Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE010820\
Data File : BE101008.D
Acq On : 8 Jan 2020 18:41
Operator : JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE010820

Quant Time: Jan 09 10:52:04 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 08 18:36:25 2020
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	-0.01
2		1,4-Dioxane	0.400	0.371	7.3	101	0.00
3		n-Nitrosodimethylamine	0.400	0.422	-5.5	127	-0.01
4	S	2-Fluorophenol	0.400	0.393	1.8	103	0.00
5	S	Phenol-d6	0.400	0.390	2.5	114	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.366	8.5	104	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	105	0.00
8	S	Nitrobenzene-d5	0.400	0.374	6.5	110	0.00
9		Naphthalene	0.400	0.411	-2.7	106	0.00
10		Hexachlorobutadiene	0.400	0.418	-4.5	106	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.408	-2.0	107	0.00
12		2-Methylnaphthalene	0.400	0.404	-1.0	108	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14	S	2,4,6-Tribromophenol	0.400	0.369	7.8	112	0.00
15	S	2-Fluorobiphenyl	0.400	0.411	-2.7	106	-0.01
16		Acenaphthylene	0.400	0.370	7.5	107	0.00
17		Acenaphthene	0.400	0.406	-1.5	109	0.00
18		Fluorene	0.400	0.408	-2.0	109	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
20		4-Bromophenyl-phenylether	0.400	0.389	2.8	112	0.00
21		Hexachlorobenzene	0.400	0.396	1.0	107	0.00
22		Pentachlorophenol	0.400	0.474	-18.5	114	0.00
23		Phenanthrene	0.400	0.400	0.0	111	0.00
24		Anthracene	0.400	0.385	3.8	111	0.00
25	SURR	Fluoranthene-d10	0.400	0.375	6.3	106	0.00
26		Fluoranthene	0.400	0.378	5.5	110	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	103	0.00
28		Pyrene	0.400	0.424	-6.0	106	0.00
29	S	Terphenyl-d14	0.400	0.421	-5.2	107	0.00
30		Benzo(a)anthracene	0.400	0.347	13.3	93	0.00
31		Chrysene	0.400	0.436	-9.0	109	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.358	10.5	106	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	103	0.00
34	I	Perylene-d12	0.400	0.400	0.0	106	0.00
35		Benzo(b)fluoranthene	0.400	0.372	7.0	105	0.00
36		Benzo(k)fluoranthene	0.400	0.367	8.3	97	0.00
37	C	Benzo(a)pyrene	0.400	0.374	6.5	102	0.00
38		Dibenzo(a,h)anthracene	0.400	0.378	5.5	108	0.00
39		Benzo(g,h,i)perylene	0.400	0.380	5.0	98	-0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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