

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
 Data File : BM022885.D
 Acq On : 02 Oct 2019 10:25
 Operator : JU
 Sample : PB123403BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK03

Quant Time: Oct 02 11:38:11 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	291183	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1328273	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.44	164	851369	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.17	188	1619630	20.00	ng/ul	-0.01
78) Chrysene-d12	21.36	240	1134392	20.00	ng/ul	-0.01
86) Perylene-d12	23.62	264	1165833	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	41148	6.11	ng/uL	0.00
5) Phenol-d5	6.97	99	740568	28.79	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	433708	30.80	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.33	132	611078	29.49	ng/ul	-0.01
13) 4-Methylphenol-d8	8.50	113	619946	29.43	ng/ul	-0.01
19) Nitrobenzene-d5	8.96	128	314257	30.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	355538	31.76	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.21	165	639069	28.67	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	920963	34.01	ng/ul	-0.01
44) Dimethylphthalate-d6	13.85	166	2027844	30.69	ng/ul	-0.01
47) Acenaphthylene-d8	14.13	160	2312708	30.32	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.63	143	322105	24.67	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1653307	29.36	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.54	200	218180	20.70	ng/ul	-0.02
71) Anthracene-d10	17.27	188	2385718	30.16	ng/ul	-0.01
79) Pyrene-d10	19.57	212	2262471	37.06	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.47	264	1866228	30.94	ng/ul	-0.02

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

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