Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042016\
Data File : BE091690.D
Acq On : 20 Apr 2016 18:12
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE042016

Quant Time: Apr 20 19:02:46 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 20 18:55:38 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	100	0.00
2		1,4-Dioxane	0.400	0.368	8.0	93	0.00
3		n-Nitrosodimethylamine	0.400	0.405	-1.3	107	0.00
4	S	2-Fluorophenol	0.400	0.391	2.3	102	0.00
5	S	Phenol-d6	0.400	0.373	6.8	100	0.00
6		bis(2-Chloroethyl)ether	0.400	0.387	3.3	100	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
8	S	Nitrobenzene-d5	0.400	0.387	3.3	107	0.00
9		Nitrobenzene	0.400	0.383	4.3	105	-0.01
10		Naphthalene	0.400	0.401	-0.3	98	0.00
11		Hexachlorobutadiene	0.400	0.405	-1.3	99	0.00
12		2-Methylnaphthalene	0.400	0.391	2.3	98	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
14	S	2,4,6-Tribromophenol	0.400	0.419	-4.7	100	-0.01
15	S	2-Fluorobiphenyl	0.400	0.401	-0.3	97	0.00
16		Acenaphthylene	0.400	0.374	6.5	100	0.00
17		Acenaphthene	0.400	0.394	1.5	98	0.00
18		Fluorene	0.400	0.393	1.8	97	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	99	0.00
20		4-Bromophenyl-phenylether	0.400	0.393	1.8	98	0.00
21		Hexachlorobenzene	0.400	0.398	0.5	96	0.00
22		Pentachlorophenol	0.400	0.414	-3.5	104	-0.02
23		Phenanthrene	0.400	0.409	-2.2	97	-0.01
24		Anthracene	0.400	0.384	4.0	99	-0.01
25		Fluoranthene	0.400	0.395	1.3	101	-0.02
26	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
27		Benzidine	0.400	0.475	-18.7	125	-0.02
28		Pyrene	0.400	0.402	-0.5	106	0.00
29	S	Terphenyl-d14	0.400	0.392	2.0	103	0.00
30		Benzo(a)anthracene	0.400	0.408	-2.0	109	0.00
31		3,3'-Dichlorobenzidine	0.400	0.460	-15.0	121	-0.02
32		Chrysene	0.400	0.417	-4.2	105	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.472	-18.0	130	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	105	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	103	-0.01
36		Benzo(b)fluoranthene	0.400	0.389	2.8	106	0.00
37		Benzo(k)fluoranthene	0.400	0.384	4.0	104	-0.01
38	C	Benzo(a)pyrene	0.400	0.377	5.8	106	-0.01
39		Dibenzo(a,h)anthracene	0.400	0.377	5.8	106	-0.01
40		Benzo(g,h,i)perylene	0.400	0.379	5.3	103	-0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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