Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE031816\
Data File : BE091477.D
Acq On : 17 Mar 2016 21:37
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE031816

Quant Time: Mar 18 04:47:47 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 04:36:48 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	77	0.00
2		1,4-Dioxane	0.400	0.510	-27.5#	103	0.00
3		n-Nitrosodimethylamine	0.400	0.475	-18.7	105	0.00
4	S	2-Fluorophenol	0.400	0.483	-20.7	103	0.00
5	S	Phenol-d6	0.400	0.460	-15.0	102	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	76	0.00
7	S	Nitrobenzene-d5	0.400	0.595	-48.7#	109	0.00
8		Naphthalene	0.400	0.494	-23.5	101	0.00
9		2-Methylnaphthalene	0.400	0.475	-18.7	98	0.00
10	I	Acenaphthene-d10	5.000	5.000	0.0	75	0.00
11	S	2,4,6-Tribromophenol	0.400	0.569	-42.2#	100	0.00
12	S	2-Fluorobiphenyl	0.400	0.475	-18.7	98	0.00
13		Acenaphthylene	0.400	0.497	-24.2	102	0.00
14		Acenaphthene	0.400	0.497	-24.2	98	0.00
15		Fluorene	0.400	0.490	-22.5	99	0.00
16	I	Phenanthrene-d10	5.000	5.000	0.0	77	0.00
17		Hexachlorobenzene	0.400	0.474	-18.5	101	0.00
18		Pentachlorophenol	0.400	0.585	-46.2#	100	0.00
19		Phenanthrene	0.400	0.491	-22.7	101	0.00
20		Anthracene	0.400	0.464	-16.0	102	0.00
21		Fluoranthene	0.400	0.481	-20.2	103	0.00
22	I	Chrysene-d12	5.000	5.000	0.0	83	0.00
23		Pyrene	0.400	0.466	-16.5	102	0.00
24	S	Terphenyl-d14	0.400	0.474	-18.5	100	0.00
25		Benzo(a)anthracene	0.400	0.475	-18.7	106	0.00
26		Chrysene	0.400	0.440	-10.0	107	0.00
27		Indeno(1,2,3-cd)pyrene	0.400	0.464	-16.0	106	-0.01
28	I	Perylene-d12	5.000	5.000	0.0	83	0.00
29		Benzo(b)fluoranthene	0.400	0.478	-19.5	110	0.00
30		Benzo(k)fluoranthene	0.400	0.448	-12.0	101	-0.01
31	C	Benzo(a)pyrene	0.400	0.439	-9.7	104	0.00
32		Dibenzo(a,h)anthracene	0.400	0.452	-13.0	106	0.00
33		Benzo(g,h,i)perylene	0.400	0.462	-15.5	106	0.00

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