

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100119\
 Data File : BM022912.D
 Acq On : 03 Oct 2019 02:52
 Operator : JU
 Sample : PB123794TB
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 03 03:34:21 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	222840	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1009132	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.43	164	632876	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1153533	20.00	ng/ul	-0.01
78) Chrysene-d12	21.35	240	799454	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	862824	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	34998	6.79	ng/uL	0.00
5) Phenol-d5	6.96	99	627196	31.86	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.13	67	360748	33.48	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.33	132	511525	32.26	ng/ul	-0.01
13) 4-Methylphenol-d8	8.50	113	516142	32.02	ng/ul	-0.01
19) Nitrobenzene-d5	8.95	128	258703	33.22	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	286111	33.64	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.21	165	534902	31.59	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	767019	37.29	ng/ul	-0.01
44) Dimethylphthalate-d6	13.85	166	1671611	34.03	ng/ul	-0.01
47) Acenaphthylene-d8	14.13	160	1898673	33.48	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.63	143	237312	24.45	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1350750	32.27	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.54	200	146591	19.53	ng/ul	-0.02
71) Anthracene-d10	17.27	188	1938865	34.42	ng/ul	-0.02
79) Pyrene-d10	19.56	212	1794602	41.71	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1584333	35.49	ng/ul	-0.03

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

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