

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101619\
 Data File : BP000773.D
 Acq On : 15 Oct 2019 18:10
 Operator : JU
 Sample : K5175-22
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPG6

Quant Time: Oct 16 02:40:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 16 02:22:01 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	226000	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	826667	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	432001	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	761189	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	670272	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	756593	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.37	96	21497	3.95	ng/uL	-0.01
5) Phenol-d5	6.39	99	367884	21.10	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	203405	23.34	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	333884	22.73	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	226334	16.99	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	158210	25.18	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	170333	25.45	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	290207	22.66	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	330975	23.24	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	784228	25.93	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	944307	25.38	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	97359	20.90	ng/ul	0.00
58) Fluorene-d10	10.32	176	652953	25.24	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	65130	17.97	ng/ul	0.00
71) Anthracene-d10	11.35	188	930819	26.40	ng/ul	0.00
79) Pyrene-d10	12.73	212	1041506	27.48	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1092159	29.24	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	56613	3.221 ng/ul	97
45) Dimethylphthalate	9.51	163	295867	9.819 ng/ul	100

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

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