

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111624\
 Data File : BM048748.D
 Acq On : 16 Nov 2024 18:53
 Operator : RC/JU
 Sample : P4803-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A0AR9

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/19/2024
 Supervised By :mohammad ahmed 11/19/2024

Quant Time: Nov 17 22:49:17 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 : Sun Nov 17 22:31:02 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.629	152	3665	0.400	ng/ul	-0.05
4) Naphthalene-d8	10.410	136	10782	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.276	164	5720	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.024	188	10697m	0.400	ng/ul	-0.02
17) Chrysene-d12	21.240	240	8280m	0.400	ng/ul	0.01
23) Perylene-d12	23.432	264	6354m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.089	96	21494	4.861	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.021	152	4912	0.352	ng/ul	-0.03
18) Fluoranthene-d10	19.059	212	10007	0.430	ng/ul	-0.01
Target Compounds						Qvalue
22) Chrysene	21.275	228	910	0.028	ng/ul#	81
24) Benzo(b)fluoranthene	22.763	252	592	0.030	ng/ul#	1
25) Benzo(k)fluoranthene	22.809	252	554	0.024	ng/ul#	1

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

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