

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102124\
 Data File : BM048165.D
 Acq On : 21 Oct 2024 21:50
 Operator : RC/JU
 Sample : PB164049BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
ClientSampleId :
 SBLK049

Manual Integrations
APPROVED
 Reviewed By :Yogesh Patel 10/22/2024
 Supervised By :mohammad ahmed 10/23/2024

Quant Time: Oct 21 23:14:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 23:11:55 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.771	152	7466	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.557	136	23449	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.405	164	11889	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.145	188	23280m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.350	240	15788m	0.400	ng/ul	0.01
23) Perylene-d12	23.589	264	12186m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.235	96	44340	4.369	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.179	152	8296	0.254	ng/ul	-0.02
18) Fluoranthene-d10	19.174	212	16563	0.360	ng/ul	-0.02

Target Compounds Qvalue

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

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