

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111918\
 Data File : BN003590.D
 Acq On : 19 Nov 2018 12:50
 Operator : MJ/SJ
 Sample : SSTD0.262
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.262

Quant Time: Nov 19 14:03:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 19 13:56:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	5624	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	20741	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11219	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	28642	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	29208	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	26617	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	6244	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	15804	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	11529	0.198	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	7984	0.191	ng/ul	98
7) Acenaphthylene	14.20	152	9774	0.200	ng/ul	99
8) Acenaphthene	14.54	153	8687	0.199	ng/ul	100
9) Fluorene	15.53	166	10086	0.197	ng/ul	99
11) Pentachlorophenol	16.87	266	733	0.128	ng/ul	98
12) Phenanthrene	17.26	178	18216	0.194	ng/ul	99
13) Anthracene	17.36	178	14482	0.189	ng/ul	99
15) Fluoranthene	19.28	202	20416	0.193	ng/ul	99
17) Pyrene	19.64	202	20355	0.192	ng/ul	98
18) Benzo(a)anthracene	21.38	228	17586	0.192	ng/ul	99
19) Chrysene	21.43	228	20895	0.201	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	21500	0.187	ng/ul	97
22) Benzo(k)fluoranthene	23.05	252	20246	0.182	ng/ul	97
23) Benzo(a)pyrene	23.61	252	18069	0.189	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.06	276	20306	0.198	ng/ul	99
25) Dibenzo(a,h)anthracene	26.08	278	16259	0.195	ng/ul	98
26) Benzo(g,h,i)perylene	26.79	276	17909	0.204	ng/ul	100

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

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