

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100924\
 Data File : BM047944.D
 Acq On : 09 Oct 2024 14:06
 Operator : RC/JU
 Sample : P4187-13
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EY072

Quant Time: Oct 09 14:51:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 08:32:31 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	184729	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	720448	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.428	164	427561	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.163	188	875814	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.386	240	944867	20.000	ng/u1	-0.01
88) Perylene-d12	24.380	264	1165467	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	24078	4.212	ng/uL	-0.01
4) Pyridine-d5	3.664	84	104689	6.255	ng/u1	0.00
7) Phenol-d5	6.958	99	131242	7.513	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	295330	23.826	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.322	132	373871	29.418	ng/u1	-0.01
15) 4-Methylphenol-d8	8.499	113	239182	18.417	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	192814	33.613	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	212135	38.591	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.204	165	360517	32.831	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.716	131	453861	26.734	ng/u1	-0.01
46) Dimethylphthalate-d6	13.834	166	1005081	32.424	ng/u1	-0.01
49) Acenaphthylene-d8	14.116	160	1192940	32.458	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.628	143	36455	5.907	ng/u1	0.00
60) Fluorene-d10	15.416	176	899986	33.519	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.533	200	143973	28.091	ng/u1	0.00
73) Anthracene-d10	17.263	188	1495808	34.734	ng/u1	-0.01
81) Pyrene-d10	19.551	212	1834579	34.211	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.174	264	2207129	36.500	ng/u1	-0.02

Target Compounds Qvalue

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

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