

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN010220\
 Data File : BN009297.D
 Acq On : 02 Jan 2020 16:31
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTD0.410

Quant Time: Jan 02 17:03:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 02 12:44:14 2020
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	66	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	55	0.00
3	Naphthalene	1.213	1.223	-0.8	54	0.00
4 SURR	2-Methylnaphthalene-d10	0.711	0.626	12.0	47	0.00
5	2-Methylnaphthalene	0.917	0.803	12.4	47	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	40	0.00
7	Acenaphthylene	2.101	2.022	3.8	39	0.00
8 C	Acenaphthene	1.649	1.580	4.2	39	0.00
9	Fluorene	1.949	1.702	12.7	36	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	33	0.00
11	Pentachlorophenol	0.107	0.105	1.9	37	0.00
12	Phenanthrene	1.391	1.395	-0.3	32	0.00
13	Anthracene	1.248	1.231	1.4	32	0.00
14 SURR	Fluoranthene-d10	1.178	1.149	2.5	32	0.00
15 C	Fluoranthene	1.591	1.520	4.5	31	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	34	0.00
17	Pyrene	2.041	1.992	2.4	32	0.00
18	Benzo(a)anthracene	1.776	1.770	0.3	33	0.00
19	Chrysene	1.785	1.905	-6.7	35	0.00
20 I	Perylene-d12	1.000	1.000	0.0	41	0.00
21	Benzo(b)fluoranthene	1.868	1.813	2.9	39	0.00
22	Benzo(k)fluoranthene	1.883	1.892	-0.5	41	0.00
23 C	Benzo(a)pyrene	1.637	1.701	-3.9	42	0.00
24	Indeno(1,2,3-cd)pyrene	1.783	1.975	-10.8	45	0.00
25	Dibenzo(a,h)anthracene	1.419	1.590	-12.1	45	0.00
26	Benzo(a,h,i)perylene	1.502	1.724	-14.8	45	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0