

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000730.D
 Acq On : 14 Oct 2019 23:44
 Operator : JU
 Sample : K5097-20
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLM82

Quant Time: Oct 15 03:56:24 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	232457	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	868448	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	470914	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	859526	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	774230	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	857758	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	21845	3.90	ng/uL	0.00
5) Phenol-d5	6.39	99	370124	20.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	194331	21.68	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	321535	21.28	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	293247	21.41	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	148520	22.50	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	157946	22.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	285036	21.19	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	351971	23.52	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	746379	22.64	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	910871	22.46	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	85316	16.80	ng/ul	0.00
58) Fluorene-d10	10.32	176	638360	22.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	61384	15.00	ng/ul	0.00
71) Anthracene-d10	11.35	188	912290	22.91	ng/ul	0.00
79) Pyrene-d10	12.73	212	1040304	23.76	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1042786	24.63	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	42431	2.347 ng/ul	99
45) Dimethylphthalate	9.51	163	172995	5.267 ng/ul	99

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

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