

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003986.D
 Acq On : 16 Dec 2018 08:16
 Operator : JU/SJ
 Sample : J6171-01DL 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y05DL

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:24:18 PM

Quant Time: Dec 17 00:48:03 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	3084	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	14082	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	7715	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	18448	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	18693	0.40	ng/ul	0.01
20) Perylene-d12	23.73	264	16346	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	932	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	2375	0.04	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.26	178	3024	0.050	ng/ul#	78
15) Fluoranthene	19.28	202	3189	0.045	ng/ul	82
17) Pyrene	19.64	202	4807	0.072	ng/ul#	86
18) Benzo(a)anthracene	21.39	228	1899	0.032	ng/ul#	1
19) Chrysene	21.44	228	3462m	0.051	ng/ul	
21) Benzo(b)fluoranthene	23.03	252	5427m	0.080	ng/ul	
22) Benzo(k)fluoranthene	23.07	252	1662m	0.025	ng/ul	
23) Benzo(a)pyrene	23.63	252	2084	0.036	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.10	276	4698	0.070	ng/ul#	90
26) Benzo(g,h,i)perylene	26.83	276	15318	0.262	ng/ul#	56

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

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M
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