

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112119\
 Data File : BM023849.D
 Acq On : 22 Nov 2019 03:44
 Operator : JU
 Sample : K5809-22
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFPW3

Quant Time: Nov 22 07:40:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 21 16:09:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	429554	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	1619896	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	946269	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1886666	20.00	ng/ul	0.00
78) Chrysene-d12	21.16	240	1646604	20.00	ng/ul	0.02
86) Perylene-d12	23.34	264	2342370	20.00	ng/ul	0.05

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL
5) Phenol-d5	0.00	99	0d	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	0d	0.00	ng/ul
19) Nitrobenzene-d5	0.00	128	0d	0.00	ng/ul
22) 2-Nitrophenol-d4	0.00	143	0d	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	
28) Naphthalene	10.34	128	190317	2.235	ng/ul# 90
34) 2-Methylnaphthalene	11.97	142	244948	3.761	ng/ul 97
70) Phenanthrene	16.97	178	445220	4.284	ng/ul# 83
73) 1,2,3,4-Tetrachlorobenzene	12.96	216	51318	1.919	ng/uL 98
77) Fluoranthene	19.02	202	223763	2.010	ng/ul# 84

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

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