

Data Path : Z:\HPCHEM1\BNA_E\Data\BE022017\
 Data File : BE092746.D
 Acq On : 20 Feb 2017 14:07
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 21 10:16:00 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE021517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 15 17:05:20 2017
 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	98	0.00
2	1,4-Dioxane	0.400	0.366	8.5	89	0.00
3	n-Nitrosodimethylamine	0.400	0.409	-2.2	72	0.02
4 S	2-Fluorophenol	0.400	0.320	20.0	81	0.01
5 S	Phenol-d6	0.400	0.348	13.0	88	0.00
6	bis(2-Chloroethyl)ether	0.400	0.359	10.3	95	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	98	0.00
8 S	Nitrobenzene-d5	0.400	0.383	4.3	95	0.00
9	Naphthalene	0.400	0.416	-4.0	101	0.00
10	Hexachlorobutadiene	0.400	0.415	-3.7	98	0.00
11	2-Methylnaphthalene	0.400	0.391	2.3	98	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
13 S	2,4,6-Tribromophenol	0.400	0.351	12.3	90	0.01
14 S	2-Fluorobiphenyl	0.400	0.408	-2.0	101	0.00
15	Acenaphthylene	0.400	0.398	0.5	100	0.00
16	Acenaphthene	0.400	0.405	-1.3	100	-0.02
17	Fluorene	0.400	0.395	1.3	100	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
19	Hexachlorobenzene	0.400	0.428	-7.0	98	0.00
20	Pentachlorophenol	0.400	0.366	8.5	46	0.12
21	Phenanthrene	0.400	0.405	-1.3	97	0.00
22	Anthracene	0.400	0.385	3.8	99	0.02
23	Fluoranthene	0.400	0.388	3.0	98	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	93	0.00
25	Pyrene	0.400	0.404	-1.0	96	0.00
26 S	Terphenyl-d14	0.400	0.400	0.0	95	0.00
27	Benzo(a)anthracene	0.400	0.411	-2.7	97	0.00
28	Chrysene	0.400	0.443	-10.7	103	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.380	5.0	87	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.410	-2.5	81	0.00
31 I	Perylene-d12	5.000	5.000	0.0	96	0.00
32	Benzo(b)fluoranthene	0.400	0.416	-4.0	89	0.00
33	Benzo(k)fluoranthene	0.400	0.367	8.3	89	0.00
34 C	Benzo(a)pyrene	0.400	0.348	13.0	89	-0.01
35	Dibenzo(a,h)anthracene	0.400	0.361	9.8	80	0.00
36	Benzo(g,h,i)perylene	0.400	0.376	6.0	83	0.00

(#) = Out of Range

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