

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042879.D
 Acq On : 18 Nov 2023 13:03
 Operator : MA/JU
 Sample : 05226-22
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 18 21:10:19 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	20157	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.628	136	78987	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.468	164	49238	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.227	188	108879	20.000	ng/u1	-0.01
79) Chrysene-d12	21.409	240	112078	20.000	ng/u1	#-0.01
88) Perylene-d12	23.768	264	141848	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	3058	4.760	ng/uL	-0.01
4) Pyridine-d5	3.622	84	163	0.102	ng/u1	0.00
7) Phenol-d5	6.998	99	6589	3.359	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	36684	26.611	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	27775	18.349	ng/u1	-0.01
15) 4-Methylphenol-d8	8.545	113	15273	9.888	ng/u1	0.00
21) Nitrobenzene-d5	9.116	128	356	0.514	ng/u1	0.09
24) 2-Nitrophenol-d4	9.733	143	15434	19.051	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	23778	15.785	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	9970	5.213	ng/u1	0.01
46) Dimethylphthalate-d6	13.898	166	128499	27.575	ng/u1	0.00
49) Acenaphthylene-d8	14.174	160	129216	25.915	ng/u1	-0.01
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/u1	
60) Fluorene-d10	15.468	176	109149	28.285	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.621	200	11529	13.979	ng/u1	0.00
73) Anthracene-d10	17.327	188	178241	29.211	ng/u1	-0.01
81) Pyrene-d10	19.621	212	246904	33.641	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.615	264	262456	30.933	ng/u1	-0.01

Target Compounds	Qvalue

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

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