

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035552.D
 Acq On : 11 Dec 2024 04:06
 Operator : RC/JU
 Sample : P5066-21MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E28N8MS

Manual Integrations
 APPROVED

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

Quant Time: Dec 11 04:28:43 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	3490	0.400	ng/ul	0.00
4) Naphthalene-d8	10.037	136	9749	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	13.949	164	6337	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	11114	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.961	240	7282	0.400	ng/ul	# 0.00
23) Perylene-d12	23.057	264	7625	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	1452	0.455	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	4625	0.324	ng/ul	0.00
18) Fluoranthene-d10	18.768	212	9129m	0.454	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.993	88	4218	1.218	ng/ul	91
5) Naphthalene	10.086	128	8256	0.336	ng/ul	99
7) 2-Methylnaphthalene	11.714	142	7063	0.415	ng/ul	97
8) 1-Methylnaphthalene	11.939	142	6814	0.396	ng/ul	96
10) Acenaphthylene	13.662	152	8194	0.326	ng/ul	93
11) Acenaphthene	14.014	153	6474	0.337	ng/ul	97
12) Fluorene	15.013	166	8818	0.391	ng/ul#	90
14) Pentachlorophenol	16.384	266	60	0.023	ng/ul#	1
15) Phenanthrene	16.755	178	29420	1.001	ng/ul#	86
16) Anthracene	16.848	178	9753	0.358	ng/ul#	71
19) Fluoranthene	18.800	202	16669	0.657	ng/ul#	90
20) Pyrene	19.167	202	17110	0.648	ng/ul#	90
21) Benzo(a)anthracene	20.947	228	10071	0.399	ng/ul#	54
22) Chrysene	20.999	228	15839	0.623	ng/ul#	62
24) Benzo(b)fluoranthene	22.443	252	11063	0.439	ng/ul#	1
25) Benzo(k)fluoranthene	22.484	252	9991	0.384	ng/ul#	1
26) Benzo(a)pyrene	22.966	252	9316	0.404	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	25.084	276	11474m	0.369	ng/ul	
28) Dibenzo(a,h)anthracene	25.104	278	8218m	0.337	ng/ul	
29) Benzo(g,h,i)perylene	25.698	276	9926	0.405	ng/ul#	67

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

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