

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101219\
 Data File : BP000660.D
 Acq On : 12 Oct 2019 12:19
 Operator : JU
 Sample : K5199-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPJ8

Quant Time: Oct 12 21:46:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 09 17:20:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	200856	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	780754	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	448048	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	844650	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	727748	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	803098	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.36	96	25154	5.19	ng/uL	-0.01
5) Phenol-d5	6.38	99	412768	26.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	226370	29.22	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	359623	27.55	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	322398	27.24	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	172163	29.01	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	184030	29.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	321660	26.59	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	400834	29.80	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	930652	29.67	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1098470	28.47	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	100677	20.84	ng/ul	0.00
58) Fluorene-d10	10.32	176	775997	28.92	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	80938	20.12	ng/ul	0.00
71) Anthracene-d10	11.35	188	1110934	28.40	ng/ul	0.00
79) Pyrene-d10	12.73	212	1196193	29.07	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1156178	29.16	ng/ul	0.00
Target Compounds						
6) Phenol	6.40	94	29716	1.903	ng/ul	98
45) Dimethylphthalate	9.51	163	323266	10.344	ng/ul	100

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

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