

Data Path : Z:\HPCHEM1\BNA_E\Data\Be102414\
 Data File : BE088225.D
 Acq On : 24 Oct 2014 22:17
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Oct 27 13:39:47 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102414.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 27 13:04:40 2014
 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	93	0.00
2	1,4-Dioxane	1.000	1.075	-7.5	95	0.00
3 I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
4 S	Nitrobenzene-d5	1.000	1.049	-4.9	94	0.00
5	Naphthalene	1.000	1.048	-4.8	92	0.00
6	2-Methylnaphthalene	1.000	1.058	-5.8	91	0.00
7 I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
8 S	2-Fluorobiphenyl	1.000	1.052	-5.2	91	0.00
9	Acenaphthylene	1.000	1.086	-8.6	91	0.00
10	Acenaphthene	1.000	1.032	-3.2	90	0.00
11	Fluorene	1.000	1.087	-8.7	91	0.00
12 I	Phenanthrene-d10	5.000	5.000	0.0	91	0.00
13	Phenanthrene	1.000	1.035	-3.5	90	0.00
14	Anthracene	1.000	1.081	-8.1	92	0.00
15	Fluoranthene	1.000	1.100	-10.0	94	0.00
16 I	Chrysene-d12	5.000	5.000	0.0	96	0.00
17	Pyrene	1.000	1.060	-6.0	95	0.00
18 S	Terphenyl-d14	1.000	1.057	-5.7	96	0.00
19	Benzo(a)anthracene	1.000	1.037	-3.7	98	0.00
20	Chrysene	1.000	1.018	-1.8	95	0.00
21	Indeno(1,2,3-cd)pyrene	1.000	0.885	11.5	89	0.00
22 I	Perylene-d12	5.000	5.000	0.0	96	0.00
23	Benzo(b)fluoranthene	1.000	0.944	5.6	90	0.00
24	Benzo(k)fluoranthene	1.000	1.047	-4.7	95	0.00
25 C	Benzo(a)pyrene	1.000	0.968	3.2	94	0.00
26	Dibenzo(a,h)anthracene	1.000	0.948	5.2	93	0.00
27	Benzo(g,h,i)perylene	1.000	1.019	-1.9	95	0.00

(#) = Out of Range

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