Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080316\
Data File : BE092182.D
Acq On : 4 Aug 2016 14:40
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 04 15:15:49 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-PAH-BE080216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 02 16:59:37 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	99	0.00
2		1,4-Dioxane	0.400	0.375	6.3	89	0.00
3		n-Nitrosodimethylamine	0.400	0.402	-0.5	98	0.00
4	S	2-Fluorophenol	0.400	0.336	16.0	94	0.01
5	S	Phenol-d6	0.400	0.375	6.3	101	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	106	-0.02
7	S	Nitrobenzene-d5	0.400	0.379	5.3	105	0.00
8		Naphthalene	0.400	0.398	0.5	103	0.00
9		2-Methylnaphthalene	0.400	0.404	-1.0	107	0.00
10	I	Acenaphthene-d10	5.000	5.000	0.0	107	-0.02
11	S	2,4,6-Tribromophenol	0.400	0.339	15.3	109	0.00
12	S	2-Fluorobiphenyl	0.400	0.386	3.5	108	0.00
13		Acenaphthylene	0.400	0.380	5.0	108	-0.02
14		Acenaphthene	0.400	0.394	1.5	109	0.00
15		Fluorene	0.400	0.376	6.0	107	-0.01
16	I	Phenanthrene-d10	5.000	5.000	0.0	114	-0.02
17		Hexachlorobenzene	0.400	0.368	8.0	104	0.00
18		Pentachlorophenol	0.400	0.398	0.5	99	0.00
19		Phenanthrene	0.400	0.363	9.3	111	0.00
20		Anthracene	0.400	0.358	10.5	109	0.00
21		Fluoranthene	0.400	0.354	11.5	105	0.00
22	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
23		Pyrene	0.400	0.402	-0.5	108	0.00
24	S	Terphenyl-d14	0.400	0.430	-7.5	116	-0.02
25		Benzo(a)anthracene	0.400	0.376	6.0	100	0.00
26		Chrysene	0.400	0.379	5.3	99	0.00
27		Indeno(1,2,3-cd)pyrene	0.400	0.358	10.5	96	0.00
28	I	Perylene-d12	5.000	5.000	0.0	95	0.00
29		Benzo(b)fluoranthene	0.400	0.367	8.3	86	0.00
30		Benzo(k)fluoranthene	0.400	0.389	2.8	107	-0.01
31	C	Benzo(a)pyrene	0.400	0.395	1.3	98	0.00
32		Dibenzo(a,h)anthracene	0.400	0.381	4.8	96	0.00
33		Benzo(g,h,i)perylene	0.400	0.375	6.3	94	0.00

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