

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012221.D
 Acq On : 15 Oct 2020 08:59
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Oct 15 10:52:14 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 : Thu Oct 15 10:47:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1767	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7563	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4361	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8571	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	7014	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	7398	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	4566	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	8877	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	8554	0.392	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	5642	0.407	ng/ul	99
7) Acenaphthylene	13.98	152	7895	0.417	ng/ul	98
8) Acenaphthene	14.32	153	6247	0.391	ng/ul	99
9) Fluorene	15.32	166	6774	0.381	ng/ul	99
11) Pentachlorophenol	16.67	266	1110	0.448	ng/ul	99
12) Phenanthrene	17.05	178	10560	0.387	ng/ul	99
13) Anthracene	17.14	178	9590	0.409	ng/ul	99
15) Fluoranthene	19.07	202	12118	0.379	ng/ul	98
17) Pyrene	19.44	202	12716	0.402	ng/ul	98
18) Benzo(a)anthracene	21.20	228	10704	0.412	ng/ul	98
19) Chrysene	21.25	228	11623	0.367	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	11557	0.387	ng/ul	95
22) Benzo(k)fluoranthene	22.79	252	11835	0.358	ng/ul#	96
23) Benzo(a)pyrene	23.30	252	10843	0.373	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.58	276	13774	0.369	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	11059	0.370	ng/ul	97
26) Benzo(g,h,i)perylene	26.25	276	12071	0.353	ng/ul	96

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

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