

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102718\
 Data File : BM017380.D
 Acq On : 27 Oct 2018 10:34
 Operator : MJ/SJ
 Sample : J5673-10
 Misc : GCMS Confirmation
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BL4

Quant Time: Oct 29 02:13:19 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	326994	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	1439564	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	829470	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	1774087	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	1890327	20.00	ng/ul	0.00
86) Perylene-d12	23.46	264	1752852	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	0d	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	12.46	216	549937	19.727 ng/ul 99
73) 1,2,3,4-Tetrachlorobenzene	13.08	216	639646	25.572 ng/uL 98
74) Pentachlorobenzene	14.62	250	36631	1.423 ng/uL 95

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

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