

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030819\
 Data File : BN004706.D
 Acq On : 09 Mar 2019 11:31
 Operator : JU/SJ
 Sample : K1762-05
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0959

Quant Time: Mar 11 01:42:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN030719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 08 00:36:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	32185	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	118582	20.00	ng/ul	0.00
36) Acenaphthene-d10	0.00	164	0m	20.00	ng/ul	-14.36
62) Phenanthrene-d10	17.11	188	42216	20.00	ng/ul	0.00
78) Chrysene-d12	21.31	240	4717	20.00	ng/ul	0.00
86) Perylene-d12	0.00	264	0	0.00	ng/ul	-23.57

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	2457	4.31	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.06	67	38528	27.77	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	4379	1.90	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	8.88	128	34910	38.62	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	11764	11.46	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	15363	7.85	ng/ul	0.01
29) 4-Chloroaniline-d4	10.65	131	4672	1.99	ng/ul	0.00
44) Dimethylphthalate-d6	13.77	166	206549	0.00	ng/ul	0.00
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.36	176	12928	0.00	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.11	188	42216	20.53	ng/ul	-0.10
79) Pyrene-d10	0.00	212	0	0.00	ng/ul	
90) Benzo(a)pyrene-d12	0.00	264	0	0.00	ng/ul	

Target Compounds

				Ovalue	
2) 1,4-Dioxane	3.26	88	4829	6.867	ng/uL 94
4) Benzaldehyde	6.88	77	10539	6.120	ng/ul 90
8) Bis(2-Chloroethyl)ether	7.13	93	98595	49.489	ng/ul# 49
14) Acetophenone	8.54	105	32118	9.260	ng/ul 99

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

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