

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM022919.D
 Acq On : 03 Oct 2019 14:24
 Operator : JU
 Sample : PB123379BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK79

Quant Time: Oct 04 01:21:31 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	244563	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1033863	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.43	164	613968	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1136316	20.00	ng/ul	-0.02
78) Chrysene-d12	21.35	240	906029	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	1008279	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	37446	6.62	ng/uL	0.00
5) Phenol-d5	6.96	99	642627	29.74	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.13	67	384797	32.54	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.33	132	542761	31.19	ng/ul	-0.02
13) 4-Methylphenol-d8	8.50	113	522472	29.53	ng/ul	-0.02
19) Nitrobenzene-d5	8.95	128	264028	33.09	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	285669	32.79	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.20	165	522822	30.14	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	748478	35.52	ng/ul	-0.01
44) Dimethylphthalate-d6	13.84	166	1586121	33.29	ng/ul	-0.02
47) Acenaphthylene-d8	14.12	160	1835908	33.37	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.62	143	245582	26.08	ng/ul	-0.02
58) Fluorene-d10	15.42	176	1307115	32.19	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.54	200	160497	21.70	ng/ul	-0.02
71) Anthracene-d10	17.27	188	1903236	34.30	ng/ul	-0.02
79) Pyrene-d10	19.56	212	1895714	38.88	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1847757	35.42	ng/ul	-0.03

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

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