

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM103118\
 Data File : BM017429.D
 Acq On : 31 Oct 2018 14:01
 Operator : MJ/SJ
 Sample : J5674-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COBN0

Quant Time: Oct 31 15:15:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 31 14:01:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	216632	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	945550	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.29	164	562556	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.04	188	1269936	20.00	ng/ul	-0.01
78) Chrysene-d12	21.24	240	1600766	20.00	ng/ul	0.00
86) Perylene-d12	23.45	264	1366249	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	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	0	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0	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	15.42	200	72	0.01	ng/ul	-0.01
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	
70) Phenanthrene	17.09	178	78728	1.080	ng/ul# 79
73) 1,2,3,4-Tetrachlorobenzene	13.08	216	46671	2.607	ng/uL 99
74) Pentachlorobenzene	14.61	250	41289	2.241	ng/uL 97
77) Fluoranthene	19.11	202	216322	2.588	ng/ul# 97
80) Pyrene	19.47	202	139418	1.500	ng/ul# 91
85) Chrysene	21.28	228	165097	1.772	ng/ul# 92

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

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