Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032116\
Data File : BE091503.D
Acq On : 21 Mar 2016 17:34
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4

Quant Time: Mar 21 18:10:36 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 14:14:30 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	159#	0.00
2		1,4-Dioxane	0.624	0.553	11.4	148	0.00
3		n-Nitrosodimethylamine	0.887	0.784	11.6	162#	0.00
4	S	2-Fluorophenol	1.149	0.937	18.5	144	0.00
5	S	Phenol-d6	1.500	1.305	13.0	159#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	158#	-0.02
7	S	Nitrobenzene-d5	0.224	0.199	11.2	167#	0.00
8		Naphthalene	1.221	1.177	3.6	163#	0.00
9		2-Methylnaphthalene	0.785	0.700	10.8	154#	-0.01
10	I	Acenaphthene-d10	1.000	1.000	0.0	151#	0.00
11	S	2,4,6-Tribromophenol	0.142	0.148	-4.2	195#	-0.01
12	S	2-Fluorobiphenyl	1.258	1.225	2.6	160#	0.00
13		Acenaphthylene	2.134	1.752	17.9	152#	-0.01
14		Acenaphthene	1.540	1.445	6.2	148	0.00
15		Fluorene	1.949	1.755	10.0	146	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	143	-0.01
17		Hexachlorobenzene	0.218	0.200	8.3	145	0.00
18		Pentachlorophenol	0.086	0.058	32.6#	118	0.00
19		Phenanthrene	1.427	1.356	5.0	146	0.00
20		Anthracene	1.252	1.119	10.6	147	-0.01
21		Fluoranthene	1.638	1.345	17.9	132	0.00
22	I	Chrysene-d12	1.000	1.000	0.0	134	0.00
23		Pyrene	1.244	1.133	8.9	129	-0.02
24	S	Terphenyl-d14	0.601	0.588	2.2	134	-0.02
25		Benzo(a)anthracene	1.376	1.331	3.3	140	0.00
26		Chrysene	1.406	1.476	-5.0	154#	-0.02
27		Indeno(1,2,3-cd)pyrene	1.434	1.280	10.7	132	-0.02
28	I	Perylene-d12	1.000	1.000	0.0	140	-0.01
29		Benzo(b)fluoranthene	1.580	1.283	18.8	127	-0.01
30		Benzo(k)fluoranthene	1.615	1.366	15.4	129	-0.02
31	C	Benzo(a)pyrene	1.392	1.098	21.1#	126	-0.01
32		Dibenzo(a,h)anthracene	1.371	1.168	14.8	135	-0.02
33		Benzo(g,h,i)perylene	1.426	1.191	16.5	129	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1