

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080519\
 Data File : BG041994.D
 Acq On : 6 Aug 2019 9:28
 Operator : HP/JU
 Sample : K4028-02
 Misc : GCMS CONFIRMATION
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BEP68

Quant Time: Aug 06 10:54:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 06 04:50:21 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	143755	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	536052	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.69	164	382017	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.45	188	823750	20.00	ng/ul	0.00
77) Chrysene-d12	21.73	240	892150	20.00	ng/ul	0.00
85) Perylene-d12	25.00	264	1054085	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul	
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
78) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
89) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds

					Ovalue	
69) Phenanthrene	17.49	178	104225	2.311	ng/ul	98
76) Fluoranthene	19.50	202	204703	3.929	ng/ul	96
79) Pyrene	19.86	202	150272	2.504	ng/ul	96
82) Benzo(a)anthracene	21.71	228	118964	1.923	ng/ul	99
84) Chrysene	21.78	228	151570	2.625	ng/ul	97
87) Benzo(b)fluoranthene	23.96	252	105532	1.595	ng/ul	99

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

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