

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000738.D
 Acq On : 15 Oct 2019 03:20
 Operator : JU
 Sample : K5175-15
 Misc : GCMS CONFIRMATION
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPG2

Quant Time: Oct 15 05:30:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 03:51:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	236605	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	930352	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	523629	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	1003695	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	898980	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	999784	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	0.00	99	0d	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	0d	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0d	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0d	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul	
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
79) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
90) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds

					Ovalue	
70) Phenanthrene	11.32	178	80739	1.476	ng/ul	98
77) Fluoranthene	12.52	202	170121	3.035	ng/ul	96
80) Pyrene	12.75	202	84891	1.331	ng/ul	98
83) Benzo(a)anthracene	13.96	228	63807	1.058	ng/ul	98
85) Chrysene	13.99	228	103486	1.845	ng/ul	99
88) Benzo(b)fluoranthene	15.02	252	73560	1.233	ng/ul	99

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

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