

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012822.D
 Acq On : 02 Dec 2020 16:32
 Operator : CG/JU
 Sample : L4865-02
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B20

Manual Integrations
 APPROVED

mohammad
 12/3/2020 3:43:40 PM

Quant Time: Dec 02 17:05:38 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 : Wed Dec 02 12:19:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	716	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	3044	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1783	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	3812	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3504	0.40	ng/ul	0.00
20) Perylene-d12	23.16	264	3237	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	679	0.15	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	1632	0.16	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.23	128	1798	0.196	ng/ul	97
5) 2-Methylnaphthalene	11.86	142	596	0.102	ng/ul	98
7) Acenaphthylene	13.79	152	2022	0.233	ng/ul	98
8) Acenaphthene	14.14	153	1263	0.172	ng/ul	99
9) Fluorene	15.13	166	1188	0.147	ng/ul#	92
12) Phenanthrene	16.87	178	75440	5.687	ng/ul	96
13) Anthracene	16.96	178	5887	0.511	ng/ul	98
15) Fluoranthene	18.90	202	197615	12.904	ng/ul	97
17) Pyrene	19.27	202	164415	10.490	ng/ul	100
18) Benzo(a)anthracene	21.03	228	65413	5.210	ng/ul	98
19) Chrysene	21.09	228	81241	5.180	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	97347m	6.537	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	33939m	1.976	ng/ul	
23) Benzo(a)pyrene	23.07	252	62517m	4.461	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.23	276	41534	2.266	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.23	278	10190	0.692	ng/ul	97
26) Benzo(a,h,i)perylene	25.85	276	35894	2.176	ng/ul	96

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

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