

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN092319\
 Data File : BN007828.D
 Acq On : 24 Sep 2019 07:40
 Operator : HP/JU
 Sample : K4977-06
 Misc : GCMS CONFIRMATION
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 24 08:16:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091619MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 05:34:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	616472	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	2348926	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	1433924	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	3014761	20.00	ng/ul	0.00
77) Chrysene-d12	21.26	240	2775624	20.00	ng/ul	0.00
85) Perylene-d12	23.47	264	2833927	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	1312	0.08	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.03	67	415	0.01	ng/ul	0.00
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	10.46	131	3273	0.07	ng/ul	-0.13
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	14.03	160	417	0.00	ng/ul	0.02
51) 4-Nitrophenol-d4	14.65	143	1312	0.06	ng/ul	0.14
57) Fluorene-d10	15.31	176	3044	0.03	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.16	188	13828	0.09	ng/ul	0.00
78) Pyrene-d10	19.46	212	5082	0.03	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.47	264	2800078	18.99	ng/ul	0.14

Target Compounds Ovalue

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

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