

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004045.D
 Acq On : 17 Dec 2018 20:46
 Operator : JU/SJ
 Sample : J6236-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E18

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:17 PM

Quant Time: Dec 18 05:51:47 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 : Tue Dec 18 04:29:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1407	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7411	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4691	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12260m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13973m	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	12291m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3271	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	10704m	0.30	ng/ul	0.00
Target Compounds				Qvalue		
5) 2-Methylnaphthalene	12.29	142	596	0.041	ng/ul	99
7) Acenaphthylene	14.19	152	1479	0.070	ng/ul#	87
12) Phenanthrene	17.26	178	14202	0.353	ng/ul	99
13) Anthracene	17.35	178	2078m	0.063	ng/ul	
15) Fluoranthene	19.27	202	14273m	0.300	ng/ul	
17) Pyrene	19.64	202	22415m	0.447	ng/ul	
18) Benzo(a)anthracene	21.39	228	13068m	0.296	ng/ul	
19) Chrysene	21.44	228	32577m	0.645	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	17274m	0.340	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	4726m	0.095	ng/ul	
23) Benzo(a)pyrene	23.62	252	10341	0.235	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.08	276	6833	0.135	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.09	278	3400	0.084	ng/ul#	36
26) Benzo(g,h,i)perylene	26.82	276	8797	0.200	ng/ul#	84

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

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