Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081015\
Data File : BE090416.D
Acq On : 11 Aug 2015 00:27
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 11 04:00:42 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 2015
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	130	0.00
2		1,4-Dioxane	0.687	0.679	1.2	120	0.00
3		n-Nitrosodimethylamine	0.773	0.825	-6.7	139	0.00
4	S	2-Fluorophenol	0.750	1.175	-56.7#	218#	0.00
5	S	Phenol-d6	1.241	1.718	-38.4#	194#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	137	0.00
7	S	Nitrobenzene-d5	0.331	0.373	-12.7	162#	0.00
8		Nitrobenzene	0.349	0.381	-9.2	166#	0.00
9		Naphthalene	1.104	1.125	-1.9	140	0.00
10		Hexachlorobutadiene	0.191	0.187	2.1	133	0.00
11		2-Methylnaphthalene	0.640	0.739	-15.5	165#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	148	0.00
13	S	2,4,6-Tribromophenol	0.095	0.166	-74.7#	331#	0.00
14	S	2-Fluorobiphenyl	1.375	1.465	-6.5	155#	0.00
15		Acenaphthylene	8.098	8.982	-10.9	165#	0.00
16		Acenaphthene	1.259	1.272	-1.0	149	0.00
17		Fluorene	1.635	1.668	-2.0	149	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	139	0.00
19		4-Bromophenyl-phenylether	0.194	0.210	-8.2	153#	0.00
20		Hexachlorobenzene	1.005	0.951	5.4	131	0.00
21		Pentachlorophenol	0.044	0.041	6.8	210#	0.00
22		Phenanthrene	1.251	1.310	-4.7	147	0.00
23		Anthracene	1.028	1.210	-17.7	168#	-0.02
24		Fluoranthene	1.231	1.438	-16.8	168#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	141	0.00
26		Pyrene	1.084	1.265	-16.7	170#	0.00
27	S	Terphenyl-d14	0.737	0.867	-17.6	167#	0.00
28		Benzo(a)anthracene	1.108	1.218	-9.9	172#	0.00
29		Chrysene	1.339	1.387	-3.6	142	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.446	-43.4#	280#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.207	-28.1#	218#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	151#	0.00
33		Benzo(b)fluoranthene	0.918	1.092	-19.0	191#	0.00
34		Benzo(k)fluoranthene	1.392	1.409	-1.2	148	0.00
35	C	Benzo(a)pyrene	0.850	1.025	-20.6#	194#	0.00
36		Dibenzo(a,h)anthracene	0.766	1.014	-32.4#	231#	0.00
37		Benzo(g,h,i)perylene	0.923	1.114	-20.7	189#	-0.01

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