

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102114\
 Data File : BE088158.D
 Acq On : 22 Oct 2014 00:09
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Oct 22 05:06:57 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 21 02:11:14 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	75	0.00
2	1,4-Dioxane	1.000	1.040	-4.0	74	-0.08
3 I	Naphthalene-d8	5.000	5.000	0.0	79	0.00
4 S	Nitrobenzene-d5	1.000	1.134	-13.4	86	0.00
5	Naphthalene	1.000	1.014	-1.4	77	0.00
6	2-Methylnaphthalene	1.000	1.041	-4.1	78	0.00
7 I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
8 S	2-Fluorobiphenyl	1.000	1.649	-64.9#	143	0.00
9	Acenaphthylene	1.000	0.935	6.5	81	0.01
10	Acenaphthene	1.000	0.916	8.4	82	0.00
11	Fluorene	1.000	0.955	4.5	83	0.00
12 I	Phenanthrene-d10	5.000	5.000	0.0	78	0.00
13	Phenanthrene	1.000	0.995	0.5	67	0.00
14	Anthracene	1.000	1.072	-7.2	82	0.00
15	Fluoranthene	1.000	0.998	0.2	78	0.00
16 I	Chrysene-d12	5.000	5.000	0.0	74	0.00
17	Pyrene	1.000	1.021	-2.1	76	0.00
18 S	Terphenyl-d14	1.000	2.481	-148.1#	182	0.00
19	Benzo(a)anthracene	1.000	0.846	15.4	65	0.00
20	Chrysene	1.000	0.938	6.2	70	0.00
21	Indeno(1,2,3-cd)pyrene	1.000	0.654	34.6#	49	0.00
22 I	Perylene-d12	5.000	5.000	0.0	62	0.00
23	Benzo(b)fluoranthene	1.000	0.796	20.4	49	0.00
24	Benzo(k)fluoranthene	1.000	0.934	6.6	58	0.00
25 C	Benzo(a)pyrene	1.000	0.805	19.5	49	0.00
26	Dibenzo(a,h)anthracene	1.000	0.770	23.0	48	0.00
27	Benzo(g,h,i)perylene	1.000	0.856	14.4	53	0.00

(#) = Out of Range

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