

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091124\
 Data File : BM047474.D
 Acq On : 11 Sep 2024 09:26
 Operator : RC/JU
 Sample : PB163193BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK193

Quant Time: Sep 12 03:52:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 11 00:35:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	388384	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	1703748	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	1087213	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2205896	20.000	ng/u1	0.00
79) Chrysene-d12	21.462	240	1907633	20.000	ng/u1	0.00
88) Perylene-d12	24.509	264	1956177	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.304	96	58078	6.031	ng/uL	0.00
4) Pyridine-d5	3.710	84	807741	26.254	ng/u1	0.00
7) Phenol-d5	7.022	99	923436	25.365	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.198	67	622860	27.486	ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	748886	27.009	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	740404	25.874	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	389652	29.615	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	364797	29.348	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	669752	26.499	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	1042250	25.197	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2390294	29.912	ng/u1	0.00
49) Acenaphthylene-d8	14.192	160	2720953	28.217	ng/u1	0.00
54) 4-Nitrophenol-d4	14.668	143	413922	26.143	ng/u1	0.00
60) Fluorene-d10	15.486	176	2032314	29.431	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	274164	23.510	ng/u1	0.00
73) Anthracene-d10	17.333	188	3196232	30.090	ng/u1	0.00
81) Pyrene-d10	19.615	212	3566473	32.079	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.303	264	3136555	30.173	ng/u1	0.00

Target Compounds	Qvalue

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

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