

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
 Data File : BM044554.D
 Acq On : 07 Mar 2024 17:06
 Operator : MA/JU
 Sample : PB159446BL
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK446

Quant Time: Mar 07 23:58:07 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	7.722	152	238172	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.498	136	1024837	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.363	164	665863	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.116	188	1365665	20.000	ng/u1	0.00
79) Chrysene-d12	21.309	240	1185483	20.000	ng/u1	0.00
88) Perylene-d12	23.568	264	1196196	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	40983	5.836	ng/uL	-0.01
4) Pyridine-d5	3.669	84	624731	30.226	ng/u1	-0.01
7) Phenol-d5	6.893	99	722542	29.507	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.069	67	485203	32.079	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.252	132	561600	31.209	ng/u1	-0.02
15) 4-Methylphenol-d8	8.422	113	569758	29.628	ng/u1	-0.02
21) Nitrobenzene-d5	8.881	128	297744	34.502	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.593	143	321539	34.562	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	524197	33.108	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.645	131	858310	32.640	ng/u1	-0.02
46) Dimethylphthalate-d6	13.786	166	1726776	33.059	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	1940840	32.490	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	325945	30.091	ng/u1	-0.01
60) Fluorene-d10	15.363	176	1477993	33.672	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	252457	28.478	ng/u1	0.00
73) Anthracene-d10	17.216	188	2344034	35.909	ng/u1	0.00
81) Pyrene-d10	19.515	212	2660199	36.587	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.427	264	2273293	36.157	ng/u1	0.00

Target Compounds Qvalue

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

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