Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be072216\
Data File : BE092009.D
Acq On : 22 Jul 2016 16:53
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 22 17:26:54 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 22 15:39:27 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	90	0.00
2		1,4-Dioxane	0.676	0.705	-4.3	88	0.00
3		n-Nitrosodimethylamine	0.753	0.689	8.5	90	0.00
4	S	2-Fluorophenol	0.969	0.941	2.9	95	0.00
5	S	Phenol-d6	1.312	1.270	3.2	94	0.00
6		bis(2-Chloroethyl)ether	1.436	1.315	8.4	89	0.00
7		n-Nitroso-di-n-propylamine	1.226	1.138	7.2	91	0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
9	S	Nitrobenzene-d5	0.301	0.294	2.3	90	0.00
10		Nitrobenzene	0.341	0.315	7.6	90	0.00
11		bis(2-Chloroethoxy)methane	0.377	0.358	5.0	91	0.00
12		Naphthalene	1.136	1.160	-2.1	87	0.00
13		Hexachlorobutadiene	0.163	0.164	-0.6	85	0.00
14		2-Methylnaphthalene	0.712	0.695	2.4	86	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	81	0.00
16	S	2,4,6-Tribromophenol	0.122	0.104	14.8	91	0.00
17	S	2-Fluorobiphenyl	1.315	1.330	-1.1	85	0.00
18		Acenaphthylene	7.691	7.026	8.6	90	0.00
19		2,6-Dinitrotoluene	0.926	0.717	22.6	93	0.00
20		Acenaphthene	1.304	1.316	-0.9	82	0.00
21		2,4-Dinitrotoluene	0.912	0.669	26.6#	89	0.00
22		Fluorene	1.556	1.575	-1.2	89	0.00
23		4-Chlorophenyl-phenylether	0.772	0.807	-4.5	88	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
25		4-Bromophenyl-phenylether	0.208	0.209	-0.5	90	0.00
26		Hexachlorobenzene	0.215	0.219	-1.9	88	0.00
27		Pentachlorophenol	0.053	0.033	37.7#	95	0.00
28		Phenanthrene	1.292	1.302	-0.8	88	-0.01
29		Anthracene	1.171	1.142	2.5	90	0.00
30		Fluoranthene	5.165	5.195	-0.6	94	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	76	0.00
32		Benzidine	0.156	0.180	-15.4	111	0.02
33		Pyrene	3.628	3.831	-5.6	86	0.00
34	S	Terphenyl-d14	0.502	0.565	-12.5	89	0.00
35		Benzo(a)anthracene	1.127	1.266	-12.3	96	0.00
36		3,3'-Dichlorobenzidine	0.169	0.157	7.1	93	0.00
37		Chrysene	1.372	1.503	-9.5	84	0.00
38		Bis(2-ethylhexyl)phthalate	0.811	1.041	-28.4#	119	0.00
39		Indeno(1,2,3-cd)pyrene	3.769	3.852	-2.2	79	0.00
40	I	Perylene-d12	1.000	1.000	0.0	85	0.00
41		Benzo(b)fluoranthene	1.312	1.228	6.4	84	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.337	1.277	4.5	81	0.00
43 C	Benzo(a)pyrene	1.221	1.111	9.0	81	0.00
44	Dibenzo(a,h)anthracene	1.089	0.990	9.1	79	0.00
45	Benzo(g,h,i)perylene	4.463	4.242	5.0	83	0.00

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