

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101719\
 Data File : BP000840.D
 Acq On : 17 Oct 2019 14:47
 Operator : JU
 Sample : PB124015BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK15

Quant Time: Oct 17 15:41:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 09:27:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	167072	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	649615	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	366630	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	687367	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	610571	20.00	ng/ul	0.00
86) Perylene-d12	15.46	264	690239	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	5158	1.28	ng/uL	-0.04
5) Phenol-d5	6.39	99	92791	7.20	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	186087	28.88	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	269305	24.80	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	172490	17.52	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	151359	30.65	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	163766	31.14	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	278070	27.63	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	29768	2.66	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	809955	31.56	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	964133	30.54	ng/ul	0.00
52) 4-Nitrophenol-d4	9.93	143	21929	5.55	ng/ul	0.01
58) Fluorene-d10	10.32	176	680402	30.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	66233	20.23	ng/ul	0.00
71) Anthracene-d10	11.35	188	1019823	32.03	ng/ul	0.00
79) Pyrene-d10	12.74	212	1081484	31.32	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1093107	32.08	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	19424	1.495 ng/ul	95
45) Dimethylphthalate	9.51	163	27507	1.076 ng/ul	99
76) Di-n-butylphthalate	11.87	149	39566	1.016 ng/ul	98

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

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