

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE102518\
 Data File : BE098056.D
 Acq On : 25 Oct 2018 15:28
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SICV67

Quant Time: Oct 25 16:41:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 25 14:48:33 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	95	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
3	Naphthalene	0.957	0.933	2.5	95	0.00
4 SURR	2-Methylnaphthalene-d10	0.713	0.707	0.8	98	0.00
5	2-Methylnaphthalene	0.724	0.710	1.9	96	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
7	Acenaphthylene	1.578	1.531	3.0	95	0.00
8 C	Acenaphthene	1.179	1.122	4.8	93	0.00
9	Fluorene	1.419	1.388	2.2	96	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
11	Pentachlorophenol	0.021	0.019	9.5	91	0.00
12	Phenanthrene	0.923	0.906	1.8	97	0.00
13	Anthracene	0.912	0.876	3.9	95	0.00
14 SURR	Fluoranthene-d10	1.300	1.270	2.3	96	0.00
15 C	Fluoranthene	1.177	1.142	3.0	96	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
17	Pyrene	0.934	0.869	7.0	85	0.00
18	Benzo(a)anthracene	1.176	1.238	-5.3	95	0.00
19	Chrysene	0.910	0.943	-3.6	97	0.00
20 I	Perylene-d12	1.000	1.000	0.0	98	0.00
21	Benzo(b)fluoranthene	1.108	1.101	0.6	97	0.00
22	Benzo(k)fluoranthene	1.085	1.071	1.3	98	0.00
23 C	Benzo(a)pyrene	1.066	1.045	2.0	98	0.00
24	Indeno(1,2,3-cd)pyrene	1.221	1.221	0.0	100	-0.01
25	Dibenzo(a,h)anthracene	1.018	1.014	0.4	99	0.00
26	Benzo(g,h,i)perylene	1.028	1.018	1.0	97	-0.01

(#) = Out of Range

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