

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
 Data File : BM022900.D
 Acq On : 02 Oct 2019 19:34
 Operator : JU
 Sample : PB123471BL
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK71

Quant Time: Oct 03 00:59:44 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	231715	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1025443	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.43	164	631742	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1117619	20.00	ng/ul	-0.01
78) Chrysene-d12	21.35	240	778720	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	888626	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	32968	6.15	ng/uL	0.00
5) Phenol-d5	6.97	99	581473	28.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	339719	30.32	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.33	132	480232	29.13	ng/ul	-0.01
13) 4-Methylphenol-d8	8.50	113	481559	28.73	ng/ul	-0.01
19) Nitrobenzene-d5	8.95	128	236021	29.82	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	262017	30.32	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.21	165	491656	28.57	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	702400	33.60	ng/ul	-0.01
44) Dimethylphthalate-d6	13.85	166	1475829	30.10	ng/ul	-0.01
47) Acenaphthylene-d8	14.13	160	1708238	30.18	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.63	143	198948	20.54	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1195774	28.62	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.54	200	126982	17.46	ng/ul	-0.02
71) Anthracene-d10	17.27	188	1661075	30.44	ng/ul	-0.01
79) Pyrene-d10	19.56	212	1483898	35.41	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1423866	30.97	ng/ul	-0.02

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

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