

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102519\
 Data File : BM023410.D
 Acq On : 25 Oct 2019 17:08
 Operator : JU
 Sample : PB124099BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK99

Quant Time: Oct 25 22:11:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 25 07:44:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	493576	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	2114990	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.28	164	1333953	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	2642540	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	1993849	20.00	ng/ul	0.00
86) Perylene-d12	23.42	264	2057073	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.16	96	78801	7.58	ng/uL	0.00
5) Phenol-d5	6.82	99	1439833	34.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.97	67	864238	35.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	1186959	36.63	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	1166864	34.66	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	602068	37.81	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	671377	38.00	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.04	165	1216424	35.04	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	7780	0.19	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	3807820	37.59	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	4395188	37.87	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.49	143	607910	34.64	ng/ul	0.00
58) Fluorene-d10	15.28	176	3199140	37.29	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.41	200	437170	26.57	ng/ul	0.00
71) Anthracene-d10	17.13	188	4752319	38.05	ng/ul	0.00
79) Pyrene-d10	19.43	212	4581329	41.78	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.29	264	3999439	38.17	ng/ul	0.00

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

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