

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092216\
 Data File : BE092406.D
 Acq On : 22 Sep 2016 21:33
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 23 03:42:56 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 18:25:56 2016
 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	91	-0.02
2	1,4-Dioxane	0.400	0.393	1.8	94	-0.01
3	n-Nitrosodimethylamine	0.400	0.441	-10.2	88	0.00
4 S	2-Fluorophenol	0.400	0.413	-3.2	88	0.00
5 S	Phenol-d6	0.400	0.335	16.3	80	0.00
6	bis(2-Chloroethyl)ether	0.400	0.342	14.5	81	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	87	0.00
8 S	Nitrobenzene-d5	0.400	0.410	-2.5	84	0.00
9	Naphthalene	0.400	0.398	0.5	87	0.00
10	Hexachlorobutadiene	0.400	0.402	-0.5	88	0.00
11	2-Methylnaphthalene	0.400	0.392	2.0	86	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	88	0.00
13 S	2,4,6-Tribromophenol	0.400	0.428	-7.0	86	0.01
14 S	2-Fluorobiphenyl	0.400	0.392	2.0	88	0.01
15	Acenaphthylene	0.400	0.384	4.0	87	0.00
16	Acenaphthene	0.400	0.393	1.8	84	0.00
17	Fluorene	0.400	0.393	1.8	88	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	89	0.02
19	Hexachlorobenzene	0.400	0.374	6.5	89	0.00
20	Pentachlorophenol	0.400	0.362	9.5	78	0.00
21	Phenanthrene	0.400	0.394	1.5	92	0.02
22	Anthracene	0.400	0.369	7.8	90	0.02
23	Fluoranthene	0.400	0.372	7.0	89	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	91	0.00
25	Pyrene	0.400	0.380	5.0	90	0.02
26 S	Terphenyl-d14	0.400	0.381	4.8	88	0.02
27	Benzo(a)anthracene	0.400	0.378	5.5	88	0.00
28	Chrysene	0.400	0.396	1.0	84	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.342	14.5	77	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.370	7.5	85	0.05
31 I	Perylene-d12	5.000	5.000	0.0	90	0.04
32	Benzo(b)fluoranthene	0.400	0.397	0.8	90	0.04
33	Benzo(k)fluoranthene	0.400	0.373	6.8	90	0.04
34 C	Benzo(a)pyrene	0.400	0.380	5.0	90	0.02
35	Dibenzo(a,h)anthracene	0.400	0.372	7.0	87	0.07
36	Benzo(g,h,i)perylene	0.400	0.387	3.3	90	0.05

(#) = Out of Range

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