

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111124\
 Data File : BM048604.D
 Acq On : 12 Nov 2024 04:14
 Operator : RC/JU
 Sample : PB164730BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK730

Quant Time: Nov 12 04:49:25 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	7.657	152	184416	20.000	ng/u1	0.00
20) Naphthalene-d8	10.445	136	759366	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	465972	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.051	188	954903	20.000	ng/u1	0.00
79) Chrysene-d12	21.280	240	926380	20.000	ng/u1	0.00
88) Perylene-d12	24.185	264	1089700	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	25519	5.516	ng/uL	0.00
4) Pyridine-d5	3.540	84	356187	27.939	ng/u1	0.00
7) Phenol-d5	6.840	99	409677	29.182	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.992	67	244804	29.032	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.192	132	352545	30.316	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	320767	29.059	ng/u1	0.00
21) Nitrobenzene-d5	8.816	128	166357	30.226	ng/u1	0.00
24) 2-Nitrophenol-d4	9.533	143	177518	30.324	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	298060	27.407	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	450708	29.177	ng/u1	0.00
46) Dimethylphthalate-d6	13.721	166	910098	30.596	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	1074943	30.652	ng/u1	0.00
54) 4-Nitrophenol-d4	14.539	143	124483	22.532	ng/u1	0.00
60) Fluorene-d10	15.304	176	795720	30.646	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	104652	21.786	ng/u1	0.00
73) Anthracene-d10	17.151	188	1261813	31.226	ng/u1	0.00
81) Pyrene-d10	19.450	212	1485669	31.956	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.985	264	1588769	31.205	ng/u1	0.00

Target Compounds Qvalue

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

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