

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008078.D
 Acq On : 11 Oct 2019 15:31
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Oct 11 19:31:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2031	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	9931	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	6552	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	15696	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	15718	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	17034	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	6071	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	19397	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	10295	0.364	ng/ul	100
5) 2-Methylnaphthalene	12.07	142	6952	0.358	ng/ul	100
7) Acenaphthylene	13.98	152	11027	0.353	ng/ul	99
8) Acenaphthene	14.33	153	9700	0.400	ng/ul	99
9) Fluorene	15.32	166	9954	0.361	ng/ul	98
11) Pentachlorophenol	16.68	266	1327	0.350	ng/ul	98
12) Phenanthrene	17.06	178	16176	0.352	ng/ul	99
13) Anthracene	17.15	178	13891	0.343	ng/ul	100
15) Fluoranthene	19.08	202	21029	0.346	ng/ul	100
17) Pyrene	19.44	202	24886	0.398	ng/ul	99
18) Benzo(a)anthracene	21.20	228	19875	0.348	ng/ul	99
19) Chrysene	21.25	228	25576	0.365	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	22036	0.397	ng/ul	99
22) Benzo(k)fluoranthene	22.79	252	24831	0.395	ng/ul	100
23) Benzo(a)pyrene	23.31	252	22670	0.389	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	26869	0.391	ng/ul	100
25) Dibenzo(a,h)anthracene	25.59	278	21114	0.389	ng/ul	99
26) Benzo(g,h,i)perylene	26.24	276	24718	0.392	ng/ul	98

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

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