

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004126.D
 Acq On : 19 Dec 2018 22:34
 Operator : JU/SJ
 Sample : J6299-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F20

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:03:38 PM

Quant Time: Dec 20 06:54:44 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	1413	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7160	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4385	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	11744m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13803m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12528m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2119	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	7655m	0.22	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	729	0.037	ng/ul#	77
12) Phenanthrene	17.26	178	1672m	0.043	ng/ul	
13) Anthracene	17.35	178	852m	0.027	ng/ul	
15) Fluoranthene	19.27	202	2866m	0.063	ng/ul	
17) Pyrene	19.64	202	2954	0.060	ng/ul#	88
18) Benzo(a)anthracene	21.38	228	2287	0.052	ng/ul#	63
19) Chrysene	21.44	228	3153	0.063	ng/ul#	70
21) Benzo(b)fluoranthene	23.02	252	3370m	0.065	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	1236m	0.024	ng/ul	
23) Benzo(a)pyrene	23.62	252	1880	0.042	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.08	276	1773	0.034	ng/ul#	99
26) Benzo(g,h,i)perylene	26.82	276	2423	0.054	ng/ul#	44

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

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Quant Time: Dec 20 06:54:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 20 05:49:27 2018
Response via : Initial Calibration

