

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101919\
 Data File : BM023309.D
 Acq On : 20 Oct 2019 04:49
 Operator : JU
 Sample : PB124101BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK01

Quant Time: Oct 21 01:55:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 19 23:34:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	447473	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	1783147	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	1057010	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	2117384	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	2195876	20.00	ng/ul	0.00
86) Perylene-d12	23.43	264	2579167	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.17	96	66602	6.55	ng/uL	0.00
5) Phenol-d5	6.82	99	1092501	28.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	627702	28.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	897151	30.08	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	846075	27.03	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	411753	34.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	459704	40.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	862859	29.54	ng/ul	0.00
29) 4-Chloroaniline-d4	10.57	131	59014	1.66	ng/ul	0.00
44) Dimethylphthalate-d6	13.71	166	2459005	30.36	ng/ul	0.00
47) Acenaphthylene-d8	13.99	160	2927597	30.81	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	377891	28.82	ng/ul	0.00
58) Fluorene-d10	15.29	176	2113254	30.62	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.41	200	284535	33.89	ng/ul	0.00
71) Anthracene-d10	17.14	188	3233446	31.67	ng/ul	0.00
79) Pyrene-d10	19.43	212	3524959	33.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.29	264	4195741	32.09	ng/ul	0.00

Target Compounds	Ovalue

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

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