

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\
 Data File : BM024587.D
 Acq On : 22 Jan 2020 18:22
 Operator : JU
 Sample : L1118-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GASK1

Quant Time: Jan 23 01:33:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 23 01:22:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	235158	20.00	ng/ul	0.00
18) Naphthalene-d8	10.21	136	1057253	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	746477	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1685009	20.00	ng/ul	0.00
78) Chrysene-d12	21.06	240	1519366	20.00	ng/ul	0.00
86) Perylene-d12	23.17	264	1395079	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
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	10.36	131	366	0.02	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	178	0.00	ng/ul	0.00
47) Acenaphthylene-d8	13.78	160	977	0.01	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.09	176	324	0.01	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
79) Pyrene-d10	19.25	212	2657	0.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.04	264	2166	0.03	ng/ul	0.00

Target Compounds Ovalue

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

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