Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061616\
Data File : BE091935.D
Acq On : 16 Jun 2016 11:19
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 16 15:47:03 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	82	0.00
2		1,4-Dioxane	0.400	0.389	2.8	78	0.00
3		n-Nitrosodimethylamine	0.400	0.343	14.2	70	0.03
4	S	2-Fluorophenol	0.400	0.379	5.3	79	0.02
5	S	Phenol-d6	0.400	0.425	-6.2	80	0.02
6		bis(2-Chloroethyl)ether	0.400	0.351	12.3	73	0.02
7		n-Nitroso-di-n-propylamine	0.400	0.425	-6.2	77	0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	85	0.03
9	S	Nitrobenzene-d5	0.400	0.372	7.0	78	0.02
10		Nitrobenzene	0.400	0.359	10.3	74	0.02
11		bis(2-Chloroethoxy)methane	0.400	0.402	-0.5	79	0.03
12		Naphthalene	0.400	0.389	2.8	77	0.00
13		Hexachlorobutadiene	0.400	0.386	3.5	77	0.00
14		2-Methylnaphthalene	0.400	0.377	5.8	77	0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	82	0.00
16	S	2,4,6-Tribromophenol	0.400	0.343	14.2	74	0.02
17	S	2-Fluorobiphenyl	0.400	0.403	-0.8	77	0.00
18		Acenaphthylene	0.400	0.366	8.5	76	0.03
19		2,6-Dinitrotoluene	0.400	0.475	-18.7	95	0.00
20		Acenaphthene	0.400	0.384	4.0	79	0.00
21		2,4-Dinitrotoluene	0.400	0.371	7.3	73	0.03
22		Fluorene	0.400	0.366	8.5	75	0.02
23		4-Chlorophenyl-phenylether	0.400	0.375	6.3	73	0.02
24	I	Phenanthrene-d10	5.000	5.000	0.0	80	0.00
25		4-Bromophenyl-phenylether	0.400	0.369	7.8	72	0.02
26		Hexachlorobenzene	0.400	0.380	5.0	73	0.00
27		Pentachlorophenol	0.400	0.338	15.5	67	0.02
28		Phenanthrene	0.400	0.383	4.3	73	0.00
29		Anthracene	0.400	0.364	9.0	74	0.00
30		Fluoranthene	0.400	0.376	6.0	74	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	82	0.00
32		Benzidine	0.400	0.373	6.8	91	0.00
33		Pyrene	0.400	0.349	12.8	73	0.00
34	S	Terphenyl-d14	0.400	0.360	10.0	71	0.00
35		Benzo(a)anthracene	0.400	0.393	1.8	82	0.00
36		3,3'-Dichlorobenzidine	0.400	0.319	20.3	79	0.03
37		Chrysene	0.400	0.398	0.5	73	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.450	-12.5	92	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.373	6.8	78	0.00
40	I	Perylene-d12	5.000	5.000	0.0	85	-0.01
41		Benzo(b)fluoranthene	0.400	0.348	13.0	71	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.404	-1.0	85	0.00
43 C	Benzo(a)pyrene	0.400	0.378	5.5	80	0.00
44	Dibenzo(a,h)anthracene	0.400	0.366	8.5	77	-0.02
45	Benzo(g,h,i)perylene	0.400	0.371	7.3	78	-0.02

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