

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003132.D
 Acq On : 12 Oct 2018 08:20
 Operator : MJ/SJ
 Sample : J5234-17
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK0

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:38:44 PM

Quant Time: Oct 12 12:52:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1840	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6774	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3994	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.01	188	5521m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	4866m	0.40	ng/ul	-0.01
20) Perylene-d12	23.43	264	4799m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	2758	0.26	ng/ul	-0.03
14) Fluoranthene-d10	19.05	212	4577m	0.29	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	25940	1.492	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	45710	3.820	ng/ul	97
7) Acenaphthylene	13.98	152	3758	0.206	ng/ul#	44
8) Acenaphthene	14.32	153	6735	0.486	ng/ul	93
9) Fluorene	15.31	166	29636	1.866	ng/ul#	99
12) Phenanthrene	17.05	178	55394	3.599	ng/ul	98
13) Anthracene	17.15	178	17814m	1.225	ng/ul	
15) Fluoranthene	19.08	202	19207m	1.064	ng/ul	
17) Pyrene	19.44	202	30021	1.588	ng/ul	99
18) Benzo(a)anthracene	21.21	228	9228m	0.589	ng/ul	
19) Chrysene	21.26	228	8895m	0.559	ng/ul	
21) Benzo(b)fluoranthene	22.78	252	4885m	0.283	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	1186m	0.066	ng/ul	
23) Benzo(a)pyrene	23.34	252	5341	0.345	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.65	276	2515	0.141	ng/ul	90
25) Dibenzo(a,h)anthracene	25.65	278	567	0.040	ng/ul#	1
26) Benzo(a,h,i)perylene	26.33	276	2231	0.149	ng/ul#	72

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

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