

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021121\
 Data File : BM028719.D
 Acq On : 11 Feb 2021 19:49
 Operator : CG/JU
 Sample : PB134593TB
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB36

Quant Time: Feb 11 23:47:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011621MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 11 16:46:31 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	26110	20.00	ng/ul	0.00
18) Naphthalene-d8	10.70	136	102890	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.54	164	55519	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.27	188	105248	20.00	ng/ul	0.00
78) Chrysene-d12	21.42	240	101657	20.00	ng/ul	0.00
86) Perylene-d12	23.66	264	108341	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.20	96	3429	5.53	ng/uL	0.00
5) Phenol-d5	7.06	99	60550	27.16	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.22	67	35842	25.69	ng/ul	0.00
9) 2-Chlorophenol-d4	7.42	132	53016	27.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.61	113	47586	26.43	ng/ul	0.00
19) Nitrobenzene-d5	9.06	128	23108	26.25	ng/ul	0.00
22) 2-Nitrophenol-d4	9.79	143	26059	27.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.32	165	45224	25.81	ng/ul	0.00
29) 4-Chloroaniline-d4	10.84	131	64429	31.59	ng/ul	0.00
44) Dimethylphthalate-d6	13.95	166	120104	27.24	ng/ul	0.00
47) Acenaphthylene-d8	14.23	160	159904	28.39	ng/ul	0.00
52) 4-Nitrophenol-d4	14.76	143	19359	23.44	ng/ul	0.00
58) Fluorene-d10	15.53	176	102471	26.89	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.66	200	15265	21.08	ng/ul	0.00
71) Anthracene-d10	17.37	188	156166	28.82	ng/ul	0.00
79) Pyrene-d10	19.65	212	160013	26.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.51	264	179101	29.13	ng/ul	0.00

Target Compounds Ovalue

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

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