

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM093019\
 Data File : BM022848.D
 Acq On : 01 Oct 2019 06:14
 Operator : JU
 Sample : K5027-05
 Misc : GCMS CONFIRMATION
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 01 06:52:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 27 18:03:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	241860	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	996388	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.44	164	647568	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.18	188	1362815	20.00	ng/ul	0.00
78) Chrysene-d12	21.36	240	1131087	20.00	ng/ul	-0.01
86) Perylene-d12	23.62	264	1017966	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
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	7.27	67	225	0.02	ng/ul	0.12
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	1190	0.06	ng/ul	-0.14
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
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	15.43	176	1124	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.18	188	1362815	20.48	ng/ul	-0.11
79) Pyrene-d10	19.57	212	256	0.00	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.62	264	1008218	19.15	ng/ul	0.12

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

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