Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081215\
Data File : BE090497.D
Acq On : 13 Aug 2015 9:11
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 13 10:51:00 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	137	0.00
2		1,4-Dioxane	0.687	0.675	1.7	125	0.00
3		n-Nitrosodimethylamine	0.773	0.854	-10.5	152#	0.00
4	S	2-Fluorophenol	0.750	1.502	-100.3#	294#	0.00
5	S	Phenol-d6	1.241	2.251	-81.4#	268#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	143	0.00
7	S	Nitrobenzene-d5	0.331	0.433	-30.8#	197#	-0.01
8		Nitrobenzene	0.349	0.451	-29.2#	206#	0.00
9		Naphthalene	1.104	1.119	-1.4	146	0.00
10		Hexachlorobutadiene	0.191	0.183	4.2	136	0.00
11		2-Methylnaphthalene	0.640	0.759	-18.6	177#	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	151#	0.00
13	S	2,4,6-Tribromophenol	0.095	0.247	-160.0#	502#	-0.02
14	S	2-Fluorobiphenyl	1.375	1.490	-8.4	161#	0.00
15		Acenaphthylene	8.098	9.240	-14.1	173#	0.00
16		Acenaphthene	1.259	1.295	-2.9	155#	0.00
17		Fluorene	1.635	1.737	-6.2	158#	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
19		4-Bromophenyl-phenylether	0.194	0.219	-12.9	161#	0.00
20		Hexachlorobenzene	1.005	0.925	8.0	128	-0.02
21		Pentachlorophenol	0.044	0.114	-159.1#	589#	-0.01
22		Phenanthrene	1.251	1.308	-4.6	147	0.00
23		Anthracene	1.028	1.259	-22.5	176#	-0.02
24		Fluoranthene	1.231	1.495	-21.4	176#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	137	0.00
26		Pyrene	1.084	1.366	-26.0#	179#	0.00
27	S	Terphenyl-d14	0.737	0.959	-30.1#	180#	-0.01
28		Benzo(a)anthracene	1.108	1.330	-20.0	183#	0.00
29		Chrysene	1.339	1.356	-1.3	135	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.891	-186.5#	545#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.340	-42.3#	235#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	159#	0.00
33		Benzo(b)fluoranthene	0.918	1.140	-24.2	210#	0.00
34		Benzo(k)fluoranthene	1.392	1.334	4.2	148	0.00
35	C	Benzo(a)pyrene	0.850	1.107	-30.2#	221#	0.00
36		Dibenzo(a,h)anthracene	0.766	1.065	-39.0#	256#	-0.01
37		Benzo(g,h,i)perylene	0.923	1.157	-25.4#	207#	0.00

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