Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081115\
Data File : BE090460.D
Acq On : 12 Aug 2015 10:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 12 11:25:24 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	121	0.00
2		1,4-Dioxane	0.687	0.674	1.9	110	0.00
3		n-Nitrosodimethylamine	0.773	0.828	-7.1	130	0.00
4	S	2-Fluorophenol	0.750	1.072	-42.9#	185#	0.00
5	S	Phenol-d6	1.241	1.668	-34.4#	175#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	128	0.00
7	S	Nitrobenzene-d5	0.331	0.359	-8.5	146	0.00
8		Nitrobenzene	0.349	0.371	-6.3	151#	0.00
9		Naphthalene	1.104	1.128	-2.2	131	0.00
10		Hexachlorobutadiene	0.191	0.192	-0.5	127	0.00
11		2-Methylnaphthalene	0.640	0.693	-8.3	144	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
13	S	2,4,6-Tribromophenol	0.095	0.133	-40.0#	244#	0.00
14	S	2-Fluorobiphenyl	1.375	1.423	-3.5	139	0.00
15		Acenaphthylene	8.098	8.406	-3.8	143	0.00
16		Acenaphthene	1.259	1.266	-0.6	137	0.00
17		Fluorene	1.635	1.648	-0.8	136	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	132	0.00
19		4-Bromophenyl-phenylether	0.194	0.195	-0.5	136	0.00
20		Hexachlorobenzene	1.005	0.931	7.4	122	0.00
21		Pentachlorophenol	0.044	0.045	-2.3	219#	0.00
22		Phenanthrene	1.251	1.259	-0.6	134	0.00
23		Anthracene	1.028	1.134	-10.3	150	-0.02
24		Fluoranthene	1.231	1.301	-5.7	145	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	129	0.00
26		Pyrene	1.084	1.197	-10.4	148	0.00
27	S	Terphenyl-d14	0.737	0.817	-10.9	145	-0.01
28		Benzo(a)anthracene	1.108	1.188	-7.2	154#	0.00
29		Chrysene	1.339	1.316	1.7	124	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.322	-3.5	186#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.016	-7.9	169#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	135	0.00
33		Benzo(b)fluoranthene	0.918	1.047	-14.1	163#	0.00
34		Benzo(k)fluoranthene	1.392	1.248	10.3	117	0.00
35	C	Benzo(a)pyrene	0.850	0.925	-8.8	156#	0.00
36		Dibenzo(a,h)anthracene	0.766	0.867	-13.2	176#	-0.01
37		Benzo(g,h,i)perylene	0.923	0.970	-5.1	146	0.00

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