

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102419\
 Data File : BM023390.D
 Acq On : 24 Oct 2019 21:40
 Operator : JU
 Sample : K5449-05RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 24 22:50:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 23 18:47:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	745364	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	2789864	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	1456315	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	2612159	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	2426073	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	3073392	20.00	ng/ul	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	413	0.03	ng/uL	-0.01
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	6.97	67	157	0.00	ng/ul	-0.01
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	8.76	128	272	0.01	ng/ul	-0.02
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	10.55	131	468	0.01	ng/ul	0.00
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.29	176	3430	0.04	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.04	188	2612159	21.16	ng/ul	-0.09
79) Pyrene-d10	19.43	212	2957	0.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.44	264	3062964	19.56	ng/ul	0.15

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

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