

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
 Data File : BE089344.D
 Acq On : 8 Jan 2015 00:19
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE010715

Quant Time: Jan 08 06:11:01 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BASE-BE010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 06:03:23 2015
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	88	0.00
2 I	Naphthalene-d8	5.000	5.000	0.0	89	0.00
3 S	Nitrobenzene-d5	1.000	1.051	-5.1	90	0.00
4	Naphthalene	1.000	1.053	-5.3	91	0.00
5	2-Methylnaphthalene	1.000	1.049	-4.9	91	0.00
6 I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
7 S	2-Fluorobiphenyl	1.000	1.039	-3.9	91	0.00
8	Acenaphthylene	1.000	1.044	-4.4	91	0.00
9 C	Acenaphthene	1.000	1.018	-1.8	91	0.00
10	Fluorene	1.000	1.047	-4.7	93	0.00
11 I	Phenanthrene-d10	5.000	5.000	0.0	88	0.00
12	Phenanthrene	1.000	1.046	-4.6	91	0.00
13	Anthracene	1.000	1.056	-5.6	90	0.00
14 C	Fluoranthene	1.000	1.051	-5.1	90	0.00
15 I	Chrysene-d12	5.000	5.000	0.0	88	0.00
16	Pyrene	1.000	1.032	-3.2	90	0.00
17 S	Terphenyl-d14	1.000	1.008	-0.8	89	0.00
18	Benzo(a)anthracene	1.000	1.025	-2.5	90	0.00
19	Chrysene	1.000	1.055	-5.5	89	0.00
20	Indeno(1,2,3-cd)pyrene	1.000	1.008	-0.8	81	0.00
21 I	Perylene-d12	5.000	5.000	0.0	84	0.00
22	Benzo(b)fluoranthene	1.000	1.055	-5.5	86	0.00
23	Benzo(k)fluoranthene	1.000	1.095	-9.5	88	0.00
24 C	Benzo(a)pyrene	1.000	1.034	-3.4	85	0.00
25	Dibenzo(a,h)anthracene	1.000	1.039	-3.9	83	0.00
26	Benzo(g,h,i)perylene	1.000	1.039	-3.9	84	0.00

(#) = Out of Range

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