

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101619\
 Data File : BP000765.D
 Acq On : 15 Oct 2019 14:57
 Operator : JU
 Sample : K5175-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPG1

Quant Time: Oct 16 04:25:53 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	208103	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	764572	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	401469	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	717531	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	655024	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	738544	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.36	96	29720	5.92	ng/uL	-0.02
5) Phenol-d5	6.39	99	469882	29.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	254973	31.77	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	420187	31.07	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	312897	25.51	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	200715	34.54	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	212479	34.33	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	362518	30.61	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	476659	36.18	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	959814	34.15	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1162855	33.64	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	110241	25.47	ng/ul	0.00
58) Fluorene-d10	10.32	176	805492	33.50	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	75000	21.95	ng/ul	0.00
71) Anthracene-d10	11.35	188	1126014	33.88	ng/ul	0.00
79) Pyrene-d10	12.73	212	1227104	33.13	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1293568	35.48	ng/ul	0.00

Target Compounds				Ovalue		
6) Phenol	6.40	94	58343	3.605	ng/ul	98
45) Dimethylphthalate	9.51	163	305301	10.902	ng/ul	100

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

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