Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100715\
Data File : BE090987.D
Acq On : 7 Oct 2015 14:04
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 08 01:44:51 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 17:11:43 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	176#	0.00
2		1,4-Dioxane	0.888	1.587	-78.7#	367#	0.00
3		n-Nitrosodimethylamine	0.879	0.561	36.2#	124	0.00
4	S	2-Fluorophenol	1.045	1.044	0.1	201#	0.00
5	S	Phenol-d6	1.280	1.226	4.2	189#	0.00
6		bis(2-Chloroethyl)ether	1.524	2.981	-95.6#	357#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	174#	0.00
8	S	Nitrobenzene-d5	0.270	0.362	-34.1#	247#	0.00
9		Nitrobenzene	0.289	0.396	-37.0#	266#	0.00
10		Naphthalene	1.075	1.019	5.2	173#	0.00
11		Hexachlorobutadiene	0.179	0.150	16.2	147	0.00
12		2-Methylnaphthalene	0.634	0.668	-5.4	195#	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	169#	0.00
14	S	2,4,6-Tribromophenol	0.158	0.135	14.6	192#	0.00
15	S	2-Fluorobiphenyl	1.355	1.360	-0.4	182#	0.00
16		Acenaphthylene	8.351	8.074	3.3	178#	0.00
17		Acenaphthene	1.334	1.277	4.3	172#	0.00
18		Fluorene	1.649	1.630	1.2	179#	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	177#	0.00
20		4-Bromophenyl-phenylether	0.189	0.185	2.1	188#	0.00
21		Hexachlorobenzene	0.970	0.202	79.2#	39#	0.00
22		Pentachlorophenol	0.070	0.045	35.7#	177#	0.00
23		Phenanthrene	1.265	1.164	8.0	179#	0.00
24		Anthracene	1.131	1.095	3.2	194#	0.00
25		Fluoranthene	1.503	1.391	7.5	201#	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	195#	0.00
27		Benzidine	0.090	0.156	-73.3#	415#	0.00
28		Pyrene	1.277	1.087	14.9	199#	0.00
29	S	Terphenyl-d14	0.631	0.621	1.6	221#	0.00
30		Benzo(a)anthracene	1.176	1.197	-1.8	221#	0.00
31		3,3'-Dichlorobenzidine	0.201	0.000	100.0#	0#	-20.84#
32		Chrysene	1.505	1.230	18.3	187#	0.00
33		Bis(2-ethylhexyl)phthalate	0.469	0.560	-19.4	367#	0.00
34		Indeno(1,2,3-cd)pyrene	1.212	1.426	-17.7	266#	0.00
35	I	Perylene-d12	1.000	1.000	0.0	239#	0.00
36		Benzo(b)fluoranthene	1.102	1.164	-5.6	281#	0.00
37		Benzo(k)fluoranthene	1.515	1.260	16.8	222#	0.00
38	C	Benzo(a)pyrene	1.022	1.122	-9.8	304#	0.00
39		Dibenzo(a,h)anthracene	1.009	1.282	-27.1#	361#	0.00
40		Benzo(g,h,i)perylene	1.096	1.323	-20.7	332#	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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