

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG112816\
 Data File : BG024938.D
 Acq On : 28 Nov 2016 18:32
 Operator : UM/SJ
 Sample : H5770-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SVOC-GPC2-BLANK

Quant Time: Nov 28 23:48:20 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG112316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 28 23:45:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.25	152	82646	20.00	ng/ul	0.00
18) Naphthalene-d8	11.07	136	357113	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.87	164	300122	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.61	188	669938	20.00	ng/ul	0.00
75) Chrysene-d12	21.90	240	854444	20.00	ng/ul	0.00
83) Perylene-d12	25.33	264	872343	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.59	96	69	0.04	ng/uL	-0.01
5) Phenol-d5	7.32	99	106	0.01	ng/ul	-0.07
7) Bis-(2-Chloroethyl)ether-d	7.55	67	184	0.04	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.85	132	84	0.02	ng/ul	0.07
13) 4-Methylphenol-d8	8.91	113	61	0.01	ng/ul	-0.04
19) Nitrobenzene-d5	9.44	128	74	0.03	ng/ul	0.02
22) 2-Nitrophenol-d4	10.14	143	111	0.03	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.74	165	117	0.02	ng/ul	0.04
29) 4-Chloroaniline-d4	11.24	131	66	0.01	ng/ul	0.03
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	14.51	160	126	0.01	ng/ul	-0.06
51) 4-Nitrophenol-d4	15.07	143	57	0.01	ng/ul	0.02
57) Fluorene-d10	15.86	176	1687	0.08	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.03	200	176	0.04	ng/ul	0.06
70) Anthracene-d10	17.72	188	2872	0.09	ng/ul	0.00
76) Pyrene-d10	19.99	212	203	0.01	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.07	264	80	0.00	ng/ul	0.00

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

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