Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE070516\
Data File : BE091982.D
Acq On : 6 Jul 2016 4:06
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 06 06:52:07 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	89	0.00
2		1,4-Dioxane	0.678	0.742	-9.4	93	0.00
3		n-Nitrosodimethylamine	0.768	0.731	4.8	90	0.00
4	S	2-Fluorophenol	1.196	1.126	5.9	89	0.00
5	S	Phenol-d6	1.662	1.454	12.5	86	0.00
6		bis(2-Chloroethyl)ether	1.449	1.420	2.0	94	0.00
7		n-Nitroso-di-n-propylamine	1.387	1.190	14.2	82	-0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
9	S	Nitrobenzene-d5	0.341	0.312	8.5	83	-0.02
10		Nitrobenzene	0.370	0.332	10.3	86	0.00
11		bis(2-Chloroethoxy)methane	0.418	0.464	-11.0	86	0.00
12		Naphthalene	1.164	1.187	-2.0	89	0.00
13		Hexachlorobutadiene	0.177	0.179	-1.1	84	0.00
14		2-Methylnaphthalene	0.748	0.740	1.1	87	-0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	84	0.00
16	S	2,4,6-Tribromophenol	0.143	0.121	15.4	84	0.00
17	S	2-Fluorobiphenyl	1.220	1.253	-2.7	84	0.00
18		Acenaphthylene	6.276	6.175	1.6	85	0.00
19		2,6-Dinitrotoluene	1.093	0.762	30.3#	81	0.00
20		Acenaphthene	1.367	1.359	0.6	81	0.00
21		2,4-Dinitrotoluene	1.115	0.715	35.9#	82	0.00
22		Fluorene	1.442	1.504	-4.3	85	0.00
23		4-Chlorophenyl-phenylether	0.719	0.745	-3.6	83	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
25		4-Bromophenyl-phenylether	0.217	0.210	3.2	85	0.00
26		Hexachlorobenzene	0.216	0.229	-6.0	93	0.00
27		Pentachlorophenol	0.078	0.053	32.1#	105	0.00
28		Phenanthrene	1.303	1.325	-1.7	88	0.00
29		Anthracene	1.218	1.182	3.0	88	0.00
30		Fluoranthene	5.599	5.631	-0.6	91	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
32		Benzidine	0.333	0.167	49.8#	60	0.00
33		Pyrene	3.640	3.334	8.4	93	0.00
34	S	Terphenyl-d14	0.590	0.526	10.8	90	0.00
35		Benzo(a)anthracene	0.837	0.955	-14.1	117	0.00
36		3,3'-Dichlorobenzidine	0.257	0.186	27.6#	73	0.00
37		Chrysene	1.187	1.376	-15.9	119	0.00
38		Bis(2-ethylhexyl)phthalate	0.970	0.761	21.5	87	0.00
39		Indeno(1,2,3-cd)pyrene	4.657	4.478	3.8	95	-0.02
40	I	Perylene-d12	1.000	1.000	0.0	101	0.00
41		Benzo(b)fluoranthene	1.381	1.376	0.4	101	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.335	1.276	4.4	98	0.00
43 C	Benzo(a)pyrene	1.304	1.242	4.8	97	0.00
44	Dibenzo(a,h)anthracene	1.227	1.148	6.4	96	0.00
45	Benzo(g,h,i)perylene	5.004	4.700	6.1	95	0.00

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