

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111119\
 Data File : BM023652.D
 Acq On : 12 Nov 2019 00:28
 Operator : JU
 Sample : PB124323TB
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB32

Quant Time: Nov 12 01:29:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 12 00:02:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	367641	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	1689233	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.20	164	1239747	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.95	188	2837403	20.00	ng/ul	0.00
78) Chrysene-d12	21.15	240	2562096	20.00	ng/ul	0.00
86) Perylene-d12	23.31	264	2341056	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.08	96	15178	1.70	ng/uL	0.00
5) Phenol-d5	6.73	99	274075	8.18	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.89	67	658098	34.18	ng/ul	0.00
9) 2-Chlorophenol-d4	7.08	132	703489	28.98	ng/ul	0.00
13) 4-Methylphenol-d8	8.26	113	516199	19.31	ng/ul	0.00
19) Nitrobenzene-d5	8.70	128	448531	37.09	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	511284	40.23	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.95	165	962663	33.22	ng/ul	0.00
29) 4-Chloroaniline-d4	10.48	131	2728	0.09	ng/ul	0.02
44) Dimethylphthalate-d6	13.62	166	3579911	37.03	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	3952516	35.82	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	106949	6.48	ng/ul	0.00
58) Fluorene-d10	15.20	176	3140182	36.78	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.33	200	508216	30.96	ng/ul	0.00
71) Anthracene-d10	17.05	188	5062750	37.61	ng/ul	0.00
79) Pyrene-d10	19.35	212	5644684	42.25	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.18	264	4749054	38.55	ng/ul	0.00

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

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