

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012831.D
 Acq On : 02 Dec 2020 22:33
 Operator : CG/JU
 Sample : L4865-03
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B23

Manual Integrations
 APPROVED

mohammad
 12/3/2020 3:43:50 PM

Quant Time: Dec 03 01:17:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 03 00:42:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	1080	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	4353	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	2550	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	5530	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	4818	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	5225	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	645	0.10	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	1647	0.11	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.22	128	1358	0.104	ng/ul	95
5) 2-Methylnaphthalene	11.86	142	951	0.114	ng/ul	97
7) Acenaphthylene	13.79	152	588	0.047	ng/ul#	79
8) Acenaphthene	14.14	153	373	0.036	ng/ul	96
9) Fluorene	15.13	166	364	0.032	ng/ul#	90
12) Phenanthrene	16.87	178	8756	0.455	ng/ul	98
13) Anthracene	16.96	178	1659	0.099	ng/ul	96
15) Fluoranthene	18.90	202	19322	0.870	ng/ul	96
17) Pyrene	19.26	202	18128	0.841	ng/ul	99
18) Benzo(a)anthracene	21.03	228	9206	0.533	ng/ul	100
19) Chrysene	21.08	228	11190	0.519	ng/ul	99
21) Benzo(b)fluoranthene	22.53	252	15226m	0.633	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	5717m	0.206	ng/ul	
23) Benzo(a)pyrene	23.06	252	10167	0.449	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.21	276	7521	0.254	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.22	278	1830	0.077	ng/ul#	76
26) Benzo(a,h,i)perylene	25.85	276	6753	0.254	ng/ul	97

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

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