

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012905.D
 Acq On : 06 Dec 2020 07:05
 Operator : CG/JU
 Sample : L4859-15
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B27

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:10:51 PM

Quant Time: Dec 07 11:13:36 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 : Fri Dec 04 10:02:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1066	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	4225	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	2404	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5221	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4825	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	4634	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	995	0.16	ng/ul	-0.01
14) Fluoranthene-d10	18.86	212	2388	0.17	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.21	128	1431	0.117	ng/ul	99
5) 2-Methylnaphthalene	11.84	142	319	0.040	ng/ul	96
7) Acenaphthylene	13.77	152	2976	0.262	ng/ul	97
8) Acenaphthene	14.12	153	288	0.031	ng/ul#	92
9) Fluorene	15.12	166	374	0.035	ng/ul#	92
12) Phenanthrene	16.86	178	11273	0.644	ng/ul	98
13) Anthracene	16.95	178	2156	0.145	ng/ul	98
15) Fluoranthene	18.89	202	49858	2.397	ng/ul	95
17) Pyrene	19.25	202	50181	2.202	ng/ul	100
18) Benzo(a)anthracene	21.02	228	27032	1.354	ng/ul	100
19) Chrysene	21.07	228	30785	1.396	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	46864m	2.186	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	16492m	0.696	ng/ul	
23) Benzo(a)pyrene	23.04	252	30464m	1.531	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.19	276	20726	0.793	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.20	278	4852	0.235	ng/ul	98
26) Benzo(a,h,i)perylene	25.82	276	18267	0.773	ng/ul	97

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

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