

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102214\
 Data File : BE088198.D
 Acq On : 23 Oct 2014 00:24
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Oct 27 04:56:57 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	AvgRF	CCRF	%Dev	Area%	Dev(min)

1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	84	0.00
2	1,4-Dioxane	0.719	0.708	1.5	79	-0.07
3 I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
4 S	Nitrobenzene-d5	0.360	0.733	-103.6#	174#	0.00
5	Naphthalene	1.031	1.006	2.4	83	0.00
6	2-Methylnaphthalene	0.617	0.614	0.5	84	0.00
7 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
8 S	2-Fluorobiphenyl	1.381	4.845	-250.8#	334#	0.00
9	Acenaphthylene	7.619	6.883	9.7	86	0.01
10	Acenaphthene	1.176	1.045	11.1	87	0.00
11	Fluorene	1.246	1.106	11.2	85	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	79	0.00
13	Phenanthrene	4.779	4.408	7.8	69	0.00
14	Anthracene	3.902	4.082	-4.6	81	0.00
15	Fluoranthene	0.951	0.943	0.8	79	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	74	0.00
17	Pyrene	1.359	1.349	0.7	73	0.00
18 S	Terphenyl-d14	0.805	3.699	-359.5#	335#	0.00
19	Benzo(a)anthracene	4.447	3.894	12.4	66	0.00
20	Chrysene	4.363	4.019	7.9	68	0.00
21	Indeno(1,2,3-cd)pyrene	4.846	4.483	7.5	69	0.00
22 I	Perylene-d12	1.000	1.000	0.0	75	0.00
23	Benzo(b)fluoranthene	1.190	1.034	13.1	65	0.00
24	Benzo(k)fluoranthene	1.274	1.096	14.0	65	0.00
25 C	Benzo(a)pyrene	1.096	0.958	12.6	64	0.00
26	Dibenzo(a,h)anthracene	1.075	0.891	17.1	62	0.00
27	Benzo(g,h,i)perylene	4.752	4.338	8.7	68	0.00

(#) = Out of Range

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