

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000726.D
 Acq On : 14 Oct 2019 22:07
 Operator : JU
 Sample : K5097-16
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLMF4

Quant Time: Oct 15 04:23:59 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	236851	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	885793	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	478582	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	889234	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	799768	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	891904	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	37410	6.55	ng/uL	0.00
5) Phenol-d5	6.39	99	618761	33.87	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	310828	34.03	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	521747	33.90	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	490137	35.11	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	248940	36.97	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	270574	37.74	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	468518	34.14	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	652057	42.72	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1233830	36.83	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1478510	35.88	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	160213	31.05	ng/ul	0.00
58) Fluorene-d10	10.32	176	1026026	35.79	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	109906	25.95	ng/ul	0.00
71) Anthracene-d10	11.35	188	1485385	36.06	ng/ul	0.00
79) Pyrene-d10	12.74	212	1563352	34.57	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1602723	36.40	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	78621	4.269	ng/ul 99
45) Dimethylphthalate	9.51	163	347872	10.421	ng/ul 100

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

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