

Data Path : Z:\HPCHEM1\BNA_M\Data\bm100615\
 Data File : BM002852.D
 Acq On : 06 Oct 2015 10:23
 Operator : UM/IZ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Oct 06 15:25:36 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM100515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 19:04:11 2015
 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	103	0.00
2 SURR	1,4-Dioxane-d8	0.422	0.420	0.5	104	0.00
3	1,4-Dioxane	0.487	0.480	1.4	103	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
5	Naphthalene	1.063	1.063	0.0	102	0.00
6 SURR	2-Methylnaphthalene-d10	0.592	0.568	4.1	101	0.00
7	2-Methylnaphthalene	0.725	0.719	0.8	102	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
9	Acenaphthylene	1.510	1.556	-3.0	107	0.00
10 C	Acenaphthene	1.516	1.499	1.1	101	0.00
11	Fluorene	1.709	1.699	0.6	101	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.02
13	Pentachlorophenol	0.025	0.017	32.0#	74	0.00
14	Phenanthrene	1.227	1.223	0.3	97	0.00
15	Anthracene	1.179	1.159	1.7	97	0.00
16 SURR	Fluoranthene-d10	0.989	0.950	3.9	94	0.00
17 C	Fluoranthene	1.383	1.327	4.0	94	-0.01
18 I	Chrysene-d12	1.000	1.000	0.0	86	-0.01
19	Pyrene	1.586	1.702	-7.3	92	0.00
20	Benzo(a)anthracene	1.336	1.332	0.3	86	-0.01
21	Chrysene	1.309	1.318	-0.7	86	0.00
22 I	Perylene-d12	1.000	1.000	0.0	82	0.00
23	Benzo(b)fluoranthene	1.473	1.568	-6.4	86	0.00
24	Benzo(k)fluoranthene	1.431	1.349	5.7	80	0.00
25 C	Benzo(a)pyrene	1.340	1.361	-1.6	83	0.00
26	Indeno(1,2,3-cd)pyrene	1.364	1.422	-4.3	87	0.00
27	Dibenzo(a,h)anthracene	1.080	1.138	-5.4	89	-0.01
28	Benzo(g,h,i)perylene	1.176	1.210	-2.9	87	-0.01

(#) = Out of Range

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