

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012921.D
 Acq On : 06 Dec 2020 17:17
 Operator : CG/JU
 Sample : L4864-14
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B72

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:11:44 PM

Quant Time: Dec 07 03:29:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 03:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1824	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	7537	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	4338	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	9127	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	7155	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	5826	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1760	0.16	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	3926	0.16	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	950	0.044	ng/ul#	94
5) 2-Methylnaphthalene	11.84	142	467	0.033	ng/ul	98
7) Acenaphthylene	13.77	152	1356	0.066	ng/ul#	95
8) Acenaphthene	14.12	153	1385	0.082	ng/ul	99
9) Fluorene	15.12	166	1479	0.077	ng/ul	99
12) Phenanthrene	16.85	178	40613	1.327	ng/ul	97
13) Anthracene	16.95	178	7894	0.303	ng/ul	96
15) Fluoranthene	18.88	202	137536	3.782	ng/ul	99
17) Pyrene	19.26	202	117383	3.473	ng/ul	100
18) Benzo(a)anthracene	21.02	228	71414	2.412	ng/ul	98
19) Chrysene	21.07	228	76941	2.352	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	96171m	3.568	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	31303m	1.051	ng/ul	
23) Benzo(a)pyrene	23.05	252	58584m	2.342	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.19	276	34376	1.046	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.20	278	9696	0.373	ng/ul#	96
26) Benzo(a,h,i)perylene	25.82	276	29235	0.984	ng/ul	97

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

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