

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022424\
 Data File : BM044308.D
 Acq On : 24 Feb 2024 23:33
 Operator : MA/JU
 Sample : PB159152BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK152

Quant Time: Feb 26 00:59:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Feb 25 22:57:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	188386	20.000	ng/u1	0.00
20) Naphthalene-d8	10.728	136	798765	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.568	164	468020	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.315	188	1031105	20.000	ng/u1	0.00
79) Chrysene-d12	21.515	240	923283	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	1036921	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	40894	6.849	ng/uL	0.00
4) Pyridine-d5	3.769	84	578215	35.459	ng/u1	0.00
7) Phenol-d5	7.087	99	600119	33.254	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	394674	33.772	ng/u1	0.00
11) 2-Chlorophenol-d4	7.451	132	475744	35.002	ng/u1	0.01
15) 4-Methylphenol-d8	8.634	113	454435	33.280	ng/u1	0.00
21) Nitrobenzene-d5	9.092	128	229709	35.369	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	240413	36.283	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.345	165	397279	33.989	ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	663034	33.776	ng/u1	0.01
46) Dimethylphthalate-d6	13.992	166	1321962	36.624	ng/u1	0.01
49) Acenaphthylene-d8	14.263	160	1503944	36.214	ng/u1	0.01
54) 4-Nitrophenol-d4	14.774	143	210603	29.321	ng/u1	0.01
60) Fluorene-d10	15.563	176	1142055	37.381	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	183571	28.693	ng/u1	0.00
73) Anthracene-d10	17.415	188	1809226	36.799	ng/u1	0.00
81) Pyrene-d10	19.715	212	2095546	37.726	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.774	264	2030553	38.216	ng/u1	0.01

Target Compounds Qvalue

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

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