

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042878.D
 Acq On : 18 Nov 2023 12:27
 Operator : MA/JU
 Sample : 05226-21
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 18 21:10:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 22:28:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	25989	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.628	136	100946	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.469	164	63496	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.227	188	139555	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.409	240	145162	20.000	ng/u1	-0.01
88) Perylene-d12	23.768	264	185485	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	2940	3.550	ng/uL	-0.01
4) Pyridine-d5	3.622	84	14602	7.088	ng/u1	0.00
7) Phenol-d5	6.993	99	10642	4.207	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67	49264	27.718	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.340	132	35132	18.001	ng/u1	-0.01
15) 4-Methylphenol-d8	8.540	113	19111	9.596	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	332	0.375	ng/u1	0.14
24) 2-Nitrophenol-d4	9.728	143	20699	19.992	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.257	165	32015	16.630	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	39967	16.351	ng/u1	0.00
46) Dimethylphthalate-d6	13.892	166	153973	25.622	ng/u1	-0.01
49) Acenaphthylene-d8	14.169	160	159043	24.735	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.869	143	61	0.090	ng/u1	0.13
60) Fluorene-d10	15.469	176	132672	26.661	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.621	200	15416	14.583	ng/u1	0.00
73) Anthracene-d10	17.327	188	232836	29.770	ng/u1	-0.01
81) Pyrene-d10	19.621	212	318763	33.534	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.615	264	341220	30.755	ng/u1	-0.01

Target Compounds Qvalue

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

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