

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035543.D
 Acq On : 10 Dec 2024 22:06
 Operator : RC/JU
 Sample : P5066-18
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E28N6

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/12/2024
 Supervised By :mohammad ahmed 12/12/2024

Quant Time: Dec 10 22:36:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 10 16:10:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.293	152	3422	0.400	ng/ul	0.00
4) Naphthalene-d8	10.036	136	9774	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	6418	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	12879	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.961	240	8324	0.400	ng/ul	# 0.00
23) Perylene-d12	23.048	264	7607m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.959	96	10377	3.318	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.642	152	2577	0.180	ng/ul	0.00
18) Fluoranthene-d10	18.767	212	5033	0.219	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	16.759	178	785	0.023	ng/ul#	81
19) Fluoranthene	18.800	202	1329	0.046	ng/ul#	81
20) Pyrene	19.167	202	1149	0.038	ng/ul#	72
22) Chrysene	20.996	228	649	0.022	ng/ul#	76
24) Benzo(b)fluoranthene	22.437	252	903m	0.036	ng/ul	

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

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