

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012874.D
 Acq On : 04 Dec 2020 13:05
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Dec 04 13:40:02 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 : Fri Dec 04 10:02:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	605	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	2365	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	1376	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	2870	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	2312	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	2409	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.77	152	1405	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	3039	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.21	128	2687	0.393	ng/ul	95
5) 2-Methylnaphthalene	11.85	142	1754	0.390	ng/ul	99
7) Acenaphthylene	13.78	152	2559	0.393	ng/ul	98
8) Acenaphthene	14.12	153	2099	0.393	ng/ul	98
9) Fluorene	15.12	166	2365	0.386	ng/ul	97
11) Pentachlorophenol	16.48	266	322	0.387	ng/ul	99
12) Phenanthrene	16.86	178	3821	0.397	ng/ul	98
13) Anthracene	16.96	178	3363	0.411	ng/ul	98
15) Fluoranthene	18.89	202	4404	0.385	ng/ul	100
17) Pyrene	19.26	202	4463	0.409	ng/ul	98
18) Benzo(a)anthracene	21.02	228	3696	0.386	ng/ul	99
19) Chrysene	21.08	228	4322	0.409	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	4323	0.388	ng/ul	97
22) Benzo(k)fluoranthene	22.56	252	4751	0.386	ng/ul	99
23) Benzo(a)pyrene	23.06	252	4017	0.388	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.21	276	5248	0.386	ng/ul#	56
25) Dibenzo(a,h)anthracene	25.22	278	4178	0.389	ng/ul	97
26) Benzo(g,h,i)perylene	25.83	276	4681	0.381	ng/ul	99

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

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