Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\be032116\
Data File : BE097503.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 :

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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	159	0.00
2		1,4-Dioxane	0.400	0.354	11.5	148	0.00
3		n-Nitrosodimethylamine	0.400	0.353	11.8	162	0.00
4	S	2-Fluorophenol	0.400	0.326	18.5	144	0.00
5	S	Phenol-d6	0.400	0.348	13.0	159	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	158	-0.02
7	S	Nitrobenzene-d5	0.400	0.498	-24.5	167	0.00
8		Naphthalene	0.400	0.386	3.5	163	0.00
9		2-Methylnaphthalene	0.400	0.357	10.8	154	-0.01
10	I	Acenaphthene-d10	5.000	5.000	0.0	151	0.00
11	S	2,4,6-Tribromophenol	0.400	0.560	-40.0#	195	-0.01
12	S	2-Fluorobiphenyl	0.400	0.389	2.8	160	0.00
13		Acenaphthylene	0.400	0.404	-1.0	152	-0.01
14		Acenaphthene	0.400	0.375	6.3	148	0.00
15		Fluorene	0.400	0.360	10.0	146	-0.01
16	I	Phenanthrene-d10	5.000	5.000	0.0	143	-0.01
17		Hexachlorobenzene	0.400	0.366	8.5	145	0.00
18		Pentachlorophenol	0.400	0.398	0.5	118	0.00
19		Phenanthrene	0.400	0.380	5.0	146	0.00
20		Anthracene	0.400	0.357	10.8	147	-0.01
21		Fluoranthene	0.400	0.328	18.0	132	0.00
22	I	Chrysene-d12	5.000	5.000	0.0	134	0.00
23		Pyrene	0.400	0.364	9.0	129	-0.02
24	S	Terphenyl-d14	0.400	0.391	2.3	134	-0.02
25		Benzo(a)anthracene	0.400	0.387	3.3	140	0.00
26		Chrysene	0.400	0.457	-14.2	154	-0.02
27		Indeno(1,2,3-cd)pyrene	0.400	0.357	10.8	132	-0.02
28	I	Perylene-d12	5.000	5.000	0.0	140	-0.01
29		Benzo(b)fluoranthene	0.400	0.325	18.8	127	-0.01
30		Benzo(k)fluoranthene	0.400	0.338	15.5	129	-0.02
31	C	Benzo(a)pyrene	0.400	0.316	21.0#	126	-0.01
32		Dibenzo(a,h)anthracene	0.400	0.341	14.8	135	-0.02
33		Benzo(g,h,i)perylene	0.400	0.334	16.5	129	-0.01

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