

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023461.D
 Acq On : 30 Oct 2019 02:06
 Operator : JU
 Sample : PB124083BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK83

Quant Time: Oct 30 07:12:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 07:08:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1044407	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.40	136	4346119	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.27	164	2613191	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.02	188	5242446	20.00	ng/ul	-0.02
78) Chrysene-d12	21.21	240	5393468	20.00	ng/ul	-0.02
86) Perylene-d12	23.40	264	6438140	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	65007	2.96	ng/uL	0.00
5) Phenol-d5	6.80	99	1097214	12.51	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.96	67	690752	13.44	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.16	132	890304	12.98	ng/ul	-0.02
13) 4-Methylphenol-d8	8.33	113	844322	11.85	ng/ul	-0.02
19) Nitrobenzene-d5	8.77	128	420444	12.85	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.49	143	452995	12.48	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.03	165	813840	11.41	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.53	131	1065592	12.99	ng/ul	-0.02
44) Dimethylphthalate-d6	13.69	166	2408466	12.14	ng/ul	-0.02
47) Acenaphthylene-d8	13.96	160	2851505	12.54	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.48	143	339793	9.88	ng/ul	-0.01
58) Fluorene-d10	15.27	176	2058310	12.25	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.39	200	254355	7.79	ng/ul	-0.02
71) Anthracene-d10	17.12	188	3163670	12.77	ng/ul	-0.02
79) Pyrene-d10	19.42	212	3409631	11.50	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.26	264	4013949	12.24	ng/ul	-0.02

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

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