

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003958.D
 Acq On : 15 Dec 2018 14:31
 Operator : JU/SJ
 Sample : J6227-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F33

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:23:51 PM

Quant Time: Dec 15 22:48:16 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 : Sat Dec 15 21:58:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1778	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8864	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5432	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	13431m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	17371m	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	15893m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2435	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	7619m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.29	142	814	0.047	ng/ul	100
12) Phenanthrene	17.25	178	3750	0.085	ng/ul	95
13) Anthracene	17.35	178	1035m	0.029	ng/ul	
15) Fluoranthene	19.27	202	5492m	0.106	ng/ul	
17) Pyrene	19.63	202	8555m	0.137	ng/ul	
18) Benzo(a)anthracene	21.38	228	4127m	0.075	ng/ul	
19) Chrysene	21.43	228	5699m	0.091	ng/ul	
21) Benzo(b)fluoranthene	23.01	252	5909m	0.090	ng/ul	
22) Benzo(k)fluoranthene	23.05	252	2055m	0.032	ng/ul	
23) Benzo(a)pyrene	23.61	252	3741m	0.066	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.06	276	2662m	0.041	ng/ul	
26) Benzo(g,h,i)perylene	26.79	276	3186m	0.056	ng/ul	

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

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