

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023458.D
 Acq On : 30 Oct 2019 00:19
 Operator : JU
 Sample : K5426-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNJ7

Quant Time: Oct 30 07:12:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 07:08:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	580803	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.40	136	2392379	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.27	164	1447206	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.02	188	2966554	20.00	ng/ul	-0.02
78) Chrysene-d12	21.21	240	3141535	20.00	ng/ul	-0.02
86) Perylene-d12	23.40	264	3748929	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	48196	3.94	ng/uL	0.00
5) Phenol-d5	6.80	99	183331	3.76	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.96	67	455873	15.94	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.16	132	499025	13.09	ng/ul	-0.02
13) 4-Methylphenol-d8	8.33	113	303984	7.67	ng/ul	-0.02
19) Nitrobenzene-d5	8.77	128	295107	16.39	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.49	143	304966	15.26	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.03	165	516355	13.15	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.54	131	40334	0.89	ng/ul	-0.02
44) Dimethylphthalate-d6	13.69	166	1870137	17.02	ng/ul	-0.02
47) Acenaphthylene-d8	13.96	160	2104397	16.72	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.49	143	47375	2.49	ng/ul	0.00
58) Fluorene-d10	15.27	176	1659202	17.83	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.39	200	228288	12.36	ng/ul	-0.02
71) Anthracene-d10	17.12	188	2761888	19.70	ng/ul	-0.02
79) Pyrene-d10	19.42	212	3117445	18.04	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.26	264	3760658	19.69	ng/ul	-0.02

Target Compounds	Ovalue

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

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