

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE101118\
 Data File : BE097856.D
 Acq On : 11 Oct 2018 14:20
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV50

Quant Time: Oct 11 14:53:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 14:16:20 2018
 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	101	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
3	Naphthalene	0.973	0.958	1.5	101	0.00
4 SURR	2-Methylnaphthalene-d10	0.679	0.669	1.5	100	0.00
5	2-Methylnaphthalene	0.704	0.702	0.3	103	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
7	Acenaphthylene	1.513	1.489	1.6	99	0.00
8 C	Acenaphthene	1.174	1.179	-0.4	101	0.00
9	Fluorene	1.532	1.515	1.1	99	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
11	Pentachlorophenol	0.056	0.047	16.1	100	-0.01
12	Phenanthrene	0.947	0.944	0.3	99	0.00
13	Anthracene	0.877	0.870	0.8	100	0.00
14 SURR	Fluoranthene-d10	1.292	1.255	2.9	98	-0.01
15 C	Fluoranthene	1.216	1.192	2.0	99	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	118	-0.01
17	Pyrene	1.192	1.062	10.9	98	-0.01
18	Benzo(a)anthracene	1.209	1.126	6.9	107	-0.01
19	Chrysene	1.087	0.950	12.6	97	-0.01
20 I	Perylene-d12	1.000	1.000	0.0	96	-0.02
21	Benzo(b)fluoranthene	1.069	1.054	1.4	98	-0.02
22	Benzo(k)fluoranthene	1.144	1.122	1.9	96	-0.02
23 C	Benzo(a)pyrene	1.050	1.016	3.2	96	-0.03
24	Indeno(1,2,3-cd)pyrene	1.192	1.153	3.3	96	-0.04
25	Dibenzo(a,h)anthracene	0.985	0.951	3.5	96	-0.03
26	Benzo(g,h,i)perylene	1.038	1.010	2.7	96	-0.04

(#) = Out of Range

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