

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
 Data File : BM022899.D
 Acq On : 02 Oct 2019 18:58
 Operator : JU
 Sample : PB123401TB
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB19

Quant Time: Oct 03 00:59:40 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	241971	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1100160	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.43	164	710948	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1370244	20.00	ng/ul	-0.01
78) Chrysene-d12	21.36	240	936593	20.00	ng/ul	-0.01
86) Perylene-d12	23.61	264	944180	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.27	96	37383	6.68	ng/uL	0.00
5) Phenol-d5	6.97	99	667837	31.24	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	386667	33.05	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.33	132	547859	31.82	ng/ul	-0.01
13) 4-Methylphenol-d8	8.50	113	557808	31.87	ng/ul	-0.01
19) Nitrobenzene-d5	8.95	128	279093	32.87	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	311231	33.57	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.21	165	577037	31.26	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	841323	37.52	ng/ul	-0.01
44) Dimethylphthalate-d6	13.84	166	1828077	33.13	ng/ul	-0.01
47) Acenaphthylene-d8	14.13	160	2101716	32.99	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.63	143	281853	25.85	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1507250	32.05	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.54	200	177301	19.88	ng/ul	-0.02
71) Anthracene-d10	17.27	188	2197066	32.83	ng/ul	-0.01
79) Pyrene-d10	19.56	212	2094717	41.55	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1639793	33.57	ng/ul	-0.02

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

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