

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020124\
 Data File : BM043946.D
 Acq On : 01 Feb 2024 14:34
 Operator : MA/JU
 Sample : P1239-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0FB9

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 02/02/2024
 Supervised By :Sohil Jodhani 02/02/2024

Quant Time: Feb 01 15:14:25 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM013124.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Jan 31 22:50:05 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.926	152	9291	0.400	ng/ul	0.00
4) Naphthalene-d8	10.726	136	27165	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.566	164	13613	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.317	188	28359m	0.400	ng/ul	0.00
17) Chrysene-d12	21.514	240	21916m	0.400	ng/ul	0.00
23) Perylene-d12	23.926	264	22836m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.365	96	49669	4.155	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.316	152	11648	0.327	ng/ul	0.00
18) Fluoranthene-d10	19.350	212	24717m	0.359	ng/ul	0.00
Target Compounds						Qvalue
14) Pentachlorophenol	16.959	266	494	0.074	ng/ul	98
24) Benzo(b)fluoranthene	23.192	252	2462m	0.028	ng/ul	
25) Benzo(k)fluoranthene	23.242	252	2427m	0.027	ng/ul	
27) Indeno(1,2,3-cd)pyrene	26.423	276	2949m	0.029	ng/ul	
28) Dibenzo(a,h)anthracene	26.457	278	2305m	0.028	ng/ul	
29) Benzo(g,h,i)perylene	27.185	276	2544m	0.029	ng/ul	

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

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