

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022424\
 Data File : BM044306.D
 Acq On : 24 Feb 2024 22:21
 Operator : MA/JU
 Sample : PB159148BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK148

Quant Time: Feb 26 00:56:13 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	228690	20.000	ng/u1	0.00
20) Naphthalene-d8	10.727	136	971482	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.563	164	570605	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.315	188	1244457	20.000	ng/u1	0.00
79) Chrysene-d12	21.515	240	1139285	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	1276737	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	40267	5.555	ng/uL	0.00
4) Pyridine-d5	3.769	84	566912	28.639	ng/u1	0.00
7) Phenol-d5	7.087	99	583807	26.648	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	381218	26.871	ng/u1	0.00
11) 2-Chlorophenol-d4	7.451	132	470202	28.497	ng/u1	0.01
15) 4-Methylphenol-d8	8.633	113	447062	26.970	ng/u1	0.00
21) Nitrobenzene-d5	9.092	128	225955	28.606	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	239470	29.715	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	390029	27.436	ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	641313	26.861	ng/u1	0.01
46) Dimethylphthalate-d6	13.986	166	1304064	29.633	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	1488707	29.402	ng/u1	0.00
54) 4-Nitrophenol-d4	14.774	143	222533	25.412	ng/u1	0.01
60) Fluorene-d10	15.562	176	1124172	30.181	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.680	200	182220	23.599	ng/u1	0.00
73) Anthracene-d10	17.415	188	1777604	29.957	ng/u1	0.00
81) Pyrene-d10	19.715	212	2053515	29.960	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.768	264	1993334	30.469	ng/u1	0.00

Target Compounds Qvalue

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

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