Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090216\
Data File : BE092350.D
Acq On : 2 Sep 2016 19:14
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE090216

Quant Time: Sep 05 09:08:43 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 02 19:34:51 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	133	0.00
2		1,4-Dioxane	0.400	0.395	1.3	135	0.00
3		n-Nitrosodimethylamine	0.400	0.427	-6.7	170	-0.01
4	S	2-Fluorophenol	0.400	0.388	3.0	152	0.00
5	S	Phenol-d6	0.400	0.376	6.0	162	0.00
6		bis(2-Chloroethyl)ether	0.400	0.419	-4.7	175	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	137	0.00
8	S	Nitrobenzene-d5	0.400	0.375	6.3	163	0.00
9		Naphthalene	0.400	0.417	-4.2	150	0.00
10		Hexachlorobutadiene	0.400	0.422	-5.5	147	0.00
11		2-Methylnaphthalene	0.400	0.410	-2.5	151	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	131	0.00
13	S	2,4,6-Tribromophenol	0.400	0.397	0.8	162	0.00
14	S	2-Fluorobiphenyl	0.400	0.415	-3.7	151	0.00
15		Acenaphthylene	0.400	0.414	-3.5	151	0.00
16		Acenaphthene	0.400	0.416	-4.0	148	0.00
17		Fluorene	0.400	0.415	-3.7	151	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	132	0.02
19		Hexachlorobenzene	0.400	0.407	-1.7	147	0.00
20		Pentachlorophenol	0.400	0.422	-5.5	174	0.00
21		Phenanthrene	0.400	0.393	1.8	142	0.00
22		Anthracene	0.400	0.390	2.5	155	0.02
23		Fluoranthene	0.400	0.416	-4.0	155	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	157	0.00
25		Pyrene	0.400	0.387	3.3	159	0.00
26	S	Terphenyl-d14	0.400	0.373	6.8	155	0.00
27		Benzo(a)anthracene	0.400	0.402	-0.5	161	0.00
28		Chrysene	0.400	0.356	11.0	180	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.320	20.0	148	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.384	4.0	154	0.02
31	I	Perylene-d12	5.000	5.000	0.0	134	0.02
32		Benzo(b)fluoranthene	0.400	0.451	-12.7	171	0.02
33		Benzo(k)fluoranthene	0.400	0.384	4.0	134	0.02
34	C	Benzo(a)pyrene	0.400	0.416	-4.0	156	0.01
35		Dibenzo(a,h)anthracene	0.400	0.410	-2.5	155	0.04
36		Benzo(g,h,i)perylene	0.400	0.406	-1.5	146	0.02

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