

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022824\
 Data File : BM044422.D
 Acq On : 29 Feb 2024 02:40
 Operator : MA/JU
 Sample : PB159338BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK338

Quant Time: Feb 29 03:40:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 29 03:37:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	178033	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	812451	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.592	164	474371	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.339	188	970439	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	762642	20.000	ng/u1	0.00
88) Perylene-d12	23.956	264	828318	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	42959	7.613	ng/uL	0.00
4) Pyridine-d5	3.787	84	629829	40.871	ng/u1	0.00
7) Phenol-d5	7.116	99	673017	39.462	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	432884	39.196	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	533805	41.557	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	518699	40.195	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	265438	40.182	ng/u1	0.00
24) 2-Nitrophenol-d4	9.845	143	286260	42.474	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	474215	39.888	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	764650	38.296	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	1463019	39.989	ng/u1	0.00
49) Acenaphthylene-d8	14.286	160	1741936	41.383	ng/u1	0.00
54) 4-Nitrophenol-d4	14.798	143	208081	28.582	ng/u1	0.00
60) Fluorene-d10	15.586	176	1286838	41.557	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	165389	27.467	ng/u1	0.00
73) Anthracene-d10	17.439	188	1936745	41.856	ng/u1	0.00
81) Pyrene-d10	19.733	212	2113806	46.071	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.797	264	1871524	44.094	ng/u1	0.00

Target Compounds Qvalue

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

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