

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003985.D
 Acq On : 16 Dec 2018 07:41
 Operator : JU/SJ
 Sample : J6171-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y05

Manual Integrations
 APPROVED

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

Quant Time: Dec 17 00:45:49 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	2973	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	13478	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	8039	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	19370m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	19592m	0.40	ng/ul	0.01
20) Perylene-d12	23.74	264	15659m	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4253	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	11791m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.30	142	1854	0.070	ng/ul	97
7) Acenaphthylene	14.20	152	1887	0.052	ng/ul#	72
12) Phenanthrene	17.26	178	14161m	0.223	ng/ul	
13) Anthracene	17.36	178	2624m	0.050	ng/ul	
15) Fluoranthene	19.28	202	13537m	0.180	ng/ul	
17) Pyrene	19.64	202	18875m	0.269	ng/ul	
18) Benzo(a)anthracene	21.39	228	9308m	0.150	ng/ul	
19) Chrysene	21.44	228	14446m	0.204	ng/ul	
21) Benzo(b)fluoranthene	23.03	252	25091m	0.387	ng/ul	
22) Benzo(k)fluoranthene	23.07	252	7439m	0.118	ng/ul	
23) Benzo(a)pyrene	23.64	252	10325	0.184	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.11	276	22727	0.353	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.10	278	4636	0.090	ng/ul#	1
26) Benzo(g,h,i)perylene	26.85	276	70831	1.267	ng/ul#	80

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

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