

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030824\
 Data File : BM044563.D
 Acq On : 08 Mar 2024 13:12
 Operator : MA/JU
 Sample : PB159474BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK474

Quant Time: Mar 08 21:55:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 05 23:05:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	224955	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.498	136	1001216	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.363	164	646597	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.115	188	1319210	20.000	ng/u1	0.00
79) Chrysene-d12	21.309	240	1108834	20.000	ng/u1	0.00
88) Perylene-d12	23.562	264	1120922	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	36414	5.490	ng/uL	-0.01
4) Pyridine-d5	3.669	84	578466	29.632	ng/u1	-0.01
7) Phenol-d5	6.887	99	674986	29.185	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.063	67	451745	31.622	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.251	132	521221	30.667	ng/u1	-0.02
15) 4-Methylphenol-d8	8.416	113	532846	29.336	ng/u1	-0.02
21) Nitrobenzene-d5	8.875	128	276445	32.790	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.592	143	299511	32.954	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	491611	31.782	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.645	131	816201	31.771	ng/u1	-0.02
46) Dimethylphthalate-d6	13.780	166	1620140	31.942	ng/u1	-0.01
49) Acenaphthylene-d8	14.051	160	1836588	31.661	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.569	143	288420	27.420	ng/u1	-0.01
60) Fluorene-d10	15.363	176	1398558	32.812	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	228015	26.627	ng/u1	0.00
73) Anthracene-d10	17.215	188	2130759	33.791	ng/u1	0.00
81) Pyrene-d10	19.515	212	2416028	35.525	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.421	264	2008356	34.088	ng/u1	0.00

Target Compounds Qvalue

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

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