

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000722.D
 Acq On : 14 Oct 2019 20:31
 Operator : JU
 Sample : K5097-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLME9

Quant Time: Oct 15 03:54:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 03:51:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	226897	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	842701	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	457450	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.30	188	837465	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	730786	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	800921	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	16851	3.08	ng/uL	0.00
5) Phenol-d5	6.39	99	297614	17.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	157039	17.95	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	259456	17.60	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	223140	16.69	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	116744	18.22	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	123064	18.04	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	218767	16.76	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	282991	19.49	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	602706	18.82	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	742326	18.84	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	66260	13.43	ng/ul	0.00
58) Fluorene-d10	10.32	176	514452	18.78	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	41810	10.48	ng/ul	0.00
71) Anthracene-d10	11.35	188	747571	19.27	ng/ul	0.00
79) Pyrene-d10	12.73	212	839047	20.30	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	818339	20.70	ng/ul	0.00

Target Compounds

				Ovalue
6) Phenol	6.40	94	53718	3.044 ng/ul 98
45) Dimethylphthalate	9.51	163	198602	6.224 ng/ul 99

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

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