

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112119\
 Data File : BM023855.D
 Acq On : 22 Nov 2019 07:20
 Operator : JU
 Sample : K5809-18
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 22 08:09:47 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	414567	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	1560464	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	850985	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1519901	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	1674210	20.00	ng/ul	0.01
86) Perylene-d12	23.32	264	2236417	20.00	ng/ul	0.03

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	376	0.04	ng/uL	0.00
5) Phenol-d5	6.71	99	149	0.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.85	67	2555	0.12	ng/ul	-0.02
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.24	113	681	0.02	ng/ul	0.00
19) Nitrobenzene-d5	8.75	128	368	0.03	ng/ul	0.08
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	9.93	165	405	0.02	ng/ul	0.00
29) 4-Chloroaniline-d4	10.42	131	1758	0.06	ng/ul	-0.02
44) Dimethylphthalate-d6	13.60	166	115	0.00	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	373	0.00	ng/ul	0.01
52) 4-Nitrophenol-d4	14.42	143	1767	0.16	ng/ul	0.01
58) Fluorene-d10	15.16	176	194	0.00	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.32	200	452	0.05	ng/ul	0.00
71) Anthracene-d10	16.93	188	1519901	21.08	ng/ul	-0.09
79) Pyrene-d10	19.34	212	2352	0.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.32	264	2230202	18.95	ng/ul	0.16

Target Compounds Ovalue

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

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