

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012838.D
 Acq On : 03 Dec 2020 02:45
 Operator : CG/JU
 Sample : L4865-19DL 10X
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AZ9DL

Manual Integrations
 APPROVED

mohammad
 12/3/2020 3:44:39 PM

Quant Time: Dec 03 03:22:28 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	799	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	3157	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1901	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	3934	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3203	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	3264	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	54	0.01	ng/ul	0.02
14) Fluoranthene-d10	18.87	212	132	0.01	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.87	178	4548	0.332	ng/ul	98
13) Anthracene	16.96	178	647	0.054	ng/ul#	84
15) Fluoranthene	18.90	202	17734	1.122	ng/ul	99
17) Pyrene	19.26	202	15492	1.081	ng/ul	100
18) Benzo(a)anthracene	21.03	228	7836	0.683	ng/ul	99
19) Chrysene	21.08	228	11131	0.776	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	15744m	1.049	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	5740m	0.331	ng/ul	
23) Benzo(a)pyrene	23.06	252	9659	0.684	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.22	276	7736	0.419	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.22	278	1813	0.122	ng/ul#	84
26) Benzo(g,h,i)perylene	25.84	276	7149	0.430	ng/ul	98

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

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