

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061819\
 Data File : BN006378.D
 Acq On : 18 Jun 2019 15:13
 Operator : JU/SJ
 Sample : K3335-22
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Y0

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:39:08 AM

Quant Time: Jun 18 17:04:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 18 15:57:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	6587	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	26243	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	14115	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	29306m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	18313m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	16303m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7186	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	17319m	0.22	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	50861	0.702	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	21558	0.427	ng/ul	99
7) Acenaphthylene	14.01	152	11266	0.144	ng/ul#	94
8) Acenaphthene	14.35	153	59320	1.090	ng/ul	99
9) Fluorene	15.35	166	76773	1.206	ng/ul	98
12) Phenanthrene	17.09	178	944202m	10.037	ng/ul	
13) Anthracene	17.18	178	267609m	2.979	ng/ul	
15) Fluoranthene	19.12	202	1025105m	9.868	ng/ul	
17) Pyrene	19.49	202	819259	8.493	ng/ul	98
18) Benzo(a)anthracene	21.25	228	370767m	4.578	ng/ul	
19) Chrysene	21.30	228	312598	4.111	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	326806m	4.807	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	139399m	2.105	ng/ul	
23) Benzo(a)pyrene	23.40	252	238672	3.770	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.73	276	135433	1.976	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.73	278	35456	0.647	ng/ul	96
26) Benzo(a,h,i)perylene	26.41	276	103235	1.799	ng/ul	94

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

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