

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100924\
 Data File : BM047964.D
 Acq On : 10 Oct 2024 04:35
 Operator : RC/JU
 Sample : PB163768BL
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK768

Quant Time: Oct 10 05:16:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 10 05:10:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	250385	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	970002	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	579966	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.163	188	1190784	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	1239599	20.000	ng/u1	0.00
88) Perylene-d12	24.374	264	1493206	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	34409	4.441	ng/uL	0.00
4) Pyridine-d5	3.658	84	484518	21.358	ng/u1	0.00
7) Phenol-d5	6.957	99	543393	22.951	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	333529	19.852	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	460191	26.715	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	415816	23.622	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	208418	26.986	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	219934	29.716	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	386414	26.136	ng/u1	0.00
31) 4-Chloroaniline-d4	10.716	131	566026	24.763	ng/u1	0.00
46) Dimethylphthalate-d6	13.833	166	1119712	26.629	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	1303007	26.136	ng/u1	0.00
54) 4-Nitrophenol-d4	14.622	143	193823	23.152	ng/u1	0.00
60) Fluorene-d10	15.410	176	977149	26.829	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.533	200	151601	21.756	ng/u1	0.00
73) Anthracene-d10	17.257	188	1626615	27.781	ng/u1	0.00
81) Pyrene-d10	19.545	212	1990166	28.288	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.168	264	2310678	29.825	ng/u1	0.00

Target Compounds Qvalue

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

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