

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM022979.D
 Acq On : 05 Oct 2019 04:04
 Operator : JU
 Sample : PB123622BL
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK22

Quant Time: Oct 05 04:41:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 16:45:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	255496	20.00	ng/ul	0.00
18) Naphthalene-d8	10.58	136	1074580	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	600242	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.16	188	1014271	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	868019	20.00	ng/ul	-0.01
86) Perylene-d12	23.60	264	1056991	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.26	96	37912	6.42	ng/uL	0.00
5) Phenol-d5	6.96	99	649055	28.75	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	377238	30.53	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	539761	29.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	517988	28.02	ng/ul	-0.01
19) Nitrobenzene-d5	8.95	128	260318	31.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	277714	30.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	517979	28.73	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.71	131	715624	32.67	ng/ul	0.00
44) Dimethylphthalate-d6	13.83	166	1413905	30.35	ng/ul	0.00
47) Acenaphthylene-d8	14.12	160	1664848	30.95	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	182793	19.86	ng/ul	0.00
58) Fluorene-d10	15.42	176	1134024	28.56	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	124621	18.88	ng/ul	0.00
71) Anthracene-d10	17.26	188	1576639	31.83	ng/ul	0.00
79) Pyrene-d10	19.56	212	1559833	33.39	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.46	264	1762830	32.24	ng/ul	-0.01

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

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