

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111919\
 Data File : BM023821.D
 Acq On : 20 Nov 2019 06:19
 Operator : JU
 Sample : K5847-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 20 10:23:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 06:55:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	466859	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	1800062	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	966990	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1679446	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	1492975	20.00	ng/ul	0.00
86) Perylene-d12	23.30	264	2005584	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.05	96	191	0.02	ng/uL	0.00
5) Phenol-d5	6.53	99	615	0.01	ng/ul	-0.18
7) Bis-(2-Chloroethyl)ether-d	6.86	67	815	0.03	ng/ul	-0.01
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.20	113	739	0.02	ng/ul	-0.04
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	9.95	165	314	0.01	ng/ul	0.02
29) 4-Chloroaniline-d4	10.42	131	150	0.00	ng/ul	-0.02
44) Dimethylphthalate-d6	13.58	166	670	0.01	ng/ul	-0.02
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	14.40	143	313	0.02	ng/ul	0.00
58) Fluorene-d10	15.17	176	167	0.00	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	362	0.04	ng/ul	0.00
71) Anthracene-d10	16.93	188	1679446	21.08	ng/ul	-0.10
79) Pyrene-d10	19.33	212	3170	0.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	1992269	18.88	ng/ul	0.14

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

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