

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101519\
 Data File : BM023182.D
 Acq On : 16 Oct 2019 00:33
 Operator : JU
 Sample : K5349-10
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 16 06:50:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 02:28:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	358796	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	1557531	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	1044280	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	2235009	20.00	ng/ul	-0.01
78) Chrysene-d12	21.26	240	2034746	20.00	ng/ul	-0.01
86) Perylene-d12	23.47	264	1843380	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.20	96	445	0.05	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.00	67	293	0.02	ng/ul	-0.02
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	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	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	10.47	131	1685	0.05	ng/ul	-0.13
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	15.33	176	2252	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.07	188	2235009	20.74	ng/ul	-0.11
79) Pyrene-d10	19.46	212	486	0.00	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.47	264	1838817	19.68	ng/ul	0.12

Target Compounds	Ovalue

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

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