Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE070516\
Data File : BE091967.D
Acq On : 5 Jul 2016 15:55
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE070516

Quant Time: Jul 05 16:51:49 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	91	0.00
2		1,4-Dioxane	0.400	0.434	-8.5	95	0.00
3		n-Nitrosodimethylamine	0.400	0.367	8.3	93	-0.02
4	S	2-Fluorophenol	0.400	0.381	4.8	92	0.00
5	S	Phenol-d6	0.400	0.380	5.0	95	0.00
6		bis(2-Chloroethyl)ether	0.400	0.409	-2.2	101	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.383	4.3	94	-0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	93	0.00
9	S	Nitrobenzene-d5	0.400	0.387	3.3	91	-0.02
10		Nitrobenzene	0.400	0.376	6.0	93	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.401	-0.3	81	0.00
12		Naphthalene	0.400	0.413	-3.2	93	0.00
13		Hexachlorobutadiene	0.400	0.419	-4.7	91	0.00
14		2-Methylnaphthalene	0.400	0.415	-3.7	94	-0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	89	0.00
16	S	2,4,6-Tribromophenol	0.400	0.423	-5.7	94	0.00
17	S	2-Fluorobiphenyl	0.400	0.439	-9.7	95	0.00
18		Acenaphthylene	0.400	0.405	-1.3	93	0.00
19		2,6-Dinitrotoluene	0.400	0.360	10.0	96	0.00
20		Acenaphthene	0.400	0.438	-9.5	95	0.00
21		2,4-Dinitrotoluene	0.400	0.420	-5.0	101	0.00
22		Fluorene	0.400	0.436	-9.0	94	0.00
23		4-Chlorophenyl-phenylether	0.400	0.454	-13.5	96	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	95	0.00
25		4-Bromophenyl-phenylether	0.400	0.411	-2.7	96	0.00
26		Hexachlorobenzene	0.400	0.422	-5.5	98	0.00
27		Pentachlorophenol	0.400	0.363	9.3	100	0.00
28		Phenanthrene	0.400	0.423	-5.7	97	0.00
29		Anthracene	0.400	0.407	-1.7	97	0.00
30		Fluoranthene	0.400	0.426	-6.5	102	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	97	0.00
32		Benzidine	0.400	0.472	-18.0	90	0.00
33		Pyrene	0.400	0.412	-3.0	99	0.00
34	S	Terphenyl-d14	0.400	0.443	-10.7	106	0.00
35		Benzo(a)anthracene	0.400	0.412	-3.0	101	0.00
36		3,3'-Dichlorobenzidine	0.400	0.398	0.5	94	0.00
37		Chrysene	0.400	0.426	-6.5	104	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.369	7.8	98	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.430	-7.5	101	0.00
40	I	Perylene-d12	5.000	5.000	0.0	97	0.00
41		Benzo(b)fluoranthene	0.400	0.402	-0.5	99	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.425	-6.2	105	0.00
43 C	Benzo(a)pyrene	0.400	0.419	-4.7	103	0.00
44	Dibenzo(a,h)anthracene	0.400	0.409	-2.2	102	0.00
45	Benzo(g,h,i)perylene	0.400	0.407	-1.7	100	0.02

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