

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011722.D
 Acq On : 05 Sep 2020 04:24
 Operator : CG/JU
 Sample : L3843-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD0

Manual Integrations
 APPROVED

mohammad
 9/8/2020 1:47:36 PM

Quant Time: Sep 05 06:33:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 05 05:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1652	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	1007	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	405	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	501m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	543	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	480m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	4097	2.87	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	8745	6.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	1831	0.624	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	296	0.160	ng/ul	100
7) Acenaphthylene	14.10	152	3814	1.871	ng/ul	98
8) Acenaphthene	14.44	153	473	0.298	ng/ul	96
9) Fluorene	15.43	166	646	0.345	ng/ul#	88
12) Phenanthrene	17.16	178	3739	2.206	ng/ul	98
13) Anthracene	17.25	178	4814m	3.453	ng/ul	
15) Fluoranthene	19.18	202	27148	13.925	ng/ul	97
17) Pyrene	19.54	202	46586	21.950	ng/ul#	95
18) Benzo(a)anthracene	21.29	228	13504	7.412	ng/ul	98
19) Chrysene	21.35	228	23930	9.739	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	43896m	21.040	ng/ul	
22) Benzo(k)fluoranthene	22.93	252	16334m	7.081	ng/ul	
23) Benzo(a)pyrene	23.46	252	23662	13.498	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.82	276	14643	6.177	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.82	278	4354	2.253	ng/ul#	83
26) Benzo(a,h,i)perylene	26.51	276	12571	5.666	ng/ul	97

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

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