

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003123.D
 Acq On : 12 Oct 2018 03:04
 Operator : MJ/SJ
 Sample : J5230-14
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 12 07:36:05 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 : Fri Oct 12 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1513	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	5465	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	2704	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5210m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	4741m	0.40	ng/ul	-0.01
20) Perylene-d12	23.43	264	4855	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	2619	0.31	ng/ul	-0.02
14) Fluoranthene-d10	19.05	212	5072m	0.34	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	17404	1.241	ng/ul	97
5) 2-Methylnaphthalene	12.07	142	10650	1.103	ng/ul	97
7) Acenaphthylene	13.98	152	22537	1.822	ng/ul	99
8) Acenaphthene	14.32	153	3483	0.371	ng/ul	98
9) Fluorene	15.32	166	4347	0.404	ng/ul#	98
12) Phenanthrene	17.05	178	94817m	6.529	ng/ul	
13) Anthracene	17.15	178	28307m	2.063	ng/ul	
15) Fluoranthene	19.08	202	183309m	10.759	ng/ul	
17) Pyrene	19.45	202	170394	9.254	ng/ul	96
18) Benzo(a)anthracene	21.21	228	86338m	5.659	ng/ul	
19) Chrysene	21.26	228	89364	5.768	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	132667m	7.588	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	47512m	2.621	ng/ul	
23) Benzo(a)pyrene	23.34	252	80435	5.133	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.64	276	62674	3.481	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.64	278	16746	1.162	ng/ul	99
26) Benzo(a,h,i)perylene	26.32	276	56549	3.724	ng/ul	93

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

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