

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
 Data File : BM047892.D
 Acq On : 07 Oct 2024 20:24
 Operator : RC/JU
 Sample : P4109-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 07 23:23:02 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	7.798	152	295216	20.000	ng/ul	0.00
20) Naphthalene-d8	10.586	136	1183381	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.433	164	741314	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.168	188	1521260	20.000	ng/ul	0.00
79) Chrysene-d12	21.397	240	1416076	20.000	ng/ul	0.00
88) Perylene-d12	24.391	264	1612203	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.251	96	22952	2.512	ng/uL	0.00
4) Pyridine-d5	3.663	84	332670	12.437	ng/ul	0.00
7) Phenol-d5	6.963	99	427252	15.305	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	253736	12.809	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	359457	17.698	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	348335	16.783	ng/ul	0.00
21) Nitrobenzene-d5	8.951	128	167836	17.813	ng/ul	0.00
24) 2-Nitrophenol-d4	9.675	143	180222	19.960	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	344617	19.106	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	472438	16.942	ng/ul	0.00
46) Dimethylphthalate-d6	13.839	166	1065254	19.820	ng/ul	0.00
49) Acenaphthylene-d8	14.121	160	1240145	19.461	ng/ul	0.00
54) 4-Nitrophenol-d4	14.627	143	160869	15.033	ng/ul	0.00
60) Fluorene-d10	15.421	176	965461	20.739	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	107357	12.060	ng/ul	0.00
73) Anthracene-d10	17.268	188	1511388	20.205	ng/ul	0.00
81) Pyrene-d10	19.556	212	1820225	22.648	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.191	264	1811426	21.655	ng/ul	0.00

Target Compounds	Qvalue

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

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