Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091515\
Data File : BE090828.D
Acq On : 15 Sep 2015 10:56
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 15 13:48:03 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	52	0.00
2		1,4-Dioxane	0.814	0.763	6.3	50	0.00
3		n-Nitrosodimethylamine	1.033	0.776	24.9	41#	0.00
4	S	2-Fluorophenol	1.039	0.979	5.8	52	0.00
5	S	Phenol-d6	1.392	1.141	18.0	45#	0.03
6	I	Naphthalene-d8	1.000	1.000	0.0	54	0.00
7	S	Nitrobenzene-d5	0.330	0.264	20.0	43#	0.00
8		Nitrobenzene	0.365	0.287	21.4	44#	0.00
9		Naphthalene	1.091	1.035	5.1	51	0.00
10		Hexachlorobutadiene	0.171	0.167	2.3	51	0.00
11		2-Methylnaphthalene	0.590	0.544	7.8	51	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	56	0.00
13	S	2,4,6-Tribromophenol	0.143	0.097	32.2#	45#	0.02
14	S	2-Fluorobiphenyl	1.385	1.241	10.4	51	0.00
15		Acenaphthylene	7.992	7.341	8.1	53	0.00
16		Acenaphthene	1.335	1.318	1.3	55	0.00
17		Fluorene	1.666	1.563	6.2	53	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	61	0.00
19		4-Bromophenyl-phenylether	0.170	0.145	14.7	54	0.00
20		Hexachlorobenzene	0.928	0.852	8.2	57	0.00
21		Pentachlorophenol	0.054	0.029	46.3#	37#	0.00
22		Phenanthrene	1.237	1.165	5.8	57	-0.02
23		Anthracene	1.035	0.911	12.0	56	0.00
24		Fluoranthene	1.304	1.135	13.0	57	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	68	0.00
26		Pyrene	1.035	0.858	17.1	60	0.00
27	S	Terphenyl-d14	0.553	0.442	20.1	57	0.00
28		Benzo(a)anthracene	1.156	0.980	15.2	61	0.00
29		Chrysene	1.279	1.220	4.6	66	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.172	50.1#	46#	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.074	9.1	67	0.00
32	I	Perylene-d12	1.000	1.000	0.0	70	-0.01
33		Benzo(b)fluoranthene	1.084	0.967	10.8	64	0.00
34		Benzo(k)fluoranthene	1.263	1.162	8.0	65	0.00
35	C	Benzo(a)pyrene	0.970	0.853	12.1	66	0.00
36		Dibenzo(a,h)anthracene	1.020	0.907	11.1	65	-0.01
37		Benzo(g,h,i)perylene	1.116	0.988	11.5	65	-0.01

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