Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022417\
Data File : BE092755.D
Acq On : 24 Feb 2017 11:53
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 24 15:34:58 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	122	-0.02
2		1,4-Dioxane	0.710	0.706	0.6	122	-0.01
3		n-Nitrosodimethylamine	0.740	0.920	-24.3	173#	-0.05
4	S	2-Fluorophenol	0.973	1.313	-34.9#	171#	-0.03
5	S	Phenol-d6	1.194	1.724	-44.4#	182#	-0.05
6		bis(2-Chloroethyl)ether	1.321	1.558	-17.9	156#	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	114	-0.02
8	S	Nitrobenzene-d5	0.279	0.321	-15.1	136	-0.05
9		Naphthalene	1.082	1.283	-18.6	134	-0.02
10		Hexachlorobutadiene	0.159	0.156	1.9	107	-0.02
11		2-Methylnaphthalene	0.680	0.736	-8.2	127	-0.03
12	I	Acenaphthene-d10	1.000	1.000	0.0	129	-0.02
13	S	2,4,6-Tribromophenol	0.106	0.183	-72.6#	228#	-0.03
14	S	2-Fluorobiphenyl	1.219	1.186	2.7	124	-0.02
15		Acenaphthylene	7.544	7.741	-2.6	134	-0.02
16		Acenaphthene	1.096	1.112	-1.5	130	-0.02
17		Fluorene	1.407	1.592	-13.1	148	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
19		Hexachlorobenzene	0.193	0.190	1.6	125	-0.02
20		Pentachlorophenol	0.042	0.117	-178.6#	456#	-0.12
21		Phenanthrene	1.188	1.226	-3.2	137	-0.02
22		Anthracene	1.094	1.165	-6.5	152#	-0.02
23		Fluoranthene	5.096	6.351	-24.6	174#	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	165#	-0.02
25		Pyrene	4.761	4.907	-3.1	175#	-0.02
26	S	Terphenyl-d14	0.638	0.651	-2.0	174#	-0.02
27		Benzo(a)anthracene	4.849	5.502	-13.5	191#	-0.02
28		Chrysene	5.752	4.831	16.0	139	-0.02
29		Bis(2-ethylhexyl)phthalate	0.674	0.692	-2.7	196#	0.00
30		Indeno(1,2,3-cd)pyrene	6.486	5.024	22.5	127	-0.03
31	I	Perylene-d12	1.000	1.000	0.0	145	-0.02
32		Benzo(b)fluoranthene	1.414	1.369	3.2	145	-0.02
33		Benzo(k)fluoranthene	1.495	1.362	8.9	145	-0.02
34	C	Benzo(a)pyrene	1.285	1.261	1.9	151#	-0.02
35		Dibenzo(a,h)anthracene	1.355	1.186	12.5	129	-0.03
36		Benzo(g,h,i)perylene	5.728	4.952	13.5	126	-0.03

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