

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011724\
 Data File : BM043862.D
 Acq On : 17 Jan 2024 12:10
 Operator : MA/JU
 Sample : PB158463BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK463

Quant Time: Jan 17 12:45:58 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	7.939	152	244786	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.751	136	1110920	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.586	164	726942	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.339	188	1566877	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1388096	20.000	ng/u1	0.00
88) Perylene-d12	23.956	264	1597205	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	36749	5.796	ng/uL	0.00
4) Pyridine-d5	3.775	84	536029	27.447	ng/u1	0.00
7) Phenol-d5	7.110	99	647033	26.119	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	463398	30.018	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.469	132	506686	28.117	ng/u1	-0.01
15) 4-Methylphenol-d8	8.657	113	522455	26.476	ng/u1	-0.02
21) Nitrobenzene-d5	9.122	128	264824	28.444	ng/u1	0.00
24) 2-Nitrophenol-d4	9.839	143	288300	27.499	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	488815	26.036	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.904	131	764596	25.714	ng/u1	0.00
46) Dimethylphthalate-d6	14.010	166	1793430	29.453	ng/u1	0.00
49) Acenaphthylene-d8	14.280	160	2036167	29.358	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.815	143	262037	22.002	ng/u1	-0.01
60) Fluorene-d10	15.580	176	1561396	30.471	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.721	200	233059	21.067	ng/u1	0.00
73) Anthracene-d10	17.439	188	2403918	30.148	ng/u1	0.00
81) Pyrene-d10	19.733	212	2739322	31.951	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.797	264	2699162	30.614	ng/u1	0.00

Target Compounds Qvalue

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

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