

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
 Data File : BM031968.D
 Acq On : 31 Aug 2021 02:33
 Operator : CG/JU
 Sample : M3422-23
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBKE2

Quant Time: Aug 31 04:22:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM082821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 09:21:30 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	89703	20.000	ng/ul	0.00
20) Naphthalene-d8	10.263	136	383263	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.145	164	285580	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.904	188	637025	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.109	240	685076	20.000	ng/ul	# 0.00
88) Perylene-d12	23.250	264	664659	20.000	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	8617	4.432	ng/uL	0.00
4) Pyridine-d5	3.540	84	20004	3.623	ng/ul	0.02
7) Phenol-d5	6.704	99	42790	5.752	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.863	67	122507	28.465	ng/ul	0.00
11) 2-Chlorophenol-d4	7.045	132	135134	23.608	ng/ul	0.00
15) 4-Methylphenol-d8	8.222	113	87518	13.713	ng/ul	0.00
21) Nitrobenzene-d5	8.657	128	89220	32.050	ng/ul	0.00
24) 2-Nitrophenol-d4	9.369	143	101406	30.781	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.898	165	189560	28.678	ng/ul	0.00
31) 4-Chloroaniline-d4	10.422	131	104152	11.171	ng/ul	0.00
46) Dimethylphthalate-d6	13.574	166	743016	34.384	ng/ul	0.00
49) Acenaphthylene-d8	13.839	160	845643	33.044	ng/ul	0.00
54) 4-Nitrophenol-d4	14.404	143	18988	4.855	ng/ul	0.02
60) Fluorene-d10	15.151	176	650079	34.223	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.298	200	128899	29.448	ng/ul	0.00
73) Anthracene-d10	17.004	188	1126926	37.271	ng/ul	0.00
81) Pyrene-d10	19.309	212	1361051	42.416	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.115	264	1324431	37.915	ng/ul	0.00

Target Compounds	Qvalue

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

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