

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101920\
 Data File : BN012373.D
 Acq On : 20 Oct 2020 04:49
 Operator : CG/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.426

Quant Time: Oct 20 08:26:05 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 : Tue Oct 20 08:18:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1499	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6488	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	3584	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	7090	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	6046	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	6320	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3904	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	7172	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	7315	0.391	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	4788	0.403	ng/ul	99
7) Acenaphthylene	13.95	152	6353	0.409	ng/ul	100
8) Acenaphthene	14.30	153	5209	0.396	ng/ul	99
9) Fluorene	15.29	166	5768	0.394	ng/ul	97
11) Pentachlorophenol	16.65	266	804	0.392	ng/ul	99
12) Phenanthrene	17.03	178	8750	0.387	ng/ul	99
13) Anthracene	17.12	178	7863	0.406	ng/ul	100
15) Fluoranthene	19.05	202	9815	0.371	ng/ul	99
17) Pyrene	19.41	202	10211	0.375	ng/ul	97
18) Benzo(a)anthracene	21.17	228	8886	0.397	ng/ul	98
19) Chrysene	21.22	228	10104	0.370	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	10028	0.393	ng/ul	96
22) Benzo(k)fluoranthene	22.75	252	10666	0.377	ng/ul	99
23) Benzo(a)pyrene	23.26	252	9710	0.391	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.52	276	11974	0.375	ng/ul	98
25) Dibenzo(a,h)anthracene	25.54	278	9712	0.381	ng/ul	98
26) Benzo(g,h,i)perylene	26.18	276	10503	0.359	ng/ul	99

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

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