

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101024\
 Data File : BM047999.D
 Acq On : 11 Oct 2024 04:27
 Operator : RC/JU
 Sample : PB163767BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK767

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/11/2024
 Supervised By :mohammad ahmed 10/14/2024

Quant Time: Oct 11 05:11:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 18:01:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	9975	0.400	ng/ul	0.00
4) Naphthalene-d8	10.578	136	29674	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.422	164	14447	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.161	188	27238m	0.400	ng/ul	0.00
17) Chrysene-d12	21.336	240	18881	0.400	ng/ul	0.00
23) Perylene-d12	23.590	264	15430m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	61698	4.814	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.184	152	9172	0.232	ng/ul	0.00
18) Fluoranthene-d10	19.182	212	17537	0.303	ng/ul	0.00

Target Compounds	Qvalue

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

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