

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101719\
 Data File : BP000817.D
 Acq On : 16 Oct 2019 16:57
 Operator : JU
 Sample : K5223-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPH5

Quant Time: Oct 17 09:19:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 07:55:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	173404	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	664397	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	370289	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	687525	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	618127	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	700667	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.35	96	17972	4.30	ng/uL	0.00
5) Phenol-d5	6.39	99	287214	21.47	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	151178	22.61	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	256952	22.80	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	217751	21.31	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	118576	23.48	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	126577	23.54	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	224930	21.85	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	290247	25.36	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	618029	23.84	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	745205	23.37	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	69132	17.32	ng/ul	0.00
58) Fluorene-d10	10.32	176	526278	23.73	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	40695	12.43	ng/ul	0.00
71) Anthracene-d10	11.35	188	753857	23.67	ng/ul	0.00
79) Pyrene-d10	12.73	212	790028	22.60	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	802543	23.20	ng/ul	0.00
Target Compounds						
6) Phenol	6.40	94	33642	2.495	ng/ul	97
45) Dimethylphthalate	9.51	163	203800	7.891	ng/ul	100

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

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