

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121823\
 Data File : BM043405.D
 Acq On : 18 Dec 2023 22:22
 Operator : MA/JU
 Sample : 05807-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AX7

Quant Time: Dec 19 01:58:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 15 22:09:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.992	152	522381	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	2267205	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.621	164	1179855	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.368	188	1982972	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	1386640	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	1566317	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	75991	4.728	ng/uL	0.00
4) Pyridine-d5	3.810	84	1153225	24.929	ng/u1	0.00
7) Phenol-d5	7.157	99	1576768	26.534	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	1007200	26.762	ng/u1	0.00
11) 2-Chlorophenol-d4	7.522	132	1189370	26.213	ng/u1	0.00
15) 4-Methylphenol-d8	8.704	113	1250252	27.113	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	613787	27.981	ng/u1	0.00
24) 2-Nitrophenol-d4	9.886	143	655913	28.940	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	1094199	27.348	ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	1657390	28.157	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2989124	28.143	ng/u1	0.00
49) Acenaphthylene-d8	14.315	160	3689713	29.285	ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	512818	25.510	ng/u1	0.00
60) Fluorene-d10	15.615	176	2480899	28.247	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	294640	22.466	ng/u1	0.00
73) Anthracene-d10	17.468	188	3310358	29.767	ng/u1	0.00
81) Pyrene-d10	19.750	212	3104980	28.363	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.809	264	2911315	30.120	ng/u1	0.00

Target Compounds Qvalue

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

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