

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061419\
 Data File : BM020967.D
 Acq On : 15 Jun 2019 06:14
 Operator : HP/JU
 Sample : K3332-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A41W5

Quant Time: Jun 15 06:51:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 14 15:40:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	315723	20.00	ng/ul	0.00
18) Naphthalene-d8	10.58	136	1257804	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	814935	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	1824142	20.00	ng/ul	0.00
77) Chrysene-d12	21.36	240	1316147	20.00	ng/ul	0.00
85) Perylene-d12	23.63	264	1312341	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	0d	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	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul	
46) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
78) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
89) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds

					Ovalue	
6) Phenol	6.99	94	27198	1.046	ng/ul	99
44) Dimethylphthalate	13.89	163	741925	11.068	ng/ul	99
76) Fluoranthene	19.23	202	130918	1.201	ng/ul	98
79) Pyrene	19.60	202	116734	1.255	ng/ul#	96
87) Benzo(b)fluoranthene	22.94	252	93701	1.127	ng/ul#	92

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

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