Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091815\
Data File : BE090892.D
Acq On : 19 Sep 2015 2:25
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 19 03:47:26 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 18 18:17:34 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	113	0.00
2		1,4-Dioxane	0.795	0.724	8.9	101	0.00
3		n-Nitrosodimethylamine	1.067	0.891	16.5	97	0.00
4	S	2-Fluorophenol	1.238	1.009	18.5	93	0.00
5	S	Phenol-d6	1.559	1.248	19.9	93	0.00
6		bis(2-Chloroethyl)ether	1.546	1.385	10.4	111	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
8	S	Nitrobenzene-d5	0.345	0.299	13.3	99	0.00
9		Nitrobenzene	0.383	0.344	10.2	105	0.00
10		Naphthalene	1.126	1.033	8.3	105	0.00
11		Hexachlorobutadiene	0.166	0.165	0.6	111	0.00
12		2-Methylnaphthalene	0.672	0.605	10.0	104	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
14	S	2,4,6-Tribromophenol	0.196	0.135	31.1#	93	-0.02
15	S	2-Fluorobiphenyl	1.418	1.305	8.0	107	0.00
16		Acenaphthylene	8.705	7.758	10.9	106	0.00
17		Acenaphthene	1.341	1.257	6.3	111	0.00
18		Fluorene	1.705	1.550	9.1	110	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	123	0.00
20		4-Bromophenyl-phenylether	0.193	0.170	11.9	109	-0.02
21		Hexachlorobenzene	0.906	0.836	7.7	113	0.00
22		Pentachlorophenol	0.087	0.056	35.6#	86	0.00
23		Phenanthrene	1.292	1.148	11.1	112	0.00
24		Anthracene	1.144	0.977	14.6	108	0.00
25		Fluoranthene	1.525	1.327	13.0	108	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	113	0.00
27		Pyrene	1.141	1.018	10.8	108	0.00
28	S	Terphenyl-d14	0.638	0.540	15.4	105	0.00
29		Benzo(a)anthracene	1.235	1.123	9.1	108	0.00
30		Chrysene	1.352	1.248	7.7	105	0.00
31		Bis(2-ethylhexyl)phthalate	0.515	0.379	26.4#	95	0.00
32		Indeno(1,2,3-cd)pyrene	1.455	1.257	13.6	97	0.00
33	I	Perylene-d12	1.000	1.000	0.0	110	0.00
34		Benzo(b)fluoranthene	1.170	1.060	9.4	102	0.00
35		Benzo(k)fluoranthene	1.347	1.214	9.9	99	0.00
36	C	Benzo(a)pyrene	1.113	0.972	12.7	99	0.00
37		Dibenzo(a,h)anthracene	1.158	1.013	12.5	97	0.00
38		Benzo(g,h,i)perylene	1.240	1.104	11.0	99	0.00

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