

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100918\
 Data File : BN003020.D
 Acq On : 09 Oct 2018 11:20
 Operator : MJ/SJ
 Sample : J5116-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC57

Quant Time: Oct 09 11:55:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 09 08:56:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1357	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	5752	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3740	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	441575	0.40	ng/ul	0.08
16) Chrysene-d12	21.26	240	3886	0.40	ng/ul	0.02
20) Perylene-d12	23.39	264	200699	0.40	ng/ul	-0.07
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3149	0.35	ng/ul	-0.03
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds					Qvalue	
3) Naphthalene	10.44	128	7359	0.499	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	2410	0.237	ng/ul	99
7) Acenaphthylene	13.98	152	2472	0.144	ng/ul#	86
8) Acenaphthene	14.32	153	32393	2.495	ng/ul	99
9) Fluorene	15.32	166	31135	2.094	ng/ul	100
12) Phenanthrene	17.15	178	140225	0.114	ng/ul	99
17) Pyrene	19.46	202	596637	39.531	ng/ul	99
18) Benzo(a)anthracene	21.24	228	209496	16.753	ng/ul#	94
19) Chrysene	21.29	228	240837	18.966	ng/ul#	94
21) Benzo(b)fluoranthene	22.85	252	480218	0.664	ng/ul#	1
22) Benzo(k)fluoranthene	22.85	252	473486	0.632	ng/ul#	1
23) Benzo(a)pyrene	23.34	252	227422	0.351	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.57	276	40729	0.055	ng/ul#	72
25) Dibenzo(a,h)anthracene	25.51	278	25110	0.042	ng/ul#	1
26) Benzo(g,h,i)perylene	26.15	276	19982	0.032	ng/ul#	1

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

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