

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
 Data File : BM048498.D
 Acq On : 07 Nov 2024 04:25
 Operator : RC/JU
 Sample : P4634-17
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CC0R7

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
 Supervised By :mohammad ahmed 11/11/2024

Quant Time: Nov 07 04:52:51 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 06 14:39:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.674	152	4656	0.400	ng/ul	0.00
4) Naphthalene-d8	10.458	136	14688	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.322	164	7803	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.065	188	15444m	0.400	ng/ul	-0.02
17) Chrysene-d12	21.257	240	12320m	0.400	ng/ul	0.00
23) Perylene-d12	23.475	264	12157m	0.400	ng/ul	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	25033	4.457	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.069	152	5679	0.299	ng/ul	0.00
18) Fluoranthene-d10	19.095	212	11290	0.326	ng/ul	-0.01

Target Compounds	Qvalue

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

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