

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121619\
 Data File : BN009006.D
 Acq On : 16 Dec 2019 19:42
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.499

Quant Time: Dec 16 20:13:17 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 : Mon Dec 16 17:48:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	2200	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	9887	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6552	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	14496	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	11551	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	9602	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	7187	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	17986	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.35	128	12246	0.408	ng/ul	99
5) 2-Methylnaphthalene	11.97	142	9179	0.405	ng/ul	100
7) Acenaphthylene	13.89	152	13877	0.403	ng/ul	100
8) Acenaphthene	14.24	153	10482	0.388	ng/ul	100
9) Fluorene	15.23	166	12301	0.385	ng/ul	100
11) Pentachlorophenol	16.59	266	1669	0.432	ng/ul	99
12) Phenanthrene	16.96	178	20517	0.407	ng/ul	99
13) Anthracene	17.06	178	19048	0.421	ng/ul	99
15) Fluoranthene	18.99	202	23869	0.414	ng/ul	97
17) Pyrene	19.35	202	24168	0.410	ng/ul	99
18) Benzo(a)anthracene	21.11	228	22121	0.431	ng/ul	99
19) Chrysene	21.16	228	21426	0.416	ng/ul	99
21) Benzo(b)fluoranthene	22.64	252	18355	0.409	ng/ul	97
22) Benzo(k)fluoranthene	22.68	252	18431	0.408	ng/ul	97
23) Benzo(a)pyrene	23.18	252	16450	0.419	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.37	276	16461	0.385	ng/ul	100
25) Dibenzo(a,h)anthracene	25.39	278	13379	0.393	ng/ul	98
26) Benzo(g,h,i)perylene	26.01	276	13800	0.383	ng/ul	99

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

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