

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100924\
 Data File : BM047943.D
 Acq On : 09 Oct 2024 12:48
 Operator : RC/JU
 Sample : PB163684BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK684

Quant Time: Oct 10 00:09:19 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	7.792	152	268595	20.000	ng/u1	0.00
20) Naphthalene-d8	10.580	136	1042142	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.427	164	633660	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.162	188	1299366	20.000	ng/u1	-0.01
79) Chrysene-d12	21.391	240	1393735	20.000	ng/u1	0.00
88) Perylene-d12	24.380	264	1669184	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	35906	4.320	ng/uL	-0.01
4) Pyridine-d5	3.657	84	507586	20.857	ng/u1	0.00
7) Phenol-d5	6.963	99	573849	22.594	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	341834	18.967	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	486290	26.316	ng/u1	0.00
15) 4-Methylphenol-d8	8.498	113	443269	23.474	ng/u1	0.00
21) Nitrobenzene-d5	8.945	128	223504	26.936	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	239743	30.150	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.204	165	417039	26.255	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.716	131	606101	24.681	ng/u1	-0.01
46) Dimethylphthalate-d6	13.833	166	1121326	24.408	ng/u1	-0.01
49) Acenaphthylene-d8	14.115	160	1365861	25.075	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.627	143	211036	23.072	ng/u1	0.00
60) Fluorene-d10	15.415	176	1007249	25.312	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.533	200	161366	21.222	ng/u1	0.00
73) Anthracene-d10	17.262	188	1613184	25.249	ng/u1	-0.01
81) Pyrene-d10	19.550	212	1972875	24.941	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.174	264	2307699	26.646	ng/u1	-0.02

Target Compounds Qvalue

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

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