

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080619\
 Data File : BG042014.D
 Acq On : 7 Aug 2019 2:08
 Operator : HP/JU
 Sample : K4029-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BENY3

Quant Time: Aug 07 03:38:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 07 02:32:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	62965	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	265766	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.68	164	202947	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.43	188	493596	20.00	ng/ul	0.00
77) Chrysene-d12	21.72	240	507001	20.00	ng/ul	0.00
85) Perylene-d12	24.98	264	589390	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.41	96	5104	3.55	ng/uL	0.00
5) Phenol-d5	7.18	99	94159	19.06	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	54834	19.84	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	87945	20.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.73	113	68662	16.55	ng/ul	0.00
19) Nitrobenzene-d5	9.19	128	46540	23.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	56617	24.37	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	96904	20.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	121142	23.28	ng/ul	0.00
43) Dimethylphthalate-d6	14.09	166	395353	25.12	ng/ul	0.00
46) Acenaphthylene-d8	14.38	160	419313	23.32	ng/ul	0.00
51) 4-Nitrophenol-d4	14.87	143	48746	18.44	ng/ul	0.00
57) Fluorene-d10	15.68	176	343329	24.52	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.79	200	59691	18.85	ng/ul	0.00
70) Anthracene-d10	17.53	188	582037	24.58	ng/ul	0.00
78) Pyrene-d10	19.82	212	681537	24.09	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.76	264	721730	23.44	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.20	94	9271	1.818 ng/ul#	92
44) Dimethylphthalate	14.13	163	165658	10.600 ng/ul	98

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

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