Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030816\
Data File : BM004579.D
Acq On : 08 Mar 2016 16:32
Operator : SJ/UM
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM030816

Quant Time: Mar 08 18:24:19 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM030816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 08 18:22:06 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	114	-0.02
2		1,4-Dioxane	0.400	0.412	-3.0	114	0.00
3		n-Nitrosodimethylamine	0.400	0.421	-5.2	125	-0.02
4	S	2-Fluorophenol	0.400	0.393	1.8	112	-0.03
5	S	Phenol-d6	0.400	0.415	-3.7	123	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	111	-0.02
7	S	Nitrobenzene-d5	0.400	0.409	-2.2	126	-0.01
8		Naphthalene	0.400	0.419	-4.7	114	-0.02
9		2-Methylnaphthalene	0.400	0.419	-4.7	119	-0.01
10	I	Acenaphthene-d10	5.000	5.000	0.0	115	0.00
11	S	2,4,6-Tribromophenol	0.400	0.465	-16.3	132	0.00
12	S	2-Fluorobiphenyl	0.400	0.426	-6.5	119	-0.01
13		Acenaphthylene	0.400	0.433	-8.2	121	0.00
14		Acenaphthene	0.400	0.399	0.3	118	0.00
15		Fluorene	0.400	0.429	-7.2	123	-0.02
16	I	Phenanthrene-d10	5.000	5.000	0.0	124	0.00
17		Hexachlorobenzene	0.400	0.320	20.0	115	0.00
18		Pentachlorophenol	0.400	0.212	47.0#	132	-0.03
19		Phenanthrene	0.400	0.341	14.8	122	0.00
20		Anthracene	0.400	0.344	14.0	130	0.00
21		Fluoranthene	0.400	0.381	4.8	137	-0.02
22	I	Chrysene-d12	5.000	5.000	0.0	146	0.00
23		Pyrene	0.400	0.382	4.5	137	0.00
24	S	Terphenyl-d14	0.400	0.388	3.0	139	0.00
25		Benzo(a)anthracene	0.400	0.389	2.8	150	0.00
26		Chrysene	0.400	0.420	-5.0	133	0.00
27		Indeno(1,2,3-cd)pyrene	0.400	0.362	9.5	126	0.00
28	I	Perylene-d12	5.000	5.000	0.0	137	0.01
29		Benzo(b)fluoranthene	0.400	0.403	-0.8	136	0.00
30		Benzo(k)fluoranthene	0.400	0.395	1.3	145	0.00
31	C	Benzo(a)pyrene	0.400	0.423	-5.7	137	0.00
32		Dibenzo(a,h)anthracene	0.400	0.365	8.8	125	0.00
33		Benzo(g,h,i)perylene	0.400	0.369	7.8	124	0.01

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