

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011821\
 Data File : BM028427.D
 Acq On : 18 Jan 2021 16:18
 Operator : CG/JU
 Sample : M1099-18
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F1C14

Quant Time: Jan 18 16:54:16 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	34345	20.00	ng/ul	0.00
18) Naphthalene-d8	10.91	136	134256	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.73	164	72780	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.46	188	139740	20.00	ng/ul	0.00
78) Chrysene-d12	21.57	240	125783	20.00	ng/ul	0.00
86) Perylene-d12	23.90	264	127734	20.00	ng/ul	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.34	96	4413	5.41	ng/uL	0.00
5) Phenol-d5	7.24	99	90812	30.97	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.41	67	56168	30.61	ng/ul	0.00
9) 2-Chlorophenol-d4	7.62	132	77558	30.51	ng/ul	0.00
13) 4-Methylphenol-d8	8.80	113	73451	31.02	ng/ul	0.00
19) Nitrobenzene-d5	9.26	128	35616	31.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.99	143	39933	31.83	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.53	165	70143	30.68	ng/ul	0.00
29) 4-Chloroaniline-d4	11.05	131	95557	35.91	ng/ul	0.00
44) Dimethylphthalate-d6	14.13	166	190774	33.01	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	248400	33.65	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	31451	29.05	ng/ul	0.01
58) Fluorene-d10	15.72	176	156408	31.31	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	21989	22.87	ng/ul	0.00
71) Anthracene-d10	17.55	188	228778	31.80	ng/ul	0.00
79) Pyrene-d10	19.82	212	231593	31.52	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.75	264	236696	32.65	ng/ul	0.01

Target Compounds					Ovalue	
6) Phenol	7.27	94	16623	5.697	ng/ul	97
31) Hexachlorobutadiene	11.24	225	6055	4.225	ng/ul	98
45) Dimethylphthalate	14.17	163	16880	2.881	ng/ul	99

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

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