

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021715\
 Data File : BM000540.D
 Acq On : 17 Feb 2015 14:36
 Operator : TP/IZ
 Sample : SSTD0.416
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC0.4EC

Quant Time: Feb 17 16:07:54 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:45:55 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	100	-0.02
2 I	Naphthalene-d8	1.000	1.000	0.0	99	-0.02
3	Naphthalene	0.938	0.935	0.3	99	-0.02
4 SURR	2-Methylnaphthalene-d10	0.510	0.486	4.7	95	-0.02
5	2-Methylnaphthalene	0.634	0.612	3.5	96	-0.02
6 I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.02
7	Acenaphthylene	1.505	1.427	5.2	92	-0.02
8 C	Acenaphthene	1.155	1.129	2.3	92	-0.03
9	Fluorene	1.396	1.302	6.7	90	-0.02
10 I	Phenanthrene-d10	1.000	1.000	0.0	88	-0.02
11	Pentachlorophenol	0.051	0.031	39.2#	59	-0.03
12	Phenanthrene	1.086	1.047	3.6	86	-0.02
13	Anthracene	0.977	0.939	3.9	85	-0.02
14 SURR	Fluoranthene-d10	0.841	0.808	3.9	83	-0.01
15 C	Fluoranthene	1.211	1.172	3.2	83	-0.02
16 I	Chrysene-d12	1.000	1.000	0.0	79	-0.02
17	Pyrene	1.439	1.441	-0.1	83	-0.02
18	Benzo(a)anthracene	1.128	1.057	6.3	75	-0.02
19	Chrysene	1.300	1.327	-2.1	79	-0.02
20 I	Perylene-d12	1.000	1.000	0.0	76	-0.03
21	Benzo(b)fluoranthene	1.352	1.308	3.3	74	-0.03
22	Benzo(k)fluoranthene	1.345	1.386	-3.0	81	-0.03
23 C	Benzo(a)pyrene	1.264	1.230	2.7	75	-0.03
24	Indeno(1,2,3-cd)pyrene	1.410	1.394	1.1	75	-0.04
25	Dibenzo(a,h)anthracene	1.114	1.123	-0.8	77	-0.05
26	Benzo(g,h,i)perylene	1.244	1.232	1.0	76	-0.05

(#) = Out of Range

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