

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008274.D
 Acq On : 21 Oct 2019 09:50
 Operator : JU
 Sample : K5199-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPJ6

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:26:51 PM

Quant Time: Oct 21 15:33:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 15:18:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1842	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8808	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5579	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	12747m	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13363	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	15853	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	4270	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	13477	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	1259	0.050	ng/ul#	86
5) 2-Methylnaphthalene	12.03	142	722	0.042	ng/ul	100
7) Acenaphthylene	13.94	152	4090	0.154	ng/ul	96
8) Acenaphthene	14.29	153	1266	0.061	ng/ul	97
9) Fluorene	15.28	166	1819	0.078	ng/ul#	96
12) Phenanthrene	17.01	178	40138	1.076	ng/ul	99
13) Anthracene	17.11	178	9254	0.282	ng/ul	99
15) Fluoranthene	19.04	202	122192	2.479	ng/ul	100
17) Pyrene	19.40	202	127188	2.390	ng/ul	99
18) Benzo(a)anthracene	21.16	228	87251	1.796	ng/ul	98
19) Chrysene	21.21	228	95272	1.598	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	114086m	2.211	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	42189m	0.722	ng/ul	
23) Benzo(a)pyrene	23.26	252	81696	1.504	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.51	276	45958	0.719	ng/ul	100
25) Dibenzo(a,h)anthracene	25.51	278	12652	0.250	ng/ul	96
26) Benzo(a,h,i)perylene	26.16	276	39886	0.680	ng/ul	98

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

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