

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
 Data File : BM045817.D
 Acq On : 04 Jun 2024 15:08
 Operator : MA/JU
 Sample : PB160913BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK913

Quant Time: Jun 05 00:10:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 31 13:01:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.463	152	258336	20.000	ng/u1	0.00
20) Naphthalene-d8	10.234	136	1086630	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.110	164	695861	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.857	188	1426568	20.000	ng/u1	0.00
79) Chrysene-d12	21.068	240	1194689	20.000	ng/u1	0.00
88) Perylene-d12	23.786	264	1274159	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.058	96	35907	5.389	ng/uL	0.00
4) Pyridine-d5	3.463	84	438210	25.542	ng/u1	0.00
7) Phenol-d5	6.722	99	659211	30.037	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.834	67	502731	31.901	ng/u1	0.00
11) 2-Chlorophenol-d4	7.028	132	514154	31.364	ng/u1	0.00
15) 4-Methylphenol-d8	8.234	113	509876	30.752	ng/u1	0.00
21) Nitrobenzene-d5	8.634	128	263252	30.985	ng/u1	0.00
24) 2-Nitrophenol-d4	9.345	143	260689	30.318	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.892	165	466435	28.927	ng/u1	0.00
31) 4-Chloroaniline-d4	10.416	131	711868	31.899	ng/u1	0.00
46) Dimethylphthalate-d6	13.563	166	1608476	33.389	ng/u1	0.00
49) Acenaphthylene-d8	13.798	160	1732321	30.967	ng/u1	0.00
54) 4-Nitrophenol-d4	14.404	143	233047	22.415	ng/u1	0.00
60) Fluorene-d10	15.110	176	1321455	32.154	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.245	200	171746	23.576	ng/u1	0.00
73) Anthracene-d10	16.951	188	1952595	31.485	ng/u1	0.00
81) Pyrene-d10	19.251	212	2321599	31.248	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	1873819	31.466	ng/u1	0.00

Target Compounds	Qvalue

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

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