

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012248.D
 Acq On : 16 Oct 2020 01:14
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Oct 16 01:54:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 16 01:19:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1563	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6672	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3818	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	7956	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	6753	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	7003	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	4011	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	8248	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	7526	0.391	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	4955	0.405	ng/ul	99
7) Acenaphthylene	13.98	152	6870	0.415	ng/ul	99
8) Acenaphthene	14.32	153	5513	0.394	ng/ul	98
9) Fluorene	15.31	166	6165	0.396	ng/ul	100
11) Pentachlorophenol	16.66	266	962	0.418	ng/ul	99
12) Phenanthrene	17.05	178	9885	0.390	ng/ul	98
13) Anthracene	17.14	178	8866	0.408	ng/ul	99
15) Fluoranthene	19.07	202	11337	0.382	ng/ul	98
17) Pyrene	19.44	202	11668	0.383	ng/ul	99
18) Benzo(a)anthracene	21.19	228	10028	0.401	ng/ul	98
19) Chrysene	21.24	228	11224	0.368	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	11223	0.397	ng/ul	94
22) Benzo(k)fluoranthene	22.78	252	11452	0.366	ng/ul	95
23) Benzo(a)pyrene	23.30	252	10503	0.382	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.57	276	13335	0.377	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	10872	0.384	ng/ul	95
26) Benzo(g,h,i)perylene	26.24	276	11650	0.360	ng/ul	99

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

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