

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092624\
 Data File : BM047666.D
 Acq On : 27 Sep 2024 05:29
 Operator : RC/JU
 Sample : PB163687BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK687

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 09/27/2024
 Supervised By :mohammad ahmed 09/28/2024

Quant Time: Sep 27 06:45:22 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 : Wed Sep 25 16:27:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.826	152	5175	0.400	ng/u1	0.00
4) Naphthalene-d8	10.618	136	15035	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.456	164	7356	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.192	188	14247m	0.400	ng/u1	0.00
17) Chrysene-d12	21.356	240	10478	0.400	ng/u1	0.00
23) Perylene-d12	23.630	264	10226m	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	41345	6.218	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.207	152	5660	0.282	ng/u1	0.00
18) Fluoranthene-d10	19.211	212	10106	0.315	ng/u1	0.00

Target Compounds	Qvalue

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

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