

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013029.D
 Acq On : 10 Dec 2020 02:09
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 99 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.436

Quant Time: Dec 10 03:26:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 09 07:33:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1070	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.14	136	4504	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.05	164	2474	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4922	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3398	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3201	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	2871	0.43	ng/ul	-0.01
14) Fluoranthene-d10	18.85	212	5346	0.40	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.19	128	5395	0.414	ng/ul	93
5) 2-Methylnaphthalene	11.82	142	3648	0.426	ng/ul	99
7) Acenaphthylene	13.76	152	5321	0.454	ng/ul	95
8) Acenaphthene	14.11	153	4045	0.421	ng/ul	99
9) Fluorene	15.11	166	4444	0.403	ng/ul	97
11) Pentachlorophenol	16.47	266	558	0.391	ng/ul	94
12) Phenanthrene	16.85	178	7015	0.425	ng/ul	98
13) Anthracene	16.94	178	6410	0.456	ng/ul	96
15) Fluoranthene	18.88	202	7796	0.398	ng/ul	99
17) Pyrene	19.25	202	7876	0.491	ng/ul	99
18) Benzo(a)anthracene	21.01	228	6248	0.444	ng/ul	98
19) Chrysene	21.06	228	6803	0.438	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	6288	0.425	ng/ul	95
22) Benzo(k)fluoranthene	22.55	252	6667	0.407	ng/ul	98
23) Benzo(a)pyrene	23.04	252	5830	0.424	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.18	276	6795	0.376	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.19	278	5450	0.382	ng/ul	96
26) Benzo(g,h,i)perylene	25.81	276	5744	0.352	ng/ul	97

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

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