Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081015\
Data File : BE090427.D
Acq On : 11 Aug 2015 7:04
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 11 07:37:47 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	126	0.00
2		1,4-Dioxane	0.687	0.700	-1.9	120	0.00
3		n-Nitrosodimethylamine	0.773	0.826	-6.9	136	0.00
4	S	2-Fluorophenol	0.750	1.138	-51.7#	206#	0.00
5	S	Phenol-d6	1.241	1.669	-34.5#	183#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
7	S	Nitrobenzene-d5	0.331	0.357	-7.9	151#	0.00
8		Nitrobenzene	0.349	0.367	-5.2	157#	0.00
9		Naphthalene	1.104	1.134	-2.7	138	0.00
10		Hexachlorobutadiene	0.191	0.189	1.0	131	0.00
11		2-Methylnaphthalene	0.640	0.723	-13.0	157#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	144	0.00
13	S	2,4,6-Tribromophenol	0.095	0.149	-56.8#	290#	0.00
14	S	2-Fluorobiphenyl	1.375	1.452	-5.6	150#	0.00
15		Acenaphthylene	8.098	8.760	-8.2	158#	0.00
16		Acenaphthene	1.259	1.275	-1.3	146	0.00
17		Fluorene	1.635	1.661	-1.6	145	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	136	0.00
19		4-Bromophenyl-phenylether	0.194	0.199	-2.6	143	0.00
20		Hexachlorobenzene	1.005	0.929	7.6	125	0.00
21		Pentachlorophenol	0.044	0.040	9.1	202#	0.00
22		Phenanthrene	1.251	1.276	-2.0	140	0.00
23		Anthracene	1.028	1.152	-12.1	156#	-0.02
24		Fluoranthene	1.231	1.374	-11.6	157#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	132	0.00
26		Pyrene	1.084	1.255	-15.8	159#	0.00
27	S	Terphenyl-d14	0.737	0.864	-17.2	156#	0.00
28		Benzo(a)anthracene	1.108	1.214	-9.6	161#	0.00
29		Chrysene	1.339	1.386	-3.5	133	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.386	-24.1	227#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.070	-13.6	181#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	141	0.00
33		Benzo(b)fluoranthene	0.918	1.081	-17.8	176#	0.00
34		Benzo(k)fluoranthene	1.392	1.316	5.5	129	0.00
35	C	Benzo(a)pyrene	0.850	0.951	-11.9	167#	0.00
36		Dibenzo(a,h)anthracene	0.766	0.897	-17.1	190#	0.00
37		Benzo(g,h,i)perylene	0.923	1.027	-11.3	162#	0.00

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