

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100819\
 Data File : BM023118.D
 Acq On : 10 Oct 2019 16:37
 Operator : JU
 Sample : SP4934
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP4934

Quant Time: Oct 10 17:15:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 08 15:02:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	202140	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	809579	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.40	164	430178	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.14	188	751732	20.00	ng/ul	0.00
78) Chrysene-d12	21.33	240	709850	20.00	ng/ul	0.00
86) Perylene-d12	23.59	264	838761	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.25	96	68615	14.68	ng/uL	0.00
5) Phenol-d5	6.94	99	1129889	63.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	658454	67.36	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	957163	66.55	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	896482	61.30	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	466929	74.73	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	524418	76.87	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.18	165	897771	66.09	ng/ul	0.00
29) 4-Chloroaniline-d4	10.69	131	1218636	73.84	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	2244504	67.23	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	2722215	70.62	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	354669	53.76	ng/ul	0.00
58) Fluorene-d10	15.40	176	1825628	64.16	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	287959	58.86	ng/ul	0.00
71) Anthracene-d10	17.24	188	2514734	68.50	ng/ul	0.00
79) Pyrene-d10	19.54	212	2605640	68.20	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.44	264	3039673	70.05	ng/ul	0.00

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

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