

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042423\
 Data File : BM039630.D
 Acq On : 25 Apr 2023 00:08
 Operator : CG/JU
 Sample : PB152352BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK352

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/25/2023
 Supervised By : Jagrut Upadhyay 04/25/2023

Quant Time: Apr 25 01:19:15 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Apr 23 00:18:16 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.807	152	195946	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.619	136	873373	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.466	164	595130	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.224	188	1269968	20.000	ng/u1	0.00
79) Chrysene-d12	21.418	240	1164648	20.000	ng/u1	0.00
88) Perylene-d12	23.777	264	1288707	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.208	96	32894	6.444	ng/uL	0.00
4) Pyridine-d5	3.631	84	383290	26.722	ng/u1	0.00
7) Phenol-d5	6.996	99	495587	27.634	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.143	67	360236	30.390	ng/u1	0.00
11) 2-Chlorophenol-d4	7.343	132	417819	29.700	ng/u1	0.00
15) 4-Methylphenol-d8	8.537	113	405720	28.085	ng/u1	-0.01
21) Nitrobenzene-d5	8.984	128	209380	29.799	ng/u1	0.00
24) 2-Nitrophenol-d4	9.707	143	229791	29.125	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.254	165	347455	25.593	ng/u1	0.00
31) 4-Chloroaniline-d4	10.772	131	489546	24.144	ng/u1	0.00
46) Dimethylphthalate-d6	13.883	166	1471844	30.666	ng/u1	0.00
49) Acenaphthylene-d8	14.160	160	1652913	30.655	ng/u1	0.00
54) 4-Nitrophenol-d4	14.730	143	145543m	20.036	ng/u1	0.00
60) Fluorene-d10	15.466	176	1236342	31.440	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.607	200	195127	23.121	ng/u1	0.00
73) Anthracene-d10	17.318	188	1952348	31.966	ng/u1	0.00
81) Pyrene-d10	19.624	212	2280322	29.465	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.624	264	2193365	31.525	ng/u1	0.00

Target Compounds Qvalue

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

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