

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
 Data File : BM039651.D
 Acq On : 25 Apr 2023 22:05
 Operator : CG/JU
 Sample : PB152398BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK398

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/26/2023
 Supervised By : Jagrut Upadhyay 04/26/2023

Quant Time: Apr 25 23:47:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 25 22:38:31 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.807	152	211216	20.000	ng/u1	0.00
20) Naphthalene-d8	10.613	136	835171	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.465	164	590093	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.218	188	1228993	20.000	ng/u1	0.00
79) Chrysene-d12	21.418	240	1075452	20.000	ng/u1	0.00
88) Perylene-d12	23.777	264	1087413	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.201	96	29149	5.297	ng/uL	0.00
4) Pyridine-d5	3.625	84	436480	28.230	ng/u1	0.00
7) Phenol-d5	6.989	99	504919	26.119	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.142	67	392224	30.697	ng/u1	0.00
11) 2-Chlorophenol-d4	7.336	132	434422	28.647	ng/u1	0.00
15) 4-Methylphenol-d8	8.536	113	408942	26.261	ng/u1	0.00
21) Nitrobenzene-d5	8.978	128	190527	28.356	ng/u1	0.00
24) 2-Nitrophenol-d4	9.701	143	213319	28.273	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.248	165	326264	25.132	ng/u1	0.00
31) 4-Chloroaniline-d4	10.766	131	495519	25.556	ng/u1	0.00
46) Dimethylphthalate-d6	13.883	166	1392988	29.270	ng/u1	0.00
49) Acenaphthylene-d8	14.160	160	1568342	29.335	ng/u1	0.00
54) 4-Nitrophenol-d4	14.730	143	129839m	18.027	ng/u1	0.02
60) Fluorene-d10	15.459	176	1170872	30.030	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.607	200	167745	20.539	ng/u1	0.00
73) Anthracene-d10	17.318	188	1814392	30.698	ng/u1	0.00
81) Pyrene-d10	19.618	212	2088079	29.219	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.624	264	1757704	29.940	ng/u1	0.00

Target Compounds Qvalue

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

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