

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042316.D
 Acq On : 17 Oct 2023 10:11
 Operator : MA/JU
 Sample : PB156333BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK333

Quant Time: Oct 18 01:42:02 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	8.010	152	168055	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	761003	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	468644	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	1020830	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	915062	20.000	ng/u1	-0.01
88) Perylene-d12	24.062	264	967550	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	31035	5.770	ng/uL	0.00
4) Pyridine-d5	3.805	84	452278	28.393	ng/u1	0.00
7) Phenol-d5	7.157	99	490544	26.338	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	355219	28.227	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	362836	28.319	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	387364	26.262	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	185021	28.628	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	194566	28.162	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	328735	26.073	ng/u1	0.00
31) 4-Chloroaniline-d4	10.992	131	551611	27.787	ng/u1	0.00
46) Dimethylphthalate-d6	14.069	166	1168767	28.770	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1290731	29.029	ng/u1	0.00
54) 4-Nitrophenol-d4	14.863	143	203283	24.878	ng/u1	-0.01
60) Fluorene-d10	15.645	176	980661	29.365	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	154605	21.620	ng/u1	0.00
73) Anthracene-d10	17.504	188	1537384	29.399	ng/u1	0.00
81) Pyrene-d10	19.798	212	1835661	32.914	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.897	264	1684044	31.710	ng/u1	-0.02

Target Compounds	Qvalue

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

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