

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112119\
 Data File : BM023853.D
 Acq On : 22 Nov 2019 06:08
 Operator : JU
 Sample : K5809-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 22 08:09:34 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	7.51	152	498560	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.29	136	1922567	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	1053894	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1943479	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	2131691	20.00	ng/ul	0.01
86) Perylene-d12	23.32	264	2821905	20.00	ng/ul	0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	228	0.02	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	6.87	67	442	0.02	ng/ul	0.00
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.06	113	1172	0.03	ng/ul	-0.18
19) Nitrobenzene-d5	8.63	128	107	0.01	ng/ul	-0.04
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	9.94	165	440	0.01	ng/ul	0.01
29) 4-Chloroaniline-d4	10.42	131	182	0.01	ng/ul	-0.02
44) Dimethylphthalate-d6	13.59	166	810	0.01	ng/ul	-0.01
47) Acenaphthylene-d8	13.86	160	79	0.00	ng/ul	0.00
52) 4-Nitrophenol-d4	14.39	143	1225	0.09	ng/ul	-0.02
58) Fluorene-d10	15.16	176	288	0.00	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.32	200	114	0.01	ng/ul	0.00
71) Anthracene-d10	16.93	188	1943479	21.08	ng/ul	-0.09
79) Pyrene-d10	19.33	212	2703	0.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.26	264	3339	0.02	ng/ul	0.11

Target Compounds	Ovalue

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

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