Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100815\
Data File : BE090996.D
Acq On : 8 Oct 2015 10:36
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 09 03:05:01 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	172#	0.00
2		1,4-Dioxane	0.888	1.761	-98.3#	398#	-0.07
3		n-Nitrosodimethylamine	0.879	0.569	35.3#	123	-0.10
4	S	2-Fluorophenol	1.045	1.100	-5.3	207#	-0.02
5	S	Phenol-d6	1.280	1.359	-6.2	204#	-0.02
6		bis(2-Chloroethyl)ether	1.524	3.618	-137.4#	424#	-0.04
7	I	Naphthalene-d8	1.000	1.000	0.0	173#	0.00
8	S	Nitrobenzene-d5	0.270	0.371	-37.4#	251#	-0.01
9		Nitrobenzene	0.289	0.415	-43.6#	277#	0.00
10		Naphthalene	1.075	1.022	4.9	172#	0.00
11		Hexachlorobutadiene	0.179	0.147	17.9	142	0.00
12		2-Methylnaphthalene	0.634	0.684	-7.9	198#	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	173#	0.00
14	S	2,4,6-Tribromophenol	0.158	0.149	5.7	218#	0.00
15	S	2-Fluorobiphenyl	1.355	1.337	1.3	184#	0.00
16		Acenaphthylene	8.351	8.213	1.7	186#	0.00
17		Acenaphthene	1.334	1.261	5.5	174#	0.00
18		Fluorene	1.649	1.618	1.9	182#	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	177#	-0.01
20		4-Bromophenyl-phenylether	0.189	0.186	1.6	189#	-0.01
21		Hexachlorobenzene	0.970	0.204	79.0#	39#	0.00
22		Pentachlorophenol	0.070	0.065	7.1	258#	0.00
23		Phenanthrene	1.265	1.181	6.6	181#	-0.01
24		Anthracene	1.131	1.104	2.4	195#	-0.01
25		Fluoranthene	1.503	1.367	9.0	197#	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	186#	0.00
27		Benzidine	0.090	0.192	-113.3#	487#	-0.02
28		Pyrene	1.277	1.114	12.8	195#	0.00
29	S	Terphenyl-d14	0.631	0.623	1.3	211#	0.00
30		Benzo(a)anthracene	1.176	1.184	-0.7	208#	0.00
31		3,3'-Dichlorobenzidine	0.201	0.000	100.0#	0#	-20.84#
32		Chrysene	1.505	1.221	18.9	177#	0.00
33		Bis(2-ethylhexyl)phthalate	0.469	0.538	-14.7	335#	0.00
34		Indeno(1,2,3-cd)pyrene	1.212	1.344	-10.9	239#	0.00
35	I	Perylene-d12	1.000	1.000	0.0	223#	0.00
36		Benzo(b)fluoranthene	1.102	1.232	-11.8	278#	0.00
37		Benzo(k)fluoranthene	1.515	1.147	24.3	188#	0.00
38	C	Benzo(a)pyrene	1.022	1.107	-8.3	280#	0.00
39		Dibenzo(a,h)anthracene	1.009	1.253	-24.2	329#	0.00
40		Benzo(g,h,i)perylene	1.096	1.287	-17.4	302#	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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