

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008942.D
 Acq On : 12 Dec 2019 13:06
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Dec 12 13:45:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 12 00:49:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2520	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10642	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	7062	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	15698	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	14039	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	13170	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	7099	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	18528	0.39	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.36	128	12444	0.388	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	9228	0.412	ng/ul	99
7) Acenaphthylene	13.90	152	13550	0.347	ng/ul	100
8) Acenaphthene	14.25	153	10526	0.363	ng/ul	99
9) Fluorene	15.24	166	12482	0.367	ng/ul	98
11) Pentachlorophenol	16.59	266	1657	0.279	ng/ul	99
12) Phenanthrene	16.97	178	21469	0.406	ng/ul	99
13) Anthracene	17.06	178	19310	0.386	ng/ul	99
15) Fluoranthene	19.00	202	24968	0.384	ng/ul	99
17) Pyrene	19.36	202	24988	0.444	ng/ul	98
18) Benzo(a)anthracene	21.12	228	24336	0.431	ng/ul	100
19) Chrysene	21.17	228	23961	0.432	ng/ul	99
21) Benzo(b)fluoranthene	22.64	252	22948	0.423	ng/ul	99
22) Benzo(k)fluoranthene	22.69	252	23102	0.429	ng/ul	97
23) Benzo(a)pyrene	23.18	252	20527	0.403	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.38	276	21021	0.353	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.40	278	16935	0.352	ng/ul	99
26) Benzo(g,h,i)perylene	26.02	276	17137	0.344	ng/ul	97

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

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