

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
 Data File : BM039541.D
 Acq On : 21 Apr 2023 01:00
 Operator : CG/JU
 Sample : PB152246BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK246

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/21/2023
 Supervised By : Jagrut Upadhyay 04/21/2023

Quant Time: Apr 21 02:33:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 19 08:18:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.819	152	216008	20.000	ng/u1	0.00
20) Naphthalene-d8	10.631	136	1008824	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.477	164	674032	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.230	188	1413011	20.000	ng/u1	0.00
79) Chrysene-d12	21.424	240	987120	20.000	ng/u1	0.00
88) Perylene-d12	23.783	264	1005810	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	33594	5.970	ng/uL	0.00
4) Pyridine-d5	3.637	84	350309	22.154	ng/u1	0.00
7) Phenol-d5	7.001	99	540749	27.352	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.154	67	372038	28.471	ng/u1	0.00
11) 2-Chlorophenol-d4	7.354	132	456351	29.426	ng/u1	0.00
15) 4-Methylphenol-d8	8.548	113	459247	28.837	ng/u1	-0.01
21) Nitrobenzene-d5	8.995	128	228579	28.163	ng/u1	0.00
24) 2-Nitrophenol-d4	9.719	143	258550	28.370	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.260	165	415253	26.480	ng/u1	0.00
31) 4-Chloroaniline-d4	10.783	131	432985	18.487	ng/u1	0.00
46) Dimethylphthalate-d6	13.895	166	1569345	28.870	ng/u1	0.00
49) Acenaphthylene-d8	14.171	160	1762428	28.860	ng/u1	0.00
54) 4-Nitrophenol-d4	14.719	143	150733m	18.322	ng/u1	0.00
60) Fluorene-d10	15.471	176	1335479	29.986	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.607	200	184996	19.702	ng/u1	0.00
73) Anthracene-d10	17.330	188	2056951	30.270	ng/u1	0.00
81) Pyrene-d10	19.630	212	2197234	33.498	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.630	264	1645930	30.310	ng/u1	-0.01

Target Compounds Qvalue

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

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