

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011821\
 Data File : BM028434.D
 Acq On : 18 Jan 2021 20:33
 Operator : CG/JU
 Sample : M1099-15
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F1C11

Quant Time: Jan 19 00:23:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011621MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 18 23:34:50 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	36996	20.00	ng/ul	0.00
18) Naphthalene-d8	10.91	136	142548	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.73	164	75548	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.45	188	140993	20.00	ng/ul	0.00
78) Chrysene-d12	21.57	240	126722	20.00	ng/ul	0.00
86) Perylene-d12	23.89	264	128690	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	5562	6.33	ng/uL	0.00
5) Phenol-d5	7.24	99	103218	32.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	64862	32.81	ng/ul	0.00
9) 2-Chlorophenol-d4	7.61	132	89783	32.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.80	113	82747	32.44	ng/ul	0.00
19) Nitrobenzene-d5	9.26	128	40691	33.36	ng/ul	0.00
22) 2-Nitrophenol-d4	9.99	143	44850	33.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.52	165	77866	32.08	ng/ul	0.00
29) 4-Chloroaniline-d4	11.05	131	104029	36.82	ng/ul	0.00
44) Dimethylphthalate-d6	14.13	166	202675	33.78	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	263011	34.32	ng/ul	0.00
52) 4-Nitrophenol-d4	14.91	143	31793	28.29	ng/ul	0.00
58) Fluorene-d10	15.71	176	163721	31.57	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	23648	24.37	ng/ul	0.00
71) Anthracene-d10	17.55	188	233756	32.20	ng/ul	0.00
79) Pyrene-d10	19.82	212	238573	32.23	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.74	264	245254	33.58	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.26	94	5078	1.616	ng/ul 95
31) Hexachlorobutadiene	11.23	225	5780	3.798	ng/ul 97
45) Dimethylphthalate	14.17	163	12896	2.121	ng/ul 98

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

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