

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101623\
 Data File : BM042302.D
 Acq On : 16 Oct 2023 17:08
 Operator : MA/JU
 Sample : PB156300BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK300

Quant Time: Oct 17 01:27:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 15:38:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	131982	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	592186	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	356897	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	754458	20.000	ng/u1	0.00
79) Chrysene-d12	21.592	240	646411	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	710287	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	30958	7.329	ng/uL	0.00
4) Pyridine-d5	3.805	84	467214	37.347	ng/u1	0.00
7) Phenol-d5	7.163	99	493923	33.768	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.357	67	364979	36.929	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	362765	36.053	ng/u1	0.00
15) 4-Methylphenol-d8	8.722	113	386083	33.330	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	182194	36.226	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	191748	35.666	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	321415	32.760	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	543432	35.178	ng/u1	0.00
46) Dimethylphthalate-d6	14.075	166	1131338	36.568	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1273636	37.614	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	202816	32.593	ng/u1	0.00
60) Fluorene-d10	15.651	176	953694	37.499	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	150878	28.548	ng/u1	0.00
73) Anthracene-d10	17.510	188	1490218	38.559	ng/u1	0.00
81) Pyrene-d10	19.804	212	1708203	43.358	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.909	264	1607851	41.241	ng/u1	0.00

Target Compounds Qvalue

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

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