Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090215\
Data File : BE090710.D
Acq On : 2 Sep 2015 18:36
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 02 23:06:26 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 23:04:24 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	76	0.00
2		1,4-Dioxane	0.814	0.790	2.9	75	0.00
3		n-Nitrosodimethylamine	1.033	0.917	11.2	71	0.00
4	S	2-Fluorophenol	1.039	1.173	-12.9	91	0.00
5	S	Phenol-d6	1.392	1.574	-13.1	90	0.01
6	I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
7	S	Nitrobenzene-d5	0.330	0.333	-0.9	81	0.00
8		Nitrobenzene	0.365	0.363	0.5	83	0.00
9		Naphthalene	1.091	1.091	0.0	81	0.00
10		Hexachlorobutadiene	0.171	0.162	5.3	74	0.00
11		2-Methylnaphthalene	0.590	0.650	-10.2	91	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	85	0.00
13	S	2,4,6-Tribromophenol	0.143	0.171	-19.6	120	0.00
14	S	2-Fluorobiphenyl	1.385	1.409	-1.7	88	0.00
15		Acenaphthylene	7.992	8.535	-6.8	94	0.00
16		Acenaphthene	1.335	1.345	-0.7	84	0.00
17		Fluorene	1.666	1.713	-2.8	88	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
19		4-Bromophenyl-phenylether	0.170	0.171	-0.6	88	0.00
20		Hexachlorobenzene	0.928	0.868	6.5	81	0.00
21		Pentachlorophenol	0.054	0.069	-27.8#	125	0.00
22		Phenanthrene	1.237	1.242	-0.4	86	0.00
23		Anthracene	1.035	1.098	-6.1	94	0.00
24		Fluoranthene	1.304	1.477	-13.3	104	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
26		Pyrene	1.035	1.180	-14.0	111	0.00
27	S	Terphenyl-d14	0.553	0.602	-8.9	105	0.00
28		Benzo(a)anthracene	1.156	1.183	-2.3	100	0.00
29		Chrysene	1.279	1.335	-4.4	97	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.413	-19.7	150#	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.342	-13.6	113	0.00
32	I	Perylene-d12	1.000	1.000	0.0	101	0.00
33		Benzo(b)fluoranthene	1.084	1.129	-4.2	107	0.00
34		Benzo(k)fluoranthene	1.263	1.371	-8.6	111	0.00
35	C	Benzo(a)pyrene	0.970	1.088	-12.2	121	0.00
36		Dibenzo(a,h)anthracene	1.020	1.107	-8.5	113	0.00
37		Benzo(g,h,i)perylene	1.116	1.181	-5.8	112	0.00

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