

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035540.D
 Acq On : 10 Dec 2024 20:18
 Operator : RC/JU
 Sample : P5066-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E28M7

Manual Integrations
 APPROVED

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

Quant Time: Dec 10 22:05:12 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	3013	0.400	ng/ul	0.00
4) Naphthalene-d8	10.037	136	8312	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	13.949	164	5253	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	11091m	0.400	ng/ul	0.00
17) Chrysene-d12	20.961	240	8277	0.400	ng/ul	# 0.00
23) Perylene-d12	23.048	264	8063m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.959	96	13259	4.815	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	3027	0.249	ng/ul	0.00
18) Fluoranthene-d10	18.768	212	6531	0.286	ng/ul	0.00
Target Compounds						
						Qvalue
7) 2-Methylnaphthalene	11.714	142	362	0.025	ng/ul	88
8) 1-Methylnaphthalene	11.939	142	359	0.024	ng/ul	85
15) Phenanthrene	16.760	178	1548	0.053	ng/ul#	84
22) Chrysene	20.994	228	1156m	0.040	ng/ul	
24) Benzo(b)fluoranthene	22.440	252	595m	0.022	ng/ul	
29) Benzo(g,h,i)perylene	25.688	276	826m	0.032	ng/ul	

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

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