

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
 Data File : BM023382.D
 Acq On : 24 Oct 2019 16:45
 Operator : JU
 Sample : K5482-01
 Misc : GCMS CONFIRMATION
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 24 17:37:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 24 13:46:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	582113	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.41	136	2204759	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.28	164	1286604	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	2509494	20.00	ng/ul	-0.01
78) Chrysene-d12	21.22	240	2607361	20.00	ng/ul	-0.01
86) Perylene-d12	23.41	264	3130295	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	358	0.03	ng/uL	-0.01
5) Phenol-d5	6.71	99	908	0.02	ng/ul	-0.11
7) Bis-(2-Chloroethyl)ether-d	6.96	67	488	0.02	ng/ul	-0.02
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.30	113	3411	0.09	ng/ul	-0.05
19) Nitrobenzene-d5	8.87	128	375	0.02	ng/ul	0.09
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	10.54	131	2312	0.06	ng/ul	-0.02
44) Dimethylphthalate-d6	13.69	166	305	0.00	ng/ul	-0.01
47) Acenaphthylene-d8	13.96	160	445	0.00	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.49	143	124	0.01	ng/ul	0.00
58) Fluorene-d10	15.27	176	2976	0.04	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.39	200	1293	0.08	ng/ul	-0.02
71) Anthracene-d10	17.02	188	2509494	21.16	ng/ul	-0.11
79) Pyrene-d10	19.43	212	6885	0.05	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.41	264	3099704	19.44	ng/ul	0.12

Target Compounds	Ovalue

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

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