul	0.00
9) 2-Nitrophenol-d4	10.05	143	19502	18.38	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.60	165	30797	16.42	ng/ul	0.00
12) 4-Chloroaniline-d4	11.30	131	338	0.23	ng/ul	0.18
16) Dimethylphthalate-d6	14.18	166	131731	19.61	ng/ul	0.00
17) Acenaphthylene-d8	14.48	160	173970	20.11	ng/ul	0.00
20) 4-Nitrophenol-d4	15.00	143	4835	4.09	ng/ul	0.01
21) Fluorene-d10	15.77	176	118231	19.56	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.88	200	12648	13.74	ng/ul	-0.01
26) Anthracene-d10	17.53	188	185509	21.70	ng/ul	-0.10
30) Pyrene-d10	19.90	212	180522	21.77	ng/ul	0.00
37) Benzo(a)pyrene-d12	24.90	264	133564	20.38	ng/ul	0.00

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

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