

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102718\
 Data File : BM017376.D
 Acq On : 27 Oct 2018 08:09
 Operator : MJ/SJ
 Sample : J5673-20
 Misc : GCMS Confirmation
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BM4

Quant Time: Oct 29 15:23:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 26 01:13:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	238566	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	1027624	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	647664	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	1587447	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	1864934	20.00	ng/ul	0.00
86) Perylene-d12	23.45	264	1845195	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	0	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
44) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul
47) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul
58) Fluorene-d10	0.00	176	0d	0.00	ng/ul
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul
79) Pyrene-d10	0.00	212	0d	0.00	ng/ul
90) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

Target Compounds

					Ovalue
37) 1,2,4,5-Tetrachlorobenzene	12.47	216	213457	9.806 ng/ul	97
70) Phenanthrene	17.09	178	116092	1.274 ng/ul#	51
73) 1,2,3,4-Tetrachlorobenzene	13.08	216	247100	11.040 ng/uL	98

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

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