

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009199.D
 Acq On : 27 Dec 2019 17:52
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.423

Quant Time: Dec 27 23:07:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 06:49:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2201	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.27	136	7740	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.14	164	4207	0.40	ng/ul	-0.02
10) Phenanthrene-d10	16.89	188	8832	0.40	ng/ul	-0.02
16) Chrysene-d12	21.09	240	7625	0.40	ng/ul	-0.02
20) Perylene-d12	23.21	264	7368	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	5136	0.37	ng/ul	-0.02
14) Fluoranthene-d10	18.93	212	10329	0.40	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.32	128	9560	0.407	ng/ul	99
5) 2-Methylnaphthalene	11.94	142	6487	0.366	ng/ul	100
7) Acenaphthylene	13.86	152	8848	0.400	ng/ul	99
8) Acenaphthene	14.20	153	6630	0.382	ng/ul	98
9) Fluorene	15.20	166	7461	0.364	ng/ul	99
11) Pentachlorophenol	16.56	266	1152	0.489	ng/ul	100
12) Phenanthrene	16.93	178	12201	0.397	ng/ul	100
13) Anthracene	17.02	178	11373	0.413	ng/ul	99
15) Fluoranthene	18.96	202	13852	0.394	ng/ul	100
17) Pyrene	19.32	202	13996	0.360	ng/ul	99
18) Benzo(a)anthracene	21.07	228	13465	0.398	ng/ul	99
19) Chrysene	21.12	228	13388	0.393	ng/ul	99
21) Benzo(b)fluoranthene	22.59	252	13334	0.388	ng/ul	98
22) Benzo(k)fluoranthene	22.63	252	14011	0.404	ng/ul	97
23) Benzo(a)pyrene	23.13	252	12343	0.409	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.31	276	14442	0.440	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.32	278	11679	0.447	ng/ul	98
26) Benzo(g,h,i)perylene	25.94	276	12383	0.448	ng/ul	99

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

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