

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
 Data File : BM023363.D
 Acq On : 24 Oct 2019 03:57
 Operator : JU
 Sample : PB123972BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK72

Quant Time: Oct 24 04:32:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 23 18:47:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	487494	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	2227342	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	1518053	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	3148447	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	2556174	20.00	ng/ul	0.00
86) Perylene-d12	23.41	264	2323400	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.16	96	65710	6.40	ng/uL	0.00
5) Phenol-d5	6.82	99	1285793	31.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.97	67	818466	34.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	1028106	32.12	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	1088276	32.73	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	552257	32.94	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	617275	33.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	1098397	30.04	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	1602171	38.11	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	3966551	34.41	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	4362610	33.03	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	625050	31.30	ng/ul	0.00
58) Fluorene-d10	15.28	176	3270358	33.50	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	490906	25.04	ng/ul	0.00
71) Anthracene-d10	17.13	188	5217692	35.07	ng/ul	0.00
79) Pyrene-d10	19.43	212	5354918	38.09	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.28	264	4224593	35.69	ng/ul	0.00

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

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