Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081019\
Data File : BN007052.D
Acq On : 10 Aug 2019 16:22
Operator : HP/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBE081019

Quant Time: Aug 10 21:57:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 10 21:44:50 2019
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	0.400	0.400	0.0	104	0.00
2		n-Nitrosodimethylamine	0.400	0.388	3.0	129	0.00
3	S	2-Fluorophenol	0.400	0.377	5.8	103	0.00
4	S	Phenol-d6	0.400	0.361	9.8	101	0.00
5		bis(2-Chloroethyl)ether	0.400	0.387	3.3	106	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	106	0.00
7	S	Nitrobenzene-d5	0.400	0.364	9.0	105	0.00
8		Naphthalene	0.400	0.373	6.8	106	0.00
9		Hexachlorobutadiene	0.400	0.387	3.3	108	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.363	9.3	104	0.00
11		2-Methylnaphthalene	0.400	0.364	9.0	104	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	104	0.00
13	S	2,4,6-Tribromophenol	0.400	0.278	30.5#	99	0.00
14	S	2-Fluorobiphenyl	0.400	0.419	-4.7	99	0.00
15		Acenaphthylene	0.400	0.368	8.0	102	0.00
16		Acenaphthene	0.400	0.375	6.3	104	0.00
17		Fluorene	0.400	0.330	17.5	103	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
19		4-Bromophenyl-phenylether	0.400	0.393	1.8	103	0.00
20		Hexachlorobenzene	0.400	0.405	-1.3	106	0.00
21		Pentachlorophenol	0.400	0.316	21.0	94	0.00
22		Phenanthrene	0.400	0.392	2.0	105	0.00
23		Anthracene	0.400	0.369	7.8	102	0.00
24	SURR	Fluoranthene-d10	0.400	0.378	5.5	102	0.00
25		Fluoranthene	0.400	0.382	4.5	102	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	105	-0.01
27		Pyrene	0.400	0.378	5.5	103	0.00
28	S	Terphenyl-d14	0.400	0.383	4.3	105	0.00
29		Benzo(a)anthracene	0.400	0.381	4.8	103	0.00
30		Chrysene	0.400	0.396	1.0	109	-0.01
31		Indeno(1,2,3-cd)pyrene	0.400	0.354	11.5	99	-0.01
32	I	Perylene-d12	0.400	0.400	0.0	99	0.00
33		Benzo(b)fluoranthene	0.400	0.397	0.8	100	0.00
34		Benzo(k)fluoranthene	0.400	0.408	-2.0	101	0.00
35	C	Benzo(a)pyrene	0.400	0.380	5.0	98	-0.01
36		Dibenzo(a,h)anthracene	0.400	0.385	3.8	99	-0.01
37		Benzo(g,h,i)perylene	0.400	0.391	2.3	98	0.00

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