

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061119\
 Data File : BM020882.D
 Acq On : 11 Jun 2019 19:52
 Operator : HP/JU
 Sample : K3200-04
 Misc : GCMS CONFIRMATION
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jun 12 04:03:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM060619MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jun 09 01:02:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	347256	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	1393439	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	817810	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	1585103	20.00	ng/ul	0.00
77) Chrysene-d12	21.37	240	1250027	20.00	ng/ul	0.00
85) Perylene-d12	23.65	264	1515663	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	308	0.04	ng/uL	-0.01
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.15	67	222	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	10.60	131	2006	0.08	ng/ul	-0.14
43) Dimethylphthalate-d6	0.00	166	0	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	15.44	176	3577	0.07	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.19	188	1585103	22.89	ng/ul	-0.10
78) Pyrene-d10	19.59	212	1228	0.02	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.65	264	1500569	21.78	ng/ul	0.14

Target Compounds Ovalue

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

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