

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM022917.D
 Acq On : 03 Oct 2019 13:12
 Operator : JU
 Sample : PB123421BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK21

Quant Time: Oct 04 01:21:23 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	218537	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	927285	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.43	164	576090	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1090315	20.00	ng/ul	-0.02
78) Chrysene-d12	21.35	240	834143	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	870290	20.00	ng/ul	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.26	96	31796	6.29	ng/uL	0.00
5) Phenol-d5	6.96	99	545994	28.28	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.13	67	320103	30.29	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.33	132	463399	29.80	ng/ul	-0.02
13) 4-Methylphenol-d8	8.50	113	441113	27.90	ng/ul	-0.02
19) Nitrobenzene-d5	8.95	128	224924	31.43	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	243034	31.10	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.21	165	448044	28.80	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	645514	34.15	ng/ul	-0.01
44) Dimethylphthalate-d6	13.84	166	1393986	31.18	ng/ul	-0.02
47) Acenaphthylene-d8	14.12	160	1609604	31.18	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.63	143	203348	23.02	ng/ul	-0.01
58) Fluorene-d10	15.42	176	1158598	30.40	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.54	200	132632	18.69	ng/ul	-0.02
71) Anthracene-d10	17.27	188	1699981	31.93	ng/ul	-0.02
79) Pyrene-d10	19.56	212	1648294	36.71	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1437135	31.92	ng/ul	-0.03

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

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