Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091015\
Data File : BE090757.D
Acq On : 10 Sep 2015 11:06
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 11 03:44:09 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	100	0.00
2		1,4-Dioxane	0.814	0.704	13.5	89	0.00
3		n-Nitrosodimethylamine	1.033	0.762	26.2#	78	0.00
4	S	2-Fluorophenol	1.039	0.745	28.3#	76	0.00
5	S	Phenol-d6	1.392	1.196	14.1	90	0.02
6	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
7	S	Nitrobenzene-d5	0.330	0.264	20.0	88	0.00
8		Nitrobenzene	0.365	0.290	20.5	91	0.00
9		Naphthalene	1.091	0.930	14.8	94	0.00
10		Hexachlorobutadiene	0.171	0.142	17.0	88	0.00
11		2-Methylnaphthalene	0.590	0.544	7.8	105	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
13	S	2,4,6-Tribromophenol	0.143	0.104	27.3#	108	0.00
14	S	2-Fluorobiphenyl	1.385	1.133	18.2	105	0.00
15		Acenaphthylene	7.992	6.798	14.9	112	0.00
16		Acenaphthene	1.335	1.165	12.7	109	0.00
17		Fluorene	1.666	1.419	14.8	108	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	126	0.00
19		4-Bromophenyl-phenylether	0.170	0.146	14.1	112	0.00
20		Hexachlorobenzene	0.928	0.759	18.2	105	0.00
21		Pentachlorophenol	0.054	0.031	42.6#	82	0.00
22		Phenanthrene	1.237	1.071	13.4	109	-0.02
23		Anthracene	1.035	0.909	12.2	115	0.00
24		Fluoranthene	1.304	1.099	15.7	115	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	126	0.00
26		Pyrene	1.035	0.932	10.0	120	0.00
27	S	Terphenyl-d14	0.553	0.485	12.3	116	0.00
28		Benzo(a)anthracene	1.156	0.974	15.7	112	0.00
29		Chrysene	1.279	1.122	12.3	112	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.212	38.6#	106	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.052	10.9	122	0.00
32	I	Perylene-d12	1.000	1.000	0.0	133	0.00
33		Benzo(b)fluoranthene	1.084	0.887	18.2	112	0.00
34		Benzo(k)fluoranthene	1.263	1.111	12.0	119	0.00
35	C	Benzo(a)pyrene	0.970	0.839	13.5	123	0.00
36		Dibenzo(a,h)anthracene	1.020	0.884	13.3	120	-0.01
37		Benzo(g,h,i)perylene	1.116	0.943	15.5	119	-0.01

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