

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111324\
 Data File : BM048643.D
 Acq On : 13 Nov 2024 20:58
 Operator : RC/JU
 Sample : PB164898BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK898

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/14/2024
 Supervised By :mohammad ahmed 11/15/2024

Quant Time: Nov 14 02:25:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 11:38:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.662	152	5857	0.400	ng/u1	-0.08
4) Naphthalene-d8	10.442	136	18284	0.400	ng/u1	#-0.05
9) Acenaphthene-d10	14.308	164	9618	0.400	ng/u1	-0.02
13) Phenanthrene-d10	17.053	188	20412m	0.400	ng/u1	-0.03
17) Chrysene-d12	21.266	240	15963m	0.400	ng/u1	0.00
23) Perylene-d12	23.460	264	12256m	0.400	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	35950	5.088	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.069	152	7189	0.304	ng/u1	-0.03
18) Fluoranthene-d10	19.086	212	15977	0.357	ng/u1	-0.02

Target Compounds	Qvalue

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

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