

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013000.D
 Acq On : 09 Dec 2020 07:06
 Operator : CG/JU
 Sample : L4862-17
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B70

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:25:57 PM

Quant Time: Dec 09 09:10:48 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 : Wed Dec 09 07:33:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1135	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.14	136	4639	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.05	164	2713	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5588m	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4099	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	3594	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	3115	0.45	ng/ul	-0.01
14) Fluoranthene-d10	18.85	212	6079	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	443	0.033	ng/ul#	77
5) 2-Methylnaphthalene	11.82	142	242	0.027	ng/ul	99
7) Acenaphthylene	13.76	152	544	0.042	ng/ul#	77
12) Phenanthrene	16.84	178	5912	0.316	ng/ul	98
13) Anthracene	16.94	178	1007	0.063	ng/ul#	90
15) Fluoranthene	18.88	202	22702	1.020	ng/ul	98
17) Pyrene	19.25	202	21112	1.090	ng/ul	99
18) Benzo(a)anthracene	21.01	228	10736	0.633	ng/ul#	87
19) Chrysene	21.06	228	13106	0.699	ng/ul#	95
21) Benzo(b)fluoranthene	22.51	252	19296m	1.161	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	6487m	0.353	ng/ul	
23) Benzo(a)pyrene	23.04	252	11847	0.768	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.18	276	8711	0.430	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.18	278	2628	0.164	ng/ul#	70
26) Benzo(g,h,i)perylene	25.81	276	10724	0.585	ng/ul	96

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

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