

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008670.D
 Acq On : 20 Nov 2019 20:56
 Operator : JU
 Sample : K5851-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA6

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:06:24 PM

Quant Time: Nov 21 01:08:39 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	3989	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	14594	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8509	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	17274	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13663	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13468	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4370	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	10025	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	26208	0.596	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	20575	0.670	ng/ul	100
7) Acenaphthylene	13.95	152	5880	0.125	ng/ul#	91
8) Acenaphthene	14.30	153	2110	0.060	ng/ul	95
9) Fluorene	15.29	166	4297	0.105	ng/ul#	88
12) Phenanthrene	17.02	178	58892	1.011	ng/ul	98
13) Anthracene	17.11	178	14275	0.259	ng/ul#	95
15) Fluoranthene	19.05	202	112190	1.568	ng/ul	95
17) Pyrene	19.41	202	124490	2.271	ng/ul	98
18) Benzo(a)anthracene	21.17	228	66324	1.207	ng/ul	95
19) Chrysene	21.22	228	99894	1.851	ng/ul	97
21) Benzo(b)fluoranthene	22.71	252	130316m	2.346	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	33367m	0.606	ng/ul	
23) Benzo(a)pyrene	23.26	252	61454	1.179	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.49	276	59474	0.976	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.50	278	18283	0.371	ng/ul#	65
26) Benzo(a,h,i)perylene	26.14	276	63861	1.252	ng/ul	97

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

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