

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100324\
 Data File : BM047864.D
 Acq On : 05 Oct 2024 06:12
 Operator : RC/JU
 Sample : PB163918BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK918

Quant Time: Oct 05 06:54:48 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	179349	20.000	ng/u1	0.00
20) Naphthalene-d8	10.592	136	708287	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.433	164	416494	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.168	188	830593	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	736955	20.000	ng/u1	0.00
88) Perylene-d12	24.385	264	849962	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	32366	5.832	ng/uL	0.00
4) Pyridine-d5	3.663	84	455553	28.034	ng/u1	0.00
7) Phenol-d5	6.963	99	473992	27.949	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	309053	25.681	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	395825	32.080	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	369082	29.271	ng/u1	0.00
21) Nitrobenzene-d5	8.951	128	179806	31.884	ng/u1	0.00
24) 2-Nitrophenol-d4	9.675	143	181705	33.623	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	346717	32.116	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	517058	30.980	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	1032932	34.208	ng/u1	0.00
49) Acenaphthylene-d8	14.127	160	1198811	33.484	ng/u1	0.00
54) 4-Nitrophenol-d4	14.621	143	143061	23.796	ng/u1	0.00
60) Fluorene-d10	15.421	176	892312	34.116	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	103525	21.299	ng/u1	0.00
73) Anthracene-d10	17.268	188	1349063	33.032	ng/u1	0.00
81) Pyrene-d10	19.556	212	1518813	36.313	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.185	264	1477633	33.507	ng/u1	0.00

Target Compounds	Qvalue

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

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