

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003012.D
 Acq On : 09 Oct 2018 06:05
 Operator : MJ/SJ
 Sample : J5116-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC56

Manual Integrations
 APPROVED

Sohil
 10/9/2018 5:38:25 PM

Quant Time: Oct 09 09:04:47 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	1134	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	4969	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3072	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7331m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7882m	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	7839	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2680	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	7982m	0.38	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	4705	0.369	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	291	0.033	ng/ul	100
7) Acenaphthylene	13.98	152	695	0.049	ng/ul#	79
8) Acenaphthene	14.32	153	656	0.062	ng/ul	95
9) Fluorene	15.32	166	944	0.077	ng/ul#	98
12) Phenanthrene	17.06	178	25813	1.263	ng/ul	99
13) Anthracene	17.15	178	2665m	0.138	ng/ul	
15) Fluoranthene	19.09	202	57056m	2.380	ng/ul	
17) Pyrene	19.45	202	49808	1.627	ng/ul	96
18) Benzo(a)anthracene	21.21	228	25039	0.987	ng/ul	98
19) Chrysene	21.26	228	33178	1.288	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	35579m	1.260	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	12169m	0.416	ng/ul	
23) Benzo(a)pyrene	23.33	252	17747	0.701	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.65	276	10865	0.374	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.65	278	2812	0.121	ng/ul#	77
26) Benzo(g,h,i)perylene	26.33	276	9140	0.373	ng/ul	99

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

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