

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100819\
 Data File : BM023092.D
 Acq On : 09 Oct 2019 20:20
 Operator : JU
 Sample : K5096-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLMJ5

Quant Time: Oct 09 23:32:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 08 15:02:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	155201	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	642243	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.40	164	381687	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.14	188	793921	20.00	ng/ul	0.00
78) Chrysene-d12	21.33	240	831671	20.00	ng/ul	0.00
86) Perylene-d12	23.58	264	962270	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.25	96	15663	4.36	ng/uL	0.00
5) Phenol-d5	6.94	99	36034	2.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	93970	12.52	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	111491	10.10	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	67094	5.98	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	62917	12.69	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	66031	12.20	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.18	165	128337	11.91	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	7281	0.56	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	419721	14.17	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	477317	13.96	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	9937	1.70	ng/ul	0.00
58) Fluorene-d10	15.40	176	366241	14.51	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	44219	8.56	ng/ul	0.00
71) Anthracene-d10	17.24	188	576369	14.87	ng/ul	0.00
79) Pyrene-d10	19.54	212	675769	15.10	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.43	264	746574	15.00	ng/ul	-0.01

Target Compounds					Ovalue	
6) Phenol	6.96	94	23552	1.638	ng/ul	97
16) 4-Methylphenol	8.53	108	20192	1.648	ng/ul	86
45) Dimethylphthalate	13.86	163	107949	3.547	ng/ul	99

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

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