

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102718\
 Data File : BM017387.D
 Acq On : 27 Oct 2018 14:46
 Operator : MJ/SJ
 Sample : J5674-09
 Misc : GCMS Confirmation
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BN2

Quant Time: Oct 29 01:14:03 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	256091	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	1071529	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	673281	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	1571208	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	1859004	20.00	ng/ul	0.00
86) Perylene-d12	23.45	264	1944059	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.23	96	241	0.04	ng/uL	0.01
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.04	67	410	0.03	ng/ul	0.04
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	10.45	131	140	0.01	ng/ul	-0.13
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	14.57	143	146	0.01	ng/ul	0.05
58) Fluorene-d10	15.30	176	3938	0.08	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.42	200	159	0.01	ng/ul	0.00
71) Anthracene-d10	17.04	188	1571208	20.20	ng/ul	-0.10
79) Pyrene-d10	19.46	212	403	0.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.45	264	1892367	18.43	ng/ul	0.14

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

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