

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030124\
 Data File : BM044474.D
 Acq On : 01 Mar 2024 17:16
 Operator : MA/JU
 Sample : PB159248BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK248

Quant Time: Mar 01 22:30:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 01 16:56:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.069	152	177328	20.000	ng/u1	0.02
20) Naphthalene-d8	10.898	136	741396	20.000	ng/u1	0.02
38) Acenaphthene-d10	14.733	164	435981	20.000	ng/u1	0.02
64) Phenanthrene-d10	17.486	188	910766	20.000	ng/u1	0.01
79) Chrysene-d12	21.680	240	773100	20.000	ng/u1	0.01
88) Perylene-d12	24.215	264	893522	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	34864	6.203	ng/uL	0.00
4) Pyridine-d5	3.834	84	525889	34.262	ng/u1	0.00
7) Phenol-d5	7.228	99	556786	32.776	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.404	67	356729	32.428	ng/u1	0.02
11) 2-Chlorophenol-d4	7.587	132	440420	34.423	ng/u1	0.02
15) 4-Methylphenol-d8	8.792	113	417489	32.481	ng/u1	0.02
21) Nitrobenzene-d5	9.251	128	208488	34.586	ng/u1	0.02
24) 2-Nitrophenol-d4	9.981	143	223001	36.259	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.522	165	363268	33.484	ng/u1	0.02
31) 4-Chloroaniline-d4	11.051	131	599703	32.913	ng/u1	0.02
46) Dimethylphthalate-d6	14.157	166	1159857	34.494	ng/u1	0.01
49) Acenaphthylene-d8	14.427	160	1355381	35.035	ng/u1	0.02
54) 4-Nitrophenol-d4	14.945	143	173584	25.943	ng/u1	0.02
60) Fluorene-d10	15.727	176	1028642	36.144	ng/u1	0.02
65) 4,6-Dinitro-2-methylph...	15.851	200	144495	25.569	ng/u1	0.01
73) Anthracene-d10	17.586	188	1637880	37.716	ng/u1	0.01
81) Pyrene-d10	19.880	212	1844838	39.665	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.050	264	1751209	38.249	ng/u1	0.02

Target Compounds Qvalue

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

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