Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081016\
Data File : BE092261.D
Acq On : 10 Aug 2016 18:58
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc : confirmation run
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 11 01:28:49 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 10 11:12:32 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	147	0.00
2		1,4-Dioxane	0.638	0.633	0.8	136	0.00
3		n-Nitrosodimethylamine	0.775	0.779	-0.5	165#	-0.01
4	S	2-Fluorophenol	1.094	1.128	-3.1	167#	0.00
5	S	Phenol-d6	1.573	1.522	3.2	182#	-0.02
6		bis(2-Chloroethyl)ether	1.275	1.365	-7.1	174#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	147	0.00
8	S	Nitrobenzene-d5	0.353	0.332	5.9	163#	-0.02
9		Naphthalene	1.121	1.116	0.4	146	-0.02
10		Hexachlorobutadiene	0.162	0.159	1.9	145	-0.02
11		2-Methylnaphthalene	0.700	0.701	-0.1	150	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	140	0.00
13	S	2,4,6-Tribromophenol	0.150	0.136	9.3	165#	0.00
14	S	2-Fluorobiphenyl	1.545	1.555	-0.6	145	-0.01
15		Acenaphthylene	8.647	8.674	-0.3	147	0.00
16		Acenaphthene	1.388	1.380	0.6	142	0.00
17		Fluorene	1.677	1.680	-0.2	144	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	141	0.00
19		Hexachlorobenzene	0.198	0.189	4.5	135	0.00
20		Pentachlorophenol	0.042	0.046	-9.5	188#	0.00
21		Phenanthrene	1.293	1.207	6.7	137	0.00
22		Anthracene	1.169	1.118	4.4	146	0.00
23		Fluoranthene	5.137	4.982	3.0	144	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	141	-0.02
25		Pyrene	5.914	6.095	-3.1	153#	0.00
26	S	Terphenyl-d14	0.845	0.874	-3.4	158#	0.00
27		Benzo(a)anthracene	6.002	6.252	-4.2	158#	-0.02
28		Chrysene	5.546	5.200	6.2	130	0.00
29		Bis(2-ethylhexyl)phthalate	0.770	0.738	4.2	166#	0.00
30		Indeno(1,2,3-cd)pyrene	5.972	6.170	-3.3	161#	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	148	0.00
32		Benzo(b)fluoranthene	1.313	1.314	-0.1	160#	0.00
33		Benzo(k)fluoranthene	1.323	1.343	-1.5	149	0.00
34	C	Benzo(a)pyrene	1.238	1.257	-1.5	164#	0.00
35		Dibenzo(a,h)anthracene	1.151	1.155	-0.3	165#	0.00
36		Benzo(g,h,i)perylene	4.954	4.879	1.5	153#	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 0