

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121118\
 Data File : BN003827.D
 Acq On : 11 Dec 2018 22:29
 Operator : JU/SJ
 Sample : J6170-17DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9M12DL

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:50:34 PM

Quant Time: Dec 12 03:41:00 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 : Wed Dec 12 02:52:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4674	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	17696	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	8031	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	14678	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	11140	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	12559	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	1083	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	1570	0.04	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.25	178	1569	0.033	ng/ul#	91
15) Fluoranthene	19.27	202	4153	0.073	ng/ul	95
17) Pyrene	19.63	202	3766	0.094	ng/ul#	94
18) Benzo(a)anthracene	21.38	228	1265	0.036	ng/ul#	87
19) Chrysene	21.43	228	1919	0.048	ng/ul#	91
21) Benzo(b)fluoranthene	23.01	252	3021m	0.058	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	1141m	0.023	ng/ul	
23) Benzo(a)pyrene	23.61	252	1636	0.036	ng/ul#	39
24) Indeno(1,2,3-cd)pyrene	26.07	276	1318	0.026	ng/ul#	100
26) Benzo(g,h,i)perylene	26.79	276	1086	0.024	ng/ul#	63

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

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