

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012866.D
 Acq On : 04 Dec 2020 02:45
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Dec 04 03:16:57 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 : Thu Dec 03 15:27:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	640	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	2507	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	1427	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	3003	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	2480	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	2539	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.78	152	1454	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	3161	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	2823	0.389	ng/ul	96
5) 2-Methylnaphthalene	11.85	142	1841	0.386	ng/ul	100
7) Acenaphthylene	13.78	152	2658	0.394	ng/ul	98
8) Acenaphthene	14.13	153	2182	0.394	ng/ul	99
9) Fluorene	15.12	166	2453	0.386	ng/ul	100
11) Pentachlorophenol	16.48	266	333	0.383	ng/ul	91
12) Phenanthrene	16.86	178	3963	0.394	ng/ul	100
13) Anthracene	16.96	178	3382	0.395	ng/ul	99
15) Fluoranthene	18.89	202	4569	0.382	ng/ul	98
17) Pyrene	19.26	202	4627	0.395	ng/ul	99
18) Benzo(a)anthracene	21.02	228	3930	0.383	ng/ul	100
19) Chrysene	21.08	228	4615	0.407	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	4622	0.393	ng/ul	97
22) Benzo(k)fluoranthene	22.56	252	5073	0.391	ng/ul	98
23) Benzo(a)pyrene	23.05	252	4414	0.405	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.20	276	5591	0.391	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.22	278	4465	0.394	ng/ul	96
26) Benzo(g,h,i)perylene	25.83	276	4938	0.381	ng/ul	98

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

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