

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004124.D
 Acq On : 19 Dec 2018 21:24
 Operator : JU/SJ
 Sample : J6297-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F28

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:03:32 PM

Quant Time: Dec 20 06:31:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 20 05:49:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1328	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	6773	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4161	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	11026m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	12719m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	11369m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2223	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	7664m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	12544	0.659	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	9952	0.749	ng/ul	100
12) Phenanthrene	17.25	178	13441m	0.371	ng/ul	
13) Anthracene	17.34	178	3986m	0.134	ng/ul	
15) Fluoranthene	19.27	202	15680m	0.367	ng/ul	
17) Pyrene	19.64	202	23624m	0.518	ng/ul	
18) Benzo(a)anthracene	21.38	228	10600	0.264	ng/ul#	88
19) Chrysene	21.44	228	13377	0.291	ng/ul#	92
21) Benzo(b)fluoranthene	23.02	252	12521m	0.266	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	4168m	0.091	ng/ul	
23) Benzo(a)pyrene	23.62	252	9047	0.222	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.08	276	7295	0.156	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.08	278	2372	0.063	ng/ul#	4
26) Benzo(g,h,i)perylene	26.81	276	17004	0.419	ng/ul#	88

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

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Quant Time: Dec 20 06:31:13 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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