

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003983.D
 Acq On : 16 Dec 2018 06:30
 Operator : JU/SJ
 Sample : J6171-16
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y26

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:24:15 PM

Quant Time: Dec 17 00:38:49 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 00:13:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	2720	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	12333	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	7976	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	18470m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18107m	0.40	ng/ul	0.01
20) Perylene-d12	23.74	264	16062m	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4587	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	14514m	0.27	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.68	128	17147	0.495	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	33366	1.379	ng/ul	100
7) Acenaphthylene	14.19	152	1974	0.055	ng/ul#	61
8) Acenaphthene	14.54	153	2384	0.076	ng/ul	93
9) Fluorene	15.52	166	4940	0.135	ng/ul#	95
12) Phenanthrene	17.26	178	331547	5.469	ng/ul	99
13) Anthracene	17.35	178	29458m	0.591	ng/ul	
15) Fluoranthene	19.28	202	58336m	0.815	ng/ul	
17) Pyrene	19.64	202	284857	4.386	ng/ul	98
18) Benzo(a)anthracene	21.39	228	87966m	1.538	ng/ul	
19) Chrysene	21.44	228	167487m	2.559	ng/ul	
21) Benzo(b)fluoranthene	23.03	252	40347m	0.607	ng/ul	
23) Benzo(a)pyrene	23.64	252	62936m	1.094	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.12	276	87135m	1.321	ng/ul	
25) Dibenzo(a,h)anthracene	26.11	278	26818m	0.505	ng/ul	
26) Benzo(a,h,i)perylene	26.87	276	478187	8.339	ng/ul	93

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

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