Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031720\
Data File : BN010263.D
Acq On : 18 Mar 2020 13:02
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 18 15:23:08 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 16 02:17:23 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	146	0.00
2		1,4-Dioxane	0.400	0.390	2.5	131	0.00
3		n-Nitrosodimethylamine	0.400	0.329	17.8	134	0.00
4	S	2-Fluorophenol	0.400	0.336	16.0	116	0.00
5	S	Phenol-d6	0.400	0.347	13.3	117	0.00
6		bis(2-Chloroethyl)ether	0.400	0.397	0.8	147	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	140	-0.01
8	S	Nitrobenzene-d5	0.400	0.401	-0.3	141	-0.01
9		Naphthalene	0.400	0.401	-0.3	139	-0.01
10		Hexachlorobutadiene	0.400	0.425	-6.2	149	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.383	4.3	136	-0.01
12		2-Methylnaphthalene	0.400	0.384	4.0	134	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	129	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.370	7.5	121	-0.02
15	S	2-Fluorobiphenyl	0.400	0.435	-8.7	138	0.00
16		Acenaphthylene	0.400	0.333	16.8	108	-0.02
17		Acenaphthene	0.400	0.402	-0.5	130	0.00
18		Fluorene	0.400	0.381	4.8	123	-0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	130	-0.01
20		4-Bromophenyl-phenylether	0.400	0.416	-4.0	135	-0.01
21		Hexachlorobenzene	0.400	0.448	-12.0	145	-0.01
22		Pentachlorophenol	0.400	0.374	6.5	133	-0.01
23		Phenanthrene	0.400	0.405	-1.3	129	-0.01
24		Anthracene	0.400	0.339	15.3	108	-0.01
25	SURR	Fluoranthene-d10	0.400	0.368	8.0	121	-0.01
26		Fluoranthene	0.400	0.387	3.3	124	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	122	0.00
28		Pyrene	0.400	0.389	2.8	119	-0.01
29	S	Terphenyl-d14	0.400	0.399	0.3	120	-0.02
30		Benzo(a)anthracene	0.400	0.364	9.0	113	-0.02
31		Chrysene	0.400	0.412	-3.0	127	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.214	46.5#	70	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	125	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	122	-0.02
35		Benzo(b)fluoranthene	0.400	0.388	3.0	118	-0.02
36		Benzo(k)fluoranthene	0.400	0.418	-4.5	127	-0.02
37	C	Benzo(a)pyrene	0.400	0.358	10.5	109	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.417	-4.2	127	-0.03
39		Benzo(g,h,i)perylene	0.400	0.397	0.8	120	-0.03

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

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