Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032516\
Data File : BE091530.D
Acq On : 25 Mar 2016 18:26
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE032516

Quant Time: Mar 25 19:06:13 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 25 18:30:57 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	94	0.00
2		1,4-Dioxane	0.400	0.414	-3.5	92	0.00
3		n-Nitrosodimethylamine	0.400	0.409	-2.2	97	0.00
4	S	2-Fluorophenol	0.400	0.394	1.5	99	0.00
5	S	Phenol-d6	0.400	0.379	5.3	99	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
7	S	Nitrobenzene-d5	0.400	0.403	-0.8	111	-0.01
8		Naphthalene	0.400	0.403	-0.8	96	0.00
9		2-Methylnaphthalene	0.400	0.389	2.8	97	0.00
10	I	Acenaphthene-d10	5.000	5.000	0.0	93	-0.01
11	S	2,4,6-Tribromophenol	0.400	0.395	1.3	107	-0.01
12	S	2-Fluorobiphenyl	0.400	0.403	-0.8	96	-0.01
13		Acenaphthylene	0.400	0.376	6.0	99	0.00
14		Acenaphthene	0.400	0.404	-1.0	95	-0.01
15		Fluorene	0.400	0.382	4.5	94	-0.01
16	I	Phenanthrene-d10	5.000	5.000	0.0	95	0.00
17		Hexachlorobenzene	0.400	0.386	3.5	96	0.00
18		Pentachlorophenol	0.400	0.434	-8.5	98	-0.02
19		Phenanthrene	0.400	0.396	1.0	94	0.00
20		Anthracene	0.400	0.372	7.0	98	0.00
21		Fluoranthene	0.400	0.370	7.5	96	-0.02
22	I	Chrysene-d12	5.000	5.000	0.0	94	0.00
23		Pyrene	0.400	0.367	8.3	95	0.00
24	S	Terphenyl-d14	0.400	0.371	7.3	94	0.00
25		Benzo(a)anthracene	0.400	0.366	8.5	95	0.00
26		Chrysene	0.400	0.406	-1.5	95	0.00
27		Indeno(1,2,3-cd)pyrene	0.400	0.362	9.5	96	-0.02
28	I	Perylene-d12	5.000	5.000	0.0	96	-0.01
29		Benzo(b)fluoranthene	0.400	0.358	10.5	97	-0.01
30		Benzo(k)fluoranthene	0.400	0.381	4.8	95	0.00
31	C	Benzo(a)pyrene	0.400	0.477	-19.2	96	-0.01
32		Dibenzo(a,h)anthracene	0.400	0.360	10.0	96	-0.02
33		Benzo(g,h,i)perylene	0.400	0.368	8.0	96	-0.01

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