Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031020\
Data File : BN010172.D
Acq On : 10 Mar 2020 22:28
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 11 04:09:51 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN021920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 19 15:57:43 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	74	0.00
2		1,4-Dioxane	0.400	0.538	-34.5#	100	0.00
3		n-Nitrosodimethylamine	0.400	0.577	-44.2#	121	0.00
4	S	2-Fluorophenol	0.400	0.485	-21.2	93	0.00
5	S	Phenol-d6	0.400	0.431	-7.7	84	0.02
6		bis(2-Chloroethyl)ether	0.400	0.424	-6.0	82	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	78	0.01
8	S	Nitrobenzene-d5	0.400	0.398	0.5	83	0.03
9		Naphthalene	0.400	0.414	-3.5	83	0.01
10		Hexachlorobutadiene	0.400	0.443	-10.7	86	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.411	-2.7	83	0.01
12		2-Methylnaphthalene	0.400	0.408	-2.0	83	0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	77	0.00
14	S	2,4,6-Tribromophenol	0.400	0.432	-8.0	93	0.01
15	S	2-Fluorobiphenyl	0.400	0.396	1.0	79	0.01
16		Acenaphthylene	0.400	0.397	0.8	85	0.00
17		Acenaphthene	0.400	0.404	-1.0	81	0.00
18		Fluorene	0.400	0.383	4.3	76	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	70	0.01
20		4-Bromophenyl-phenylether	0.400	0.202	49.5#	46	0.00
21		Hexachlorobenzene	0.400	0.422	-5.5	78	0.01
22		Pentachlorophenol	0.400	0.475	-18.7	91	0.01
23		Phenanthrene	0.400	0.385	3.8	73	0.01
24		Anthracene	0.400	0.388	3.0	76	0.01
25	SURR	Fluoranthene-d10	0.400	0.404	-1.0	76	0.00
26		Fluoranthene	0.400	0.396	1.0	75	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	78	0.00
28		Pyrene	0.400	0.377	5.8	76	0.00
29	S	Terphenyl-d14	0.400	0.378	5.5	76	0.00
30		Benzo(a)anthracene	0.400	0.360	10.0	79	0.00
31		Chrysene	0.400	0.407	-1.7	82	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.500	-25.0	109	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.410	-2.5	80	0.00
34	I	Perylene-d12	0.400	0.400	0.0	79	0.00
35		Benzo(b)fluoranthene	0.400	0.367	8.3	80	0.00
36		Benzo(k)fluoranthene	0.400	0.398	0.5	83	0.00
37	C	Benzo(a)pyrene	0.400	0.388	3.0	79	0.00
38		Dibenzo(a,h)anthracene	0.400	0.380	5.0	80	0.02
39		Benzo(g,h,i)perylene	0.400	0.399	0.3	84	0.00

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

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