

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080320\
 Data File : BN011351.D
 Acq On : 04 Aug 2020 03:14
 Operator : CG/JU
 Sample : L3419-18
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AG7

Manual Integrations
 APPROVED

mohammad
 8/5/2020 10:28:10 AM

Quant Time: Aug 04 07:22:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 04 06:45:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1065	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4193	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	2611	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	6035	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	6810	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	8622	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	536	0.05	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	1502	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.26	128	302	0.024	ng/ul#	52
5) 2-Methylnaphthalene	11.88	142	187	0.022	ng/ul	99
7) Acenaphthylene	13.81	152	9494	0.761	ng/ul	99
8) Acenaphthene	14.16	153	457	0.046	ng/ul	93
9) Fluorene	15.15	166	593	0.051	ng/ul#	96
12) Phenanthrene	16.89	178	14196	0.710	ng/ul	100
13) Anthracene	16.98	178	8835	0.506	ng/ul	99
15) Fluoranthene	18.92	202	52681	2.137	ng/ul	99
17) Pyrene	19.28	202	55382	1.791	ng/ul	100
18) Benzo(a)anthracene	21.05	228	29739	1.137	ng/ul	94
19) Chrysene	21.10	228	77398	2.580	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	206899m	5.749	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	70873m	1.757	ng/ul	
23) Benzo(a)pyrene	23.09	252	95426m	3.050	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.24	276	241809	7.353	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.24	278	52158	1.992	ng/ul	95
26) Benzo(g,h,i)perylene	25.87	276	434697	14.630	ng/ul	98

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

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