

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111523\
 Data File : BN028783.D
 Acq On : 19 Nov 2023 19:26
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4322

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/20/2023
 Supervised By :mohammad ahmed 11/20/2023

Quant Time: Nov 20 02:49:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 20 01:23:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.896	152	5720m	0.400	ng/ul	0.10
4) Naphthalene-d8	10.528	136	14946m	0.400	ng/ul	0.04
9) Acenaphthene-d10	14.325	164	9637	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.102	188	12123	0.400	ng/ul	# 0.01
17) Chrysene-d12	21.301	240	10803	0.400	ng/ul	# 0.00
23) Perylene-d12	23.581	264	11324	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.125	96	2905m	0.460	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.151	152	6310m	0.280	ng/ul	0.04
18) Fluoranthene-d10	19.122	212	15516	0.446	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.154	88	3002m	0.430	ng/ul	
5) Naphthalene	10.572	128	12978	0.310	ng/ul#	93
7) 2-Methylnaphthalene	12.239	142	6969m	0.253	ng/ul	
8) 1-Methylnaphthalene	12.376	142	8604	0.307	ng/ul	98
10) Acenaphthylene	14.057	152	18250	0.379	ng/ul	97
11) Acenaphthene	14.385	153	15470	0.437	ng/ul	98
12) Fluorene	15.412	166	12408	0.301	ng/ul	98
14) Pentachlorophenol	16.769	266	753	0.259	ng/ul	99
15) Phenanthrene	17.140	178	18938	0.508	ng/ul	95
16) Anthracene	17.258	178	9084	0.266	ng/ul	93
19) Fluoranthene	19.154	202	22224	0.484	ng/ul	97
20) Pyrene	19.512	202	26349	0.553	ng/ul#	94
21) Benzo(a)anthracene	21.290	228	11155	0.261	ng/ul	98
22) Chrysene	21.336	228	17444	0.418	ng/ul	99
24) Benzo(b)fluoranthene	22.891	252	16425m	0.386	ng/ul	
25) Benzo(k)fluoranthene	22.938	252	19248	0.432	ng/ul#	88
26) Benzo(a)pyrene	23.485	252	17822	0.444	ng/ul#	61
27) Indeno(1,2,3-cd)pyrene	25.931	276	21259	0.477	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.967	278	15763	0.443	ng/ul#	93
29) Benzo(g,h,i)perylene	26.635	276	19503	0.534	ng/ul	96

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111523\
Data File : BN028783.D
Acq On : 19 Nov 2023 19:26
Operator : MA/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTD0.4322

Quant Time: Nov 20 02:49:38 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 20 01:23:26 2023
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Yogesh Patel 11/20/2023
Supervised By :mohammad ahmed 11/20/2023

