

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110824\
 Data File : BM048571.D
 Acq On : 09 Nov 2024 18:31
 Operator : RC/JU
 Sample : PB164674BL
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK674

Quant Time: Nov 09 22:44:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 07 22:50:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	230778	20.000	ng/u1	0.00
20) Naphthalene-d8	10.445	136	949258	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	581156	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.057	188	1163425	20.000	ng/u1	0.00
79) Chrysene-d12	21.280	240	1042778	20.000	ng/u1	0.00
88) Perylene-d12	24.191	264	1225199	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	31349	5.415	ng/uL	0.00
4) Pyridine-d5	3.546	84	427379	26.789	ng/u1	0.00
7) Phenol-d5	6.846	99	511506	29.116	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	309925	29.371	ng/u1	0.00
11) 2-Chlorophenol-d4	7.198	132	441312	30.326	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	407531	29.502	ng/u1	0.00
21) Nitrobenzene-d5	8.816	128	209530	30.454	ng/u1	0.00
24) 2-Nitrophenol-d4	9.539	143	231847	31.681	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	382839	28.161	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	527453	27.315	ng/u1	0.00
46) Dimethylphthalate-d6	13.727	166	1122186	30.249	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	1327446	30.349	ng/u1	0.00
54) 4-Nitrophenol-d4	14.533	143	167937	24.373	ng/u1	0.00
60) Fluorene-d10	15.304	176	979525	30.248	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	138246	23.621	ng/u1	0.00
73) Anthracene-d10	17.157	188	1519646	30.866	ng/u1	0.00
81) Pyrene-d10	19.451	212	1727458	33.009	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.991	264	1825599	31.891	ng/u1	0.00

Target Compounds Qvalue

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

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