

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM093019\
 Data File : BM022839.D
 Acq On : 30 Sep 2019 23:34
 Operator : JU
 Sample : PB123342BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK42

Quant Time: Oct 01 00:55:29 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	260138	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	1180234	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.44	164	762679	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.18	188	1609591	20.00	ng/ul	0.00
78) Chrysene-d12	21.36	240	1542134	20.00	ng/ul	-0.01
86) Perylene-d12	23.62	264	1361015	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	38303	6.37	ng/uL	0.00
5) Phenol-d5	6.97	99	665306	28.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	404143	32.13	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	568392	30.71	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	553239	29.40	ng/ul	-0.01
19) Nitrobenzene-d5	8.96	128	280928	30.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	307737	30.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	566708	28.62	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.73	131	843503	35.06	ng/ul	0.00
44) Dimethylphthalate-d6	13.85	166	1835921	31.02	ng/ul	0.00
47) Acenaphthylene-d8	14.13	160	2133609	31.22	ng/ul	0.00
52) 4-Nitrophenol-d4	14.63	143	326372	27.91	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1574566	31.21	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.55	200	245662	23.45	ng/ul	-0.01
71) Anthracene-d10	17.28	188	2524775	32.12	ng/ul	0.00
79) Pyrene-d10	19.57	212	2812987	33.89	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.48	264	2284725	32.45	ng/ul	-0.02

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

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