

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048265.D
 Acq On : 28 Oct 2024 21:52
 Operator : RC/JU
 Sample : PB164389BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK389

Quant Time: Oct 28 23:21:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Oct 27 23:39:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	295326	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1246607	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	797276	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1553380	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1224161	20.000	ng/u1	0.00
88) Perylene-d12	24.274	264	1318155	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	33006	4.400	ng/uL	0.00
4) Pyridine-d5	3.587	84	467122	21.484	ng/u1	0.00
7) Phenol-d5	6.898	99	533316	21.020	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	353257	22.477	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	454332	22.020	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	431326	21.175	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	224700	21.985	ng/u1	0.00
24) 2-Nitrophenol-d4	9.598	143	245708	22.192	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.133	165	403839	19.988	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	615566	21.941	ng/u1	0.00
46) Dimethylphthalate-d6	13.774	166	1334663	22.401	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	1499384	22.319	ng/u1	0.00
54) 4-Nitrophenol-d4	14.580	143	174651	16.029	ng/u1	0.00
60) Fluorene-d10	15.357	176	1137614	22.596	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	182489	17.777	ng/u1	0.00
73) Anthracene-d10	17.204	188	1719312	23.407	ng/u1	0.00
81) Pyrene-d10	19.497	212	1814435	24.425	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.068	264	1600587	23.022	ng/u1	-0.01

Target Compounds	Qvalue

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

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