

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006986.D
 Acq On : 07 Aug 2019 14:18
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV26

Quant Time: Aug 07 17:40:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	103	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
3	Naphthalene	1.151	2.261	-96.4#	202#	0.00
4 SURR	2-Methylnaphthalene-d10	0.711	1.398	-96.6#	201#	0.00
5	2-Methylnaphthalene	0.864	1.693	-95.9#	202#	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
7	Acenaphthylene	2.318	4.480	-93.3#	200#	0.00
8 C	Acenaphthene	1.586	3.097	-95.3#	200#	0.00
9	Fluorene	1.922	3.788	-97.1#	201#	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
11	Pentachlorophenol	0.110	0.194	-76.4#	167#	0.00
12	Phenanthrene	1.275	2.473	-94.0#	202#	0.00
13	Anthracene	1.304	2.511	-92.6#	200#	0.00
14 SURR	Fluoranthene-d10	1.289	2.536	-96.7#	202#	0.00
15 C	Fluoranthene	1.585	3.114	-96.5#	202#	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	107	0.00
17	Pyrene	1.786	3.409	-90.9#	202#	0.00
18	Benzo(a)anthracene	1.757	3.435	-95.5#	207#	0.00
19	Chrysene	1.645	3.236	-96.7#	206#	0.00
20 I	Perylene-d12	1.000	1.000	0.0	112	0.00
21	Benzo(b)fluoranthene	1.622	3.118	-92.2#	208#	0.00
22	Benzo(k)fluoranthene	1.572	3.025	-92.4#	213#	0.00
23 C	Benzo(a)pyrene	1.527	2.935	-92.2#	212#	0.00
24	Indeno(1,2,3-cd)pyrene	1.646	3.212	-95.1#	221#	0.00
25	Dibenzo(a,h)anthracene	1.315	2.576	-95.9#	222#	0.00
26	Benzo(g,h,i)perylene	1.381	2.697	-95.3#	221#	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 3