

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061119\
 Data File : BM020893.D
 Acq On : 12 Jun 2019 02:31
 Operator : HP/JU
 Sample : K3199-18
 Misc : GCMS CONFIRMATION
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jun 12 04:04:26 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	313901	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	1257538	20.00	ng/ul	-0.01
35) Acenaphthene-d10	14.44	164	776011	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	1813653	20.00	ng/ul	0.00
77) Chrysene-d12	21.36	240	1441642	20.00	ng/ul	-0.01
85) Perylene-d12	23.64	264	1545469	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.30	96	176	0.03	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.09	67	244	0.02	ng/ul	-0.06
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	10.60	131	1869	0.08	ng/ul	-0.14
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	15.43	176	3477	0.07	ng/ul	-0.01
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.29	188	5821	0.07	ng/ul	0.00
78) Pyrene-d10	19.58	212	384	0.01	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.64	264	1525285	21.71	ng/ul	0.14

Target Compounds	Ovalue

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

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