

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030824\
 Data File : BM044562.D
 Acq On : 08 Mar 2024 12:36
 Operator : MA/JU
 Sample : PB159472BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK472

Quant Time: Mar 08 21:54:04 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	230615	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.504	136	1020703	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.363	164	655545	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.115	188	1354090	20.000	ng/u1	0.00
79) Chrysene-d12	21.309	240	1217363	20.000	ng/u1	0.00
88) Perylene-d12	23.568	264	1279710	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	38340	5.639	ng/uL	-0.01
4) Pyridine-d5	3.669	84	574751	28.719	ng/u1	-0.01
7) Phenol-d5	6.887	99	670640	28.285	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.069	67	450677	30.773	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.251	132	510550	29.302	ng/u1	-0.02
15) 4-Methylphenol-d8	8.422	113	529903	28.458	ng/u1	-0.02
21) Nitrobenzene-d5	8.881	128	274124	31.894	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	295539	31.896	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	484306	30.712	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.651	131	805466	30.754	ng/u1	-0.01
46) Dimethylphthalate-d6	13.786	166	1610103	31.311	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	1834842	31.199	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	296742	27.826	ng/u1	-0.01
60) Fluorene-d10	15.363	176	1391164	32.193	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	228421	25.987	ng/u1	0.00
73) Anthracene-d10	17.215	188	2135235	32.990	ng/u1	0.00
81) Pyrene-d10	19.515	212	2457994	32.920	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.427	264	2245515	33.384	ng/u1	0.00

Target Compounds Qvalue

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

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