

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042423\
 Data File : BM039617.D
 Acq On : 24 Apr 2023 16:20
 Operator : CG/JU
 Sample : PB152350BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK350

Manual Integrations
 APPROVED

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

Quant Time: Apr 25 00:51:36 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	7.813	152	203768	20.000	ng/u1	0.00
20) Naphthalene-d8	10.619	136	891384	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.472	164	562708	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.224	188	1189109	20.000	ng/u1	0.00
79) Chrysene-d12	21.424	240	1014462	20.000	ng/u1	0.00
88) Perylene-d12	23.783	264	1000024	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.208	96	32266	6.078	ng/uL	0.00
4) Pyridine-d5	3.631	84	365327	24.492	ng/u1	0.00
7) Phenol-d5	6.996	99	500396	26.831	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.148	67	370084	30.022	ng/u1	0.00
11) 2-Chlorophenol-d4	7.343	132	425863	29.110	ng/u1	0.00
15) 4-Methylphenol-d8	8.542	113	417078	27.763	ng/u1	0.00
21) Nitrobenzene-d5	8.984	128	212333	29.608	ng/u1	0.00
24) 2-Nitrophenol-d4	9.707	143	236110	29.321	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.254	165	350924	25.327	ng/u1	0.00
31) 4-Chloroaniline-d4	10.772	131	494591	23.900	ng/u1	0.00
46) Dimethylphthalate-d6	13.883	166	1391543	30.663	ng/u1	0.00
49) Acenaphthylene-d8	14.166	160	1549518	30.393	ng/u1	0.00
54) 4-Nitrophenol-d4	14.736	143	117530m	17.112	ng/u1	0.00
60) Fluorene-d10	15.466	176	1157252	31.125	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.607	200	179244	22.683	ng/u1	0.00
73) Anthracene-d10	17.324	188	1823340	31.884	ng/u1	0.00
81) Pyrene-d10	19.624	212	2085450	30.937	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.630	264	1700026	31.488	ng/u1	0.00

Target Compounds Qvalue

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

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