

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM103119\
 Data File : BM023531.D
 Acq On : 31 Oct 2019 22:04
 Operator : JU
 Sample : K5426-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNJ6

Quant Time: Nov 01 07:16:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 01 07:12:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	394325	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	1606997	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	967809	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.01	188	2050990	20.00	ng/ul	0.00
78) Chrysene-d12	21.20	240	2080346	20.00	ng/ul	0.00
86) Perylene-d12	23.39	264	2516243	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	47521	5.72	ng/uL	0.00
5) Phenol-d5	6.79	99	179924	5.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	504294	25.98	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	535745	20.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	305296	11.35	ng/ul	0.00
19) Nitrobenzene-d5	8.76	128	327511	27.07	ng/ul	0.00
22) 2-Nitrophenol-d4	9.47	143	341400	25.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	575336	21.81	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	31423	1.04	ng/ul	0.00
44) Dimethylphthalate-d6	13.68	166	2002830	27.25	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	2300564	27.32	ng/ul	0.00
52) 4-Nitrophenol-d4	14.48	143	42155	3.31	ng/ul	0.00
58) Fluorene-d10	15.26	176	1731798	27.82	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	221164	17.32	ng/ul	0.00
71) Anthracene-d10	17.10	188	2651296	27.35	ng/ul	0.00
79) Pyrene-d10	19.41	212	3282817	28.69	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.26	264	3898251	30.41	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.82	94	40982	1.202	ng/ul 97
45) Dimethylphthalate	13.72	163	157944	2.140	ng/ul# 97

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

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