Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081315\
Data File : BE090512.D
Acq On : 13 Aug 2015 22:28
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 14 04:35:45 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	138	0.00
2		1,4-Dioxane	0.687	0.550	19.9	103	0.00
3		n-Nitrosodimethylamine	0.773	0.727	6.0	130	0.00
4	S	2-Fluorophenol	0.750	1.334	-77.9#	263#	0.00
5	S	Phenol-d6	1.241	1.935	-55.9#	232#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	155#	0.00
7	S	Nitrobenzene-d5	0.331	0.330	0.3	163#	0.00
8		Nitrobenzene	0.349	0.343	1.7	170#	0.00
9		Naphthalene	1.104	0.999	9.5	141	0.00
10		Hexachlorobutadiene	0.191	0.144	24.6	116	0.00
11		2-Methylnaphthalene	0.640	0.631	1.4	160#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	155#	0.00
13	S	2,4,6-Tribromophenol	0.095	0.170	-78.9#	355#	0.00
14	S	2-Fluorobiphenyl	1.375	1.263	8.1	140	0.00
15		Acenaphthylene	8.098	8.007	1.1	155#	0.00
16		Acenaphthene	1.259	1.218	3.3	150	0.00
17		Fluorene	1.635	1.484	9.2	139	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	144	0.00
19		4-Bromophenyl-phenylether	0.194	0.169	12.9	129	0.00
20		Hexachlorobenzene	1.005	0.769	23.5	110	0.00
21		Pentachlorophenol	0.044	0.062	-40.9#	328#	0.00
22		Phenanthrene	1.251	1.111	11.2	129	0.00
23		Anthracene	1.028	1.042	-1.4	150#	-0.02
24		Fluoranthene	1.231	1.197	2.8	146	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	123	0.00
26		Pyrene	1.084	1.247	-15.0	147	0.00
27	S	Terphenyl-d14	0.737	0.773	-4.9	130	-0.01
28		Benzo(a)anthracene	1.108	1.124	-1.4	139	0.00
29		Chrysene	1.339	1.124	16.1	101	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.629	-102.3#	346#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.075	-14.1	170#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	132	0.00
33		Benzo(b)fluoranthene	0.918	1.033	-12.5	158#	0.00
34		Benzo(k)fluoranthene	1.392	1.166	16.2	107	0.00
35	C	Benzo(a)pyrene	0.850	0.947	-11.4	157#	0.00
36		Dibenzo(a,h)anthracene	0.766	0.894	-16.7	178#	-0.01
37		Benzo(g,h,i)perylene	0.923	1.014	-9.9	150#	0.00

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