

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011524\
 Data File : BM043835.D
 Acq On : 15 Jan 2024 10:50
 Operator : MA/JU
 Sample : PB158368BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK368

Quant Time: Jan 16 01:15:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 10 01:07:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	286610	20.000	ng/u1	0.00
20) Naphthalene-d8	10.751	136	1290433	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.592	164	828887	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.339	188	1753771	20.000	ng/u1	0.00
79) Chrysene-d12	21.527	240	1417374	20.000	ng/u1	0.00
88) Perylene-d12	23.950	264	1506436	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	46926	6.322	ng/uL	0.00
4) Pyridine-d5	3.775	84	696076	30.441	ng/u1	0.00
7) Phenol-d5	7.116	99	837276	28.867	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	580961	32.142	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	664112	31.475	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	674491	29.193	ng/u1	-0.01
21) Nitrobenzene-d5	9.122	128	342620	31.681	ng/u1	0.00
24) 2-Nitrophenol-d4	9.845	143	380288	31.227	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	645263	29.588	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	1011327	29.280	ng/u1	0.00
46) Dimethylphthalate-d6	14.010	166	2307430	33.234	ng/u1	0.00
49) Acenaphthylene-d8	14.286	160	2604976	32.940	ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	360812	26.570	ng/u1	0.00
60) Fluorene-d10	15.586	176	1985869	33.988	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	311544	25.160	ng/u1	0.00
73) Anthracene-d10	17.445	188	3089360	34.615	ng/u1	0.00
81) Pyrene-d10	19.733	212	3378742	38.595	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.791	264	2927928	35.210	ng/u1	-0.01

Target Compounds Qvalue

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

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