

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008672.D
 Acq On : 20 Nov 2019 22:08
 Operator : JU
 Sample : K5851-11
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA9

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:06:30 PM

Quant Time: Nov 21 01:13:25 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	4001	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	15081	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8308	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	17075	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13975	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13940	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	3859	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	8524	0.16	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.42	128	1133	0.025	ng/ul#	80
7) Acenaphthylene	13.95	152	2601	0.057	ng/ul#	92
12) Phenanthrene	17.02	178	13667	0.237	ng/ul	100
13) Anthracene	17.11	178	3294m	0.061	ng/ul	
15) Fluoranthene	19.04	202	35522	0.502	ng/ul	92
17) Pyrene	19.41	202	32032	0.571	ng/ul	99
18) Benzo(a)anthracene	21.16	228	14742	0.262	ng/ul	95
19) Chrysene	21.22	228	15202	0.275	ng/ul	96
21) Benzo(b)fluoranthene	22.71	252	18896m	0.329	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	6826m	0.120	ng/ul	
23) Benzo(a)pyrene	23.25	252	13038	0.242	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.49	276	9105	0.144	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.49	278	2115	0.042	ng/ul#	1
26) Benzo(g,h,i)perylene	26.14	276	8832	0.167	ng/ul#	88

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

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