

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
 Data File : BM047939.D
 Acq On : 09 Oct 2024 08:53
 Operator : RC/JU
 Sample : SP6637
 Misc : SFAM SURROGATES
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP6637

Quant Time: Oct 09 09:27:50 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	304178	20.000	ng/u1	0.00
20) Naphthalene-d8	10.586	136	1241370	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.427	164	774436	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.168	188	1553638	20.000	ng/u1	0.00
79) Chrysene-d12	21.397	240	1416515	20.000	ng/u1	0.00
88) Perylene-d12	24.397	264	1579854	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	108172	11.492	ng/uL	0.00
4) Pyridine-d5	3.663	84	1608816	58.375	ng/u1	0.00
7) Phenol-d5	6.969	99	1841997	64.041	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	1108883	54.329	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	1556485	74.377	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	1481329	69.269	ng/u1	0.00
21) Nitrobenzene-d5	8.951	128	772899	78.198	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	856147	90.390	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	1429626	75.557	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	2110494	72.149	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	3857007	68.695	ng/u1	0.00
49) Acenaphthylene-d8	14.121	160	4728757	71.033	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	779005	69.686	ng/u1	0.00
60) Fluorene-d10	15.421	176	3397459	69.859	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	554626	61.004	ng/u1	0.00
73) Anthracene-d10	17.268	188	5229032	68.448	ng/u1	0.00
81) Pyrene-d10	19.556	212	5838224	72.620	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.203	264	5856665	71.449	ng/u1	0.01

Target Compounds Qvalue

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

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