

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN004018.D
 Acq On : 17 Dec 2018 04:48
 Operator : JU/SJ
 Sample : J6236-06
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E16

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:28:26 PM

Quant Time: Dec 17 07:50:23 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 : Mon Dec 17 06:22:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1949	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9697	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6090	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	15160m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	16944m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	15021m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4136	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	12109m	0.27	ng/ul	0.00
Target Compounds						
7) Acenaphthylene	14.19	152	804	0.029	ng/ul#	10
12) Phenanthrene	17.26	178	2028m	0.041	ng/ul	
15) Fluoranthene	19.27	202	3573m	0.061	ng/ul	
17) Pyrene	19.64	202	4800m	0.079	ng/ul	
18) Benzo(a)anthracene	21.39	228	2582m	0.048	ng/ul	
19) Chrysene	21.43	228	4175m	0.068	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	4084m	0.066	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	1786m	0.030	ng/ul	
23) Benzo(a)pyrene	23.62	252	2434	0.045	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.07	276	2019	0.033	ng/ul#	95
26) Benzo(g,h,i)perylene	26.81	276	2231	0.042	ng/ul#	28

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

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