

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
 Data File : BM022910.D
 Acq On : 03 Oct 2019 01:39
 Operator : JU
 Sample : PB123344TB
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB17

Quant Time: Oct 03 02:59:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 27 18:03:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	230635	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1033544	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.43	164	658839	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1222033	20.00	ng/ul	-0.01
78) Chrysene-d12	21.34	240	839720	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	891510	20.00	ng/ul	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.27	96	35933	6.74	ng/uL	0.00
5) Phenol-d5	6.97	99	657219	32.25	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	381115	34.17	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.33	132	544096	33.16	ng/ul	-0.01
13) 4-Methylphenol-d8	8.50	113	547767	32.83	ng/ul	-0.01
19) Nitrobenzene-d5	8.95	128	275877	34.59	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	308445	35.41	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.21	165	565071	32.58	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	813775	38.63	ng/ul	-0.01
44) Dimethylphthalate-d6	13.84	166	1797829	35.16	ng/ul	-0.01
47) Acenaphthylene-d8	14.13	160	2027122	34.34	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.63	143	259703	25.70	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1453657	33.36	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.54	200	155471	19.55	ng/ul	-0.02
71) Anthracene-d10	17.27	188	2094352	35.10	ng/ul	-0.02
79) Pyrene-d10	19.56	212	1926664	42.63	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.46	264	1650671	35.79	ng/ul	-0.04

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

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