

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM023002.D
 Acq On : 05 Oct 2019 21:28
 Operator : JU
 Sample : K5050-20
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLMEO

Quant Time: Oct 06 23:35:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 16:45:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	247982	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	1078310	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.42	164	630500	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.16	188	1086781	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	839158	20.00	ng/ul	-0.01
86) Perylene-d12	23.60	264	963540	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	25569	4.46	ng/uL	0.00
5) Phenol-d5	6.96	99	473439	21.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	276864	23.09	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.32	132	393009	22.27	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	388048	21.63	ng/ul	-0.01
19) Nitrobenzene-d5	8.94	128	191416	23.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	202412	22.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	391046	21.61	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.71	131	532378	24.22	ng/ul	0.00
44) Dimethylphthalate-d6	13.83	166	1111676	22.72	ng/ul	0.00
47) Acenaphthylene-d8	14.12	160	1326643	23.48	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	103515	10.71	ng/ul	-0.01
58) Fluorene-d10	15.42	176	916606	21.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	81688	11.55	ng/ul	-0.01
71) Anthracene-d10	17.26	188	1218059	22.95	ng/ul	-0.01
79) Pyrene-d10	19.55	212	1153572	25.54	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.46	264	1146872	23.01	ng/ul	-0.01

Target Compounds

					Ovalue
6) Phenol	6.98	94	41840	1.822	ng/ul 97
45) Dimethylphthalate	13.88	163	303068	6.029	ng/ul 100

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

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