

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003066.D
 Acq On : 10 Oct 2018 16:15
 Operator : MJ/SJ
 Sample : J5230-17
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YX3

Quant Time: Oct 11 07:07:33 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 : Thu Oct 11 07:04:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1191	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.41	136	4851	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	2923	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	7203	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7022	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	5738	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	2735	0.36	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	8613	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	4053	0.326	ng/ul	96
5) 2-Methylnaphthalene	12.08	142	5901	0.689	ng/ul	97
7) Acenaphthylene	13.99	152	660	0.049	ng/ul#	75
8) Acenaphthene	14.32	153	4493	0.443	ng/ul	99
9) Fluorene	15.33	166	7515	0.647	ng/ul	99
11) Pentachlorophenol	16.72	266	868	1.242	ng/ul	99
12) Phenanthrene	17.07	178	9667	0.481	ng/ul	100
13) Anthracene	17.17	178	7078	0.373	ng/ul	99
15) Fluoranthene	19.09	202	11874	0.504	ng/ul	96
17) Pyrene	19.46	202	11995	0.440	ng/ul	99
18) Benzo(a)anthracene	21.22	228	12439	0.550	ng/ul	99
19) Chrysene	21.27	228	13173	0.574	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	15720	0.761	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	8167	0.381	ng/ul	98
23) Benzo(a)pyrene	23.36	252	4289	0.232	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.67	276	3434	0.161	ng/ul#	47
25) Dibenzo(a,h)anthracene	25.67	278	13754	0.807	ng/ul#	91
26) Benzo(g,h,i)perylene	26.35	276	5791	0.323	ng/ul	99

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

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