Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE061416\
Data File : BE091917.D
Acq On : 14 Jun 2016 19:37
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE061416

Quant Time: Jun 15 12:18:22 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 11:58:50 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	105	0.00
2		1,4-Dioxane	0.400	0.417	-4.2	106	0.00
3		n-Nitrosodimethylamine	0.400	0.403	-0.8	105	0.02
4	S	2-Fluorophenol	0.400	0.370	7.5	98	0.00
5	S	Phenol-d6	0.400	0.420	-5.0	100	0.02
6		bis(2-Chloroethyl)ether	0.400	0.406	-1.5	107	0.02
7		n-Nitroso-di-n-propylamine	0.400	0.438	-9.5	103	0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	108	0.00
9	S	Nitrobenzene-d5	0.400	0.394	1.5	106	0.00
10		Nitrobenzene	0.400	0.390	2.5	104	0.02
11		bis(2-Chloroethoxy)methane	0.400	0.434	-8.5	109	0.00
12		Naphthalene	0.400	0.431	-7.7	109	0.00
13		Hexachlorobutadiene	0.400	0.423	-5.7	107	0.00
14		2-Methylnaphthalene	0.400	0.410	-2.5	107	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	109	0.00
16	S	2,4,6-Tribromophenol	0.400	0.351	12.3	101	0.00
17	S	2-Fluorobiphenyl	0.400	0.422	-5.5	107	0.00
18		Acenaphthylene	0.400	0.372	7.0	102	0.03
19		2,6-Dinitrotoluene	0.400	0.458	-14.5	117	0.00
20		Acenaphthene	0.400	0.408	-2.0	111	0.00
21		2,4-Dinitrotoluene	0.400	0.373	6.8	97	0.03
22		Fluorene	0.400	0.400	0.0	108	0.02
23		4-Chlorophenyl-phenylether	0.400	0.414	-3.5	107	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	107	0.00
25		4-Bromophenyl-phenylether	0.400	0.420	-5.0	110	0.02
26		Hexachlorobenzene	0.400	0.421	-5.2	108	0.00
27		Pentachlorophenol	0.400	0.367	8.3	99	0.02
28		Phenanthrene	0.400	0.426	-6.5	109	0.00
29		Anthracene	0.400	0.403	-0.8	109	0.00
30		Fluoranthene	0.400	0.409	-2.2	107	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	95	0.00
32		Benzidine	0.400	0.328	18.0	88	0.00
33		Pyrene	0.400	0.457	-14.2	111	0.00
34	S	Terphenyl-d14	0.400	0.478	-19.5	110	0.00
35		Benzo(a)anthracene	0.400	0.386	3.5	93	0.00
36		3,3'-Dichlorobenzidine	0.400	0.335	16.3	95	0.03
37		Chrysene	0.400	0.413	-3.2	88	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.441	-10.2	104	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.418	-4.5	100	0.00
40	I	Perylene-d12	5.000	5.000	0.0	100	0.00
41		Benzo(b)fluoranthene	0.400	0.409	-2.2	98	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.413	-3.2	103	0.00
43 C	Benzo(a)pyrene	0.400	0.390	2.5	98	0.00
44	Dibenzo(a,h)anthracene	0.400	0.406	-1.5	101	0.00
45	Benzo(g,h,i)perylene	0.400	0.398	0.5	98	0.00

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