

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012256.D
 Acq On : 16 Oct 2020 06:39
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Oct 16 07:24:30 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 06:21:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1506	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6413	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3698	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	7722	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	6498	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	6636	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3924	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	7908	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	7276	0.393	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	4793	0.408	ng/ul	100
7) Acenaphthylene	13.98	152	6591	0.411	ng/ul	98
8) Acenaphthene	14.32	153	5344	0.394	ng/ul	99
9) Fluorene	15.31	166	5987	0.397	ng/ul	100
11) Pentachlorophenol	16.66	266	873	0.391	ng/ul	98
12) Phenanthrene	17.05	178	9541	0.388	ng/ul	99
13) Anthracene	17.14	178	8500	0.403	ng/ul	98
15) Fluoranthene	19.07	202	10847	0.377	ng/ul	99
17) Pyrene	19.44	202	11184	0.382	ng/ul	100
18) Benzo(a)anthracene	21.19	228	9490	0.394	ng/ul	99
19) Chrysene	21.24	228	10961	0.374	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	10689	0.399	ng/ul	95
22) Benzo(k)fluoranthene	22.78	252	11279	0.380	ng/ul	95
23) Benzo(a)pyrene	23.30	252	10136	0.389	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.57	276	12930	0.386	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	10502	0.392	ng/ul	95
26) Benzo(g,h,i)perylene	26.24	276	11410	0.372	ng/ul	99

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

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