Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE070716\
Data File : BE091989.D
Acq On : 7 Jul 2016 20:20
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 08 01:31:07 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE070516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 05 16:14:58 2016
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	95	0.00
2		1,4-Dioxane	0.678	0.664	2.1	89	0.00
3		n-Nitrosodimethylamine	0.768	0.729	5.1	96	-0.02
4	S	2-Fluorophenol	1.196	1.116	6.7	94	0.00
5	S	Phenol-d6	1.662	1.458	12.3	92	0.00
6		bis(2-Chloroethyl)ether	1.449	1.387	4.3	98	-0.02
7		n-Nitroso-di-n-propylamine	1.387	1.216	12.3	89	-0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
9	S	Nitrobenzene-d5	0.341	0.309	9.4	91	-0.02
10		Nitrobenzene	0.370	0.327	11.6	93	0.00
11		bis(2-Chloroethoxy)methane	0.418	0.438	-4.8	90	-0.03
12		Naphthalene	1.164	1.139	2.1	94	0.00
13		Hexachlorobutadiene	0.177	0.168	5.1	87	0.00
14		2-Methylnaphthalene	0.748	0.708	5.3	91	-0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
16	S	2,4,6-Tribromophenol	0.143	0.121	15.4	91	0.00
17	S	2-Fluorobiphenyl	1.220	1.225	-0.4	89	0.00
18		Acenaphthylene	6.276	6.215	1.0	93	0.00
19		2,6-Dinitrotoluene	1.093	0.808	26.1#	93	0.00
20		Acenaphthene	1.367	1.267	7.3	83	0.00
21		2,4-Dinitrotoluene	1.115	0.744	33.3#	93	0.00
22		Fluorene	1.442	1.430	0.8	89	0.00
23		4-Chlorophenyl-phenylether	0.719	0.741	-3.1	90	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
25		4-Bromophenyl-phenylether	0.217	0.198	8.8	88	0.00
26		Hexachlorobenzene	0.216	0.207	4.2	93	0.00
27		Pentachlorophenol	0.078	0.051	34.6#	111	0.00
28		Phenanthrene	1.303	1.275	2.1	93	0.00
29		Anthracene	1.218	1.120	8.0	91	0.00
30		Fluoranthene	5.599	5.204	7.1	92	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
32		Benzidine	0.333	0.176	47.1#	69	0.00
33		Pyrene	3.640	3.041	16.5	92	0.00
34	S	Terphenyl-d14	0.590	0.492	16.6	92	0.00
35		Benzo(a)anthracene	0.837	0.794	5.1	106	0.00
36		3,3'-Dichlorobenzidine	0.257	0.173	32.7#	75	0.00
37		Chrysene	1.187	1.215	-2.4	115	0.00
38		Bis(2-ethylhexyl)phthalate	0.970	0.595	38.7#	74	-0.03
39		Indeno(1,2,3-cd)pyrene	4.657	3.634	22.0	84	-0.02
40	I	Perylene-d12	1.000	1.000	0.0	96	0.00
41		Benzo(b)fluoranthene	1.381	1.335	3.3	94	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.335	1.197	10.3	88	0.00
43 C	Benzo(a)pyrene	1.304	1.166	10.6	87	0.00
44	Dibenzo(a,h)anthracene	1.227	1.054	14.1	85	0.00
45	Benzo(g,h,i)perylene	5.004	4.398	12.1	85	0.00

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