

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\
 Data File : BM024248.D
 Acq On : 24 Dec 2019 11:07
 Operator : JU
 Sample : K6236-20
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFEW2

Quant Time: Dec 24 12:28:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 05:54:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	253335	20.00	ng/ul	0.00
18) Naphthalene-d8	10.20	136	1095884	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.09	164	679939	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1391091	20.00	ng/ul	0.00
78) Chrysene-d12	21.06	240	1215369	20.00	ng/ul	0.00
86) Perylene-d12	23.19	264	1381771	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.00	96	14896	2.60	ng/uL	0.00
5) Phenol-d5	6.63	99	234345	9.77	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.78	67	180058	12.52	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	196348	11.48	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	120950	6.17	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	92828	11.79	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	99545	11.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	169588	10.08	ng/ul	0.01
29) 4-Chloroaniline-d4	10.36	131	201051	9.83	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	681466	13.55	ng/ul	0.00
47) Acenaphthylene-d8	13.78	160	757468	12.57	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	41808	4.63	ng/ul	0.03
58) Fluorene-d10	15.10	176	577869	13.11	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.25	200	50647	5.97	ng/ul	0.00
71) Anthracene-d10	16.95	188	833820	13.17	ng/ul	0.00
79) Pyrene-d10	19.26	212	948232	16.38	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.06	264	975597	14.21	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.66	94	36429	1.471	ng/ul 96
45) Dimethylphthalate	13.57	163	80662	1.544	ng/ul 99

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

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