

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042326.D
 Acq On : 17 Oct 2023 18:23
 Operator : MA/JU
 Sample : PB156421BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK421

Quant Time: Oct 18 02:12:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 18 02:04:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	157414	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	714793	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	437851	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	933419	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	815983	20.000	ng/u1	0.00
88) Perylene-d12	24.056	264	850282	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	34251	6.799	ng/uL	0.00
4) Pyridine-d5	3.804	84	479296	32.123	ng/u1	0.00
7) Phenol-d5	7.157	99	567628	32.537	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	395860	33.583	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	404580	33.712	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	440038	31.850	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	206334	33.989	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	219759	33.865	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	372920	31.490	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	604073	32.397	ng/u1	0.00
46) Dimethylphthalate-d6	14.068	166	1289991	33.987	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1411826	33.986	ng/u1	0.00
54) 4-Nitrophenol-d4	14.862	143	220255	28.851	ng/u1	0.00
60) Fluorene-d10	15.645	176	1078275	34.559	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	173184	26.486	ng/u1	0.00
73) Anthracene-d10	17.503	188	1711061	35.785	ng/u1	0.00
81) Pyrene-d10	19.797	212	1996154	40.138	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.897	264	1744871	37.386	ng/u1	0.00

Target Compounds	Qvalue

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

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