

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008658.D
 Acq On : 20 Nov 2019 13:43
 Operator : JU
 Sample : K5784-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLNBO

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:06:07 PM

Quant Time: Nov 20 14:34:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 20 04:45:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3967	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	14728	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8232	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	16757	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13209	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13191	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4165	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	9603	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	10192	0.230	ng/ul	98
5) 2-Methylnaphthalene	12.04	142	4149	0.134	ng/ul	100
7) Acenaphthylene	13.96	152	6913	0.152	ng/ul	96
8) Acenaphthene	14.30	153	3810	0.113	ng/ul	99
9) Fluorene	15.29	166	3509	0.088	ng/ul#	92
12) Phenanthrene	17.03	178	77615	1.373	ng/ul	98
13) Anthracene	17.11	178	16743	0.314	ng/ul	98
15) Fluoranthene	19.05	202	180923	2.606	ng/ul	97
17) Pyrene	19.41	202	172147	3.249	ng/ul	100
18) Benzo(a)anthracene	21.16	228	88208	1.661	ng/ul	93
19) Chrysene	21.21	228	90644	1.737	ng/ul	97
21) Benzo(b)fluoranthene	22.71	252	102659m	1.887	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	37247m	0.690	ng/ul	
23) Benzo(a)pyrene	23.25	252	74199	1.453	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.49	276	50856	0.852	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.50	278	11832	0.245	ng/ul#	70
26) Benzo(a,h,i)perylene	26.14	276	52683	1.055	ng/ul	97

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

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