

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100918\
 Data File : BN003021.D
 Acq On : 09 Oct 2018 11:55
 Operator : MJ/SJ
 Sample : J5116-09DL 5X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC58DL

Quant Time: Oct 09 12:38:55 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	7.64	152	1202	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	5153	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.27	164	3067	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	85494	0.40	ng/ul	0.09
16) Chrysene-d12	21.32	240	52	0.40	ng/ul	0.08
20) Perylene-d12	23.53	264	680	0.40	ng/ul	0.07
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	508	0.06	ng/ul	0.01
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	423	0.032	ng/ul#	57
8) Acenaphthene	14.33	153	583	0.055	ng/ul	86
9) Fluorene	15.33	166	335	0.027	ng/ul#	28
17) Pyrene	19.45	202	15224	75.379	ng/ul#	94
18) Benzo(a)anthracene	21.27	228	9327	55.738	ng/ul#	75
19) Chrysene	21.38	228	864	5.085	ng/ul#	1
21) Benzo(b)fluoranthene	22.78	252	16719	6.827	ng/ul#	1
22) Benzo(k)fluoranthene	22.96	252	1397	0.550	ng/ul#	1
23) Benzo(a)pyrene	23.49	252	2404	1.095	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.66	276	5163	2.047	ng/ul#	49
25) Dibenzo(a,h)anthracene	25.75	278	341	0.169	ng/ul#	1
26) Benzo(g,h,i)perylene	26.34	276	4835	2.273	ng/ul#	1

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

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