

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003971.D
 Acq On : 15 Dec 2018 22:46
 Operator : JU/SJ
 Sample : J6171-06DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y10DL

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:24:00 PM

Quant Time: Dec 15 23:27:06 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 23:04:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	2058	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	10406	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6224	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	15553	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	20164	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	18268	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	478	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	1469	0.03	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.29	142	539	0.026	ng/ul	97
12) Phenanthrene	17.26	178	2169	0.042	ng/ul#	91
15) Fluoranthene	19.27	202	3509	0.058	ng/ul	91
17) Pyrene	19.63	202	4844	0.067	ng/ul#	91
18) Benzo(a)anthracene	21.38	228	2314	0.036	ng/ul#	68
19) Chrysene	21.43	228	4112	0.056	ng/ul#	79
21) Benzo(b)fluoranthene	23.01	252	4562m	0.060	ng/ul	
22) Benzo(k)fluoranthene	23.05	252	1598m	0.022	ng/ul	
23) Benzo(a)pyrene	23.61	252	2169	0.033	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.07	276	2108	0.028	ng/ul#	97
26) Benzo(g,h,i)perylene	26.79	276	2639	0.040	ng/ul#	49

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

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