Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111816\
Data File : BE092570.D
Acq On : 18 Nov 2016 17:35
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 21 05:39:35 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE111616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 16 15:43:18 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	105	-0.02
2		1,4-Dioxane	0.681	0.727	-6.8	117	0.00
3		n-Nitrosodimethylamine	0.654	0.576	11.9	109	0.00
4	S	2-Fluorophenol	0.844	0.870	-3.1	109	0.00
5	S	Phenol-d6	1.104	1.123	-1.7	116	0.00
6		bis(2-Chloroethyl)ether	1.341	1.184	11.7	105	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
8	S	Nitrobenzene-d5	0.269	0.272	-1.1	111	-0.02
9		Naphthalene	1.032	1.027	0.5	107	-0.02
10		Hexachlorobutadiene	0.157	0.159	-1.3	108	0.00
11		2-Methylnaphthalene	0.658	0.651	1.1	112	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
13	S	2,4,6-Tribromophenol	0.112	0.114	-1.8	124	0.00
14	S	2-Fluorobiphenyl	1.219	1.199	1.6	111	0.00
15		Acenaphthylene	6.962	6.676	4.1	114	0.00
16		Acenaphthene	1.136	1.121	1.3	112	0.00
17		Fluorene	1.408	1.399	0.6	114	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
19		Hexachlorobenzene	0.197	0.184	6.6	111	0.00
20		Pentachlorophenol	0.036	0.038	-5.6	146	0.00
21		Phenanthrene	1.119	1.140	-1.9	113	0.00
22		Anthracene	1.053	1.041	1.1	112	0.00
23		Fluoranthene	5.219	5.278	-1.1	115	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
25		Pyrene	3.975	3.976	-0.0	113	0.00
26	S	Terphenyl-d14	0.572	0.556	2.8	110	0.00
27		Benzo(a)anthracene	4.692	4.559	2.8	121	0.00
28		Chrysene	4.250	4.930	-16.0	118	-0.02
29		Bis(2-ethylhexyl)phthalate	0.305	0.283	7.2	116	-0.02
30		Indeno(1,2,3-cd)pyrene	4.722	4.475	5.2	105	0.00
31	I	Perylene-d12	1.000	1.000	0.0	103	-0.01
32		Benzo(b)fluoranthene	1.213	1.135	6.4	100	-0.01
33		Benzo(k)fluoranthene	1.270	1.271	-0.1	109	-0.01
34	C	Benzo(a)pyrene	1.122	1.083	3.5	108	-0.01
35		Dibenzo(a,h)anthracene	1.111	1.043	6.1	104	-0.02
36		Benzo(g,h,i)perylene	4.702	4.440	5.6	102	0.00

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