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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	84	0.00
2		1,4-Dioxane	0.400	0.425	-6.2	87	0.00
3		n-Nitrosodimethylamine	0.400	0.376	6.0	78	0.02
4	S	2-Fluorophenol	0.400	0.394	1.5	84	0.00
5	S	Phenol-d6	0.400	0.426	-6.5	82	0.02
6		bis(2-Chloroethyl)ether	0.400	0.392	2.0	83	0.00
7		n-Nitroso-di-n-propylamine	0.400	0.439	-9.7	82	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	87	0.00
9	S	Nitrobenzene-d5	0.400	0.362	9.5	79	0.02
10		Nitrobenzene	0.400	0.369	7.8	79	0.02
11		bis(2-Chloroethoxy)methane	0.400	0.421	-5.2	86	0.00
12		Naphthalene	0.400	0.417	-4.2	85	0.00
13		Hexachlorobutadiene	0.400	0.423	-5.7	87	0.00
14		2-Methylnaphthalene	0.400	0.401	-0.3	85	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	88	0.00
16	S	2,4,6-Tribromophenol	0.400	0.334	16.5	76	0.00
17	S	2-Fluorobiphenyl	0.400	0.400	0.0	81	0.00
18		Acenaphthylene	0.400	0.388	3.0	85	0.00
19		2,6-Dinitrotoluene	0.400	0.437	-9.2	84	0.00
20		Acenaphthene	0.400	0.397	0.8	87	0.00
21		2,4-Dinitrotoluene	0.400	0.330	17.5	64	0.03
22		Fluorene	0.400	0.390	2.5	85	0.00
23		4-Chlorophenyl-phenylether	0.400	0.401	-0.3	83	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	83	0.00
25		4-Bromophenyl-phenylether	0.400	0.417	-4.2	85	0.00
26		Hexachlorobenzene	0.400	0.413	-3.2	82	-0.02
27		Pentachlorophenol	0.400	0.379	5.3	80	0.02
28		Phenanthrene	0.400	0.422	-5.5	84	0.00
29		Anthracene	0.400	0.400	0.0	84	0.00
30		Fluoranthene	0.400	0.399	0.3	82	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	82	0.00
32		Benzidine	0.400	0.419	-4.7	106	0.00
33		Pyrene	0.400	0.372	7.0	78	0.00
34	S	Terphenyl-d14	0.400	0.362	9.5	71	0.00
35		Benzo(a)anthracene	0.400	0.427	-6.7	89	0.00
36		3,3'-Dichlorobenzidine	0.400	0.316	21.0	77	0.03
37		Chrysene	0.400	0.447	-11.7	82	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.440	-10.0	89	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.370	7.5	76	0.00
40	I	Perylene-d12	5.000	5.000	0.0	80	-0.01
41		Benzo(b)fluoranthene	0.400	0.372	7.0	72	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.436	-9.0	87	0.00
43 C	Benzo(a)pyrene	0.400	0.383	4.3	77	0.00
44	Dibenzo(a,h)anthracene	0.400	0.381	4.8	76	-0.02
45	Benzo(g,h,i)perylene	0.400	0.401	-0.3	79	0.00

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