

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN091719\
 Data File : BN007695.D
 Acq On : 17 Sep 2019 11:58
 Operator : HP/JU
 Sample : K4899-10
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 18 01:33:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 16 18:09:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	545609	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	2314413	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.33	164	1464495	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	3100263	20.00	ng/ul	0.00
77) Chrysene-d12	21.27	240	3129712	20.00	ng/ul	0.00
85) Perylene-d12	23.49	264	3260490	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	294	0.02	ng/uL	0.01
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.02	67	312	0.01	ng/ul	-0.03
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	10.48	131	3278	0.07	ng/ul	-0.12
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	14.02	160	366	0.00	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.32	176	3090	0.03	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.17	188	14623	0.09	ng/ul	0.00
78) Pyrene-d10	19.47	212	5256	0.03	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.36	264	13656	0.08	ng/ul	0.00

Target Compounds Ovalue

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

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