Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062416\
Data File : BE091942.D
Acq On : 24 Jun 2016 11:34
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 24 16:14:56 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE061416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 15 12:00:38 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	84	0.00
2		1,4-Dioxane	0.705	0.750	-6.4	87	0.00
3		n-Nitrosodimethylamine	0.802	0.753	6.1	78	0.02
4	S	2-Fluorophenol	0.977	0.962	1.5	84	0.00
5	S	Phenol-d6	1.267	1.183	6.6	82	0.02
6		bis(2-Chloroethyl)ether	1.431	1.402	2.0	83	0.00
7		n-Nitroso-di-n-propylamine	1.137	1.102	3.1	82	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	87	0.00
9	S	Nitrobenzene-d5	0.345	0.313	9.3	79	0.02
10		Nitrobenzene	0.343	0.314	8.5	79	0.02
11		bis(2-Chloroethoxy)methane	0.438	0.460	-5.0	86	0.00
12		Naphthalene	1.092	1.139	-4.3	85	0.00
13		Hexachlorobutadiene	0.177	0.187	-5.6	87	0.00
14		2-Methylnaphthalene	0.741	0.742	-0.1	85	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	88	0.00
16	S	2,4,6-Tribromophenol	0.102	0.083	18.6	76	0.00
17	S	2-Fluorobiphenyl	1.269	1.268	0.1	81	0.00
18		Acenaphthylene	5.075	4.919	3.1	85	0.00
19		2,6-Dinitrotoluene	0.811	0.552	31.9#	84	0.00
20		Acenaphthene	1.398	1.389	0.6	87	0.00
21		2,4-Dinitrotoluene	0.758	0.520	31.4#	64	0.03
22		Fluorene	1.391	1.356	2.5	85	0.00
23		4-Chlorophenyl-phenylether	0.685	0.688	-0.4	83	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	83	0.00
25		4-Bromophenyl-phenylether	0.193	0.201	-4.1	85	0.00
26		Hexachlorobenzene	0.198	0.205	-3.5	82	-0.02
27		Pentachlorophenol	0.054	0.041	24.1	80	0.02
28		Phenanthrene	1.250	1.319	-5.5	84	0.00
29		Anthracene	1.106	1.105	0.1	84	0.00
30		Fluoranthene	5.354	5.343	0.2	82	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	82	0.00
32		Benzidine	0.170	0.158	7.1	106	0.00
33		Pyrene	3.401	3.163	7.0	78	0.00
34	S	Terphenyl-d14	0.612	0.554	9.5	71	0.00
35		Benzo(a)anthracene	0.786	0.838	-6.6	89	0.00
36		3,3'-Dichlorobenzidine	0.127	0.101	20.5	77	0.03
37		Chrysene	1.147	1.283	-11.9	82	0.00
38		Bis(2-ethylhexyl)phthalate	0.580	0.552	4.8	89	-0.03
39		Indeno(1,2,3-cd)pyrene	3.925	3.627	7.6	76	0.00
40	I	Perylene-d12	1.000	1.000	0.0	80	-0.01
41		Benzo(b)fluoranthene	1.284	1.192	7.2	72	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.175	1.281	-9.0	87	0.00
43 C	Benzo(a)pyrene	1.122	1.073	4.4	77	0.00
44	Dibenzo(a,h)anthracene	1.092	1.039	4.9	76	-0.02
45	Benzo(g,h,i)perylene	4.507	4.515	-0.2	79	0.00

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