

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM061119\
 Data File : BM020887.D
 Acq On : 11 Jun 2019 22:54
 Operator : HP/JU
 Sample : K3199-10
 Misc : GCMS CONFIRMATION
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETNC3

Quant Time: Jun 12 04:41:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM060619MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jun 09 01:02:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	408046	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	1626247	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	933349	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	1754422	20.00	ng/ul	0.00
77) Chrysene-d12	21.37	240	1260080	20.00	ng/ul	0.00
85) Perylene-d12	23.65	264	1533439	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL
5) Phenol-d5	0.00	99	0	0.00	ng/ul
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0d	0.00	ng/ul
9) 2-Chlorophenol-d4	0.00	132	0	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	0d	0.00	ng/ul
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul
46) Acenaphthylene-d8	0.00	160	0	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

					Qvalue
69) Phenanthrene	17.23	178	734736	8.321 ng/ul	99
76) Fluoranthene	19.24	202	939179	9.282 ng/ul	100
79) Pyrene	19.61	202	428979	5.515 ng/ul	98
82) Benzo(a)anthracene	21.35	228	214104	2.724 ng/ul	99
84) Chrysene	21.40	228	252661	3.464 ng/ul	98
87) Benzo(b)fluoranthene	22.96	252	180612	2.025 ng/ul#	96

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

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