

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101623\
 Data File : BM042295.D
 Acq On : 16 Oct 2023 12:50
 Operator : MA/JU
 Sample : PB156264BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK264

Quant Time: Oct 17 01:17:31 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	156195	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	710898	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	427792	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.409	188	908029	20.000	ng/u1	0.00
79) Chrysene-d12	21.591	240	803064	20.000	ng/u1	0.00
88) Perylene-d12	24.074	264	868995	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	29025	5.806	ng/uL	0.00
4) Pyridine-d5	3.804	84	421610	28.477	ng/u1	0.00
7) Phenol-d5	7.157	99	471804	27.256	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	336288	28.751	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	331677	27.853	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	363753	26.534	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	170445	28.231	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	179467	27.807	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	305144	25.908	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	494508	26.666	ng/u1	0.00
46) Dimethylphthalate-d6	14.074	166	1060216	28.590	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1155206	28.462	ng/u1	0.00
54) 4-Nitrophenol-d4	14.868	143	176798	23.703	ng/u1	0.00
60) Fluorene-d10	15.651	176	882008	28.933	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	133255	20.949	ng/u1	0.00
73) Anthracene-d10	17.509	188	1359003	29.216	ng/u1	0.00
81) Pyrene-d10	19.803	212	1567305	32.022	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.909	264	1430897	29.999	ng/u1	0.00

Target Compounds Qvalue

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

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