

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082619\
 Data File : BM022326.D
 Acq On : 26 Aug 2019 12:15
 Operator : HP/JU
 Sample : K4369-05RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CD8RE

Quant Time: Aug 27 15:04:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM082219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 27 10:56:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	64255	20.00	ng/ul	0.05
18) Naphthalene-d8	10.34	136	141613	20.00	ng/ul	0.01
35) Acenaphthene-d10	14.20	164	92564	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	216882	20.00	ng/ul	0.00
77) Chrysene-d12	21.18	240	203771	20.00	ng/ul	0.00
85) Perylene-d12	23.37	264	193781	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	2831	1.97	ng/uL	0.00
5) Phenol-d5	6.76	99	17380	3.13	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.92	67	50540	15.67	ng/ul	0.01
9) 2-Chlorophenol-d4	7.09	132	58765	13.12	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	38229	7.95	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	39267	37.19	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	47327	38.51	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	80157	32.04	ng/ul	0.00
29) 4-Chloroaniline-d4	10.50	131	5000	1.64	ng/ul	0.01
43) Dimethylphthalate-d6	13.63	166	292767	35.91	ng/ul	0.00
46) Acenaphthylene-d8	13.90	160	319322	36.37	ng/ul	0.00
51) 4-Nitrophenol-d4	14.52	143	7670	6.35	ng/ul	0.05
57) Fluorene-d10	15.21	176	248586	36.77	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.38	200	47130	28.16	ng/ul	0.00
70) Anthracene-d10	17.07	188	427902	37.26	ng/ul	0.00
78) Pyrene-d10	19.37	212	529124	49.96	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.23	264	406302	38.17	ng/ul	0.00

Target Compounds

					Ovalue
20) Nitrobenzene	8.77	77	379721	122.119	ng/ul 98
36) 1,2,4,5-Tetrachlorobenzene	12.37	216	142744	37.288	ng/ul 98
44) Dimethylphthalate	13.68	163	47115	5.680	ng/ul 99
72) 1,2,3,4-Tetrachlorobenzene	12.99	216	587343	152.252	ng/uL 98
73) Pentachlorobenzene	14.52	250	13108	3.132	ng/uL 91

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

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