

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011821\
 Data File : BM028441.D
 Acq On : 19 Jan 2021 01:24
 Operator : CG/JU
 Sample : M1099-12
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F1C08

Quant Time: Jan 19 02:17:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011621MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 18 10:13:28 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	36157	20.00	ng/ul	0.00
18) Naphthalene-d8	10.90	136	143798	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.72	164	79512	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.45	188	160717	20.00	ng/ul	0.00
78) Chrysene-d12	21.57	240	154921	20.00	ng/ul	0.00
86) Perylene-d12	23.89	264	152675	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	4543	5.29	ng/uL	0.00
5) Phenol-d5	7.23	99	86357	27.97	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	54427	28.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.61	132	74702	27.91	ng/ul	0.00
13) 4-Methylphenol-d8	8.79	113	70200	28.16	ng/ul	0.00
19) Nitrobenzene-d5	9.26	128	33934	27.58	ng/ul	0.00
22) 2-Nitrophenol-d4	9.98	143	37190	27.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.52	165	67612	27.61	ng/ul	0.00
29) 4-Chloroaniline-d4	11.04	131	92349	32.40	ng/ul	0.00
44) Dimethylphthalate-d6	14.13	166	180831	28.64	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	237810	29.49	ng/ul	0.00
52) 4-Nitrophenol-d4	14.90	143	28084	23.74	ng/ul	0.00
58) Fluorene-d10	15.71	176	154311	28.28	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	19427	17.56	ng/ul	0.00
71) Anthracene-d10	17.54	188	232770	28.13	ng/ul	0.00
79) Pyrene-d10	19.81	212	249269	27.55	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.74	264	258294	29.81	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	7.26	94	5666	1.845	ng/ul 97
31) Hexachlorobutadiene	11.23	225	2297	1.496	ng/ul 95
45) Dimethylphthalate	14.17	163	8228	1.286	ng/ul 99

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

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