

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021119\
 Data File : BN004371.D
 Acq On : 11 Feb 2019 10:27
 Operator : JU/SJ
 Sample : SSTD0.203
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.203

Quant Time: Feb 11 12:10:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 11:59:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	1405	0.40	ng/ul	0.00
2) Naphthalene-d8	10.54	136	5308	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	2980	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	8084	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	9667	0.40	ng/ul	0.00
20) Perylene-d12	23.61	264	10014	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	1625	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.17	212	4928	0.21	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.59	128	3075	0.206	ng/ul	98
5) 2-Methylnaphthalene	12.20	142	2148	0.207	ng/ul	99
7) Acenaphthylene	14.11	152	2713	0.182	ng/ul	98
8) Acenaphthene	14.45	153	2380	0.199	ng/ul	99
9) Fluorene	15.44	166	2800	0.206	ng/ul	99
11) Pentachlorophenol	16.78	266	283	0.242	ng/ul	95
12) Phenanthrene	17.18	178	5282	0.201	ng/ul	99
13) Anthracene	17.27	178	4330	0.181	ng/ul	99
15) Fluoranthene	19.20	202	6274	0.198	ng/ul	99
17) Pyrene	19.56	202	6434	0.195	ng/ul	98
18) Benzo(a)anthracene	21.32	228	6184	0.188	ng/ul	99
19) Chrysene	21.37	228	7383	0.210	ng/ul	99
21) Benzo(b)fluoranthene	22.92	252	7986	0.210	ng/ul	95
22) Benzo(k)fluoranthene	22.96	252	7715	0.191	ng/ul	96
23) Benzo(a)pyrene	23.50	252	7057	0.195	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.90	276	8859	0.237	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	7064	0.230	ng/ul	96
26) Benzo(g,h,i)perylene	26.60	276	8209	0.261	ng/ul	98

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

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