Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE090415\
Data File : BE090721.D
Acq On : 4 Sep 2015 9:46
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 04 13:13:17 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	79	0.00
2		1,4-Dioxane	0.814	0.766	5.9	77	0.00
3		n-Nitrosodimethylamine	1.033	0.869	15.9	70	0.00
4	S	2-Fluorophenol	1.039	0.973	6.4	79	0.00
5	S	Phenol-d6	1.392	1.501	-7.8	90	0.01
6	I	Naphthalene-d8	1.000	1.000	0.0	86	0.00
7	S	Nitrobenzene-d5	0.330	0.323	2.1	84	0.00
8		Nitrobenzene	0.365	0.345	5.5	85	0.00
9		Naphthalene	1.091	1.090	0.1	87	0.00
10		Hexachlorobutadiene	0.171	0.164	4.1	80	0.00
11		2-Methylnaphthalene	0.590	0.658	-11.5	99	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
13	S	2,4,6-Tribromophenol	0.143	0.159	-11.2	130	0.00
14	S	2-Fluorobiphenyl	1.385	1.349	2.6	98	0.00
15		Acenaphthylene	7.992	8.367	-4.7	108	0.00
16		Acenaphthene	1.335	1.360	-1.9	99	0.00
17		Fluorene	1.666	1.732	-4.0	104	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
19		4-Bromophenyl-phenylether	0.170	0.183	-7.6	111	0.00
20		Hexachlorobenzene	0.928	0.922	0.6	100	0.00
21		Pentachlorophenol	0.054	0.051	5.6	107	0.00
22		Phenanthrene	1.237	1.295	-4.7	104	-0.02
23		Anthracene	1.035	1.144	-10.5	114	0.00
24		Fluoranthene	1.304	1.460	-12.0	120	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
26		Pyrene	1.035	1.173	-13.3	124	0.00
27	S	Terphenyl-d14	0.553	0.603	-9.0	119	0.00
28		Benzo(a)anthracene	1.156	1.188	-2.8	113	0.00
29		Chrysene	1.279	1.354	-5.9	111	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.378	-9.6	155#	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.355	-14.7	129	-0.01
32	I	Perylene-d12	1.000	1.000	0.0	113	0.00
33		Benzo(b)fluoranthene	1.084	1.147	-5.8	122	0.00
34		Benzo(k)fluoranthene	1.263	1.327	-5.1	120	0.00
35	C	Benzo(a)pyrene	0.970	1.074	-10.7	133	0.00
36		Dibenzo(a,h)anthracene	1.020	1.110	-8.8	128	-0.01
37		Benzo(g,h,i)perylene	1.116	1.179	-5.6	126	-0.01

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