Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032720\
Data File : BN010349.D
Acq On : 27 Mar 2020 13:27
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 27 14:16:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 19 15:26:45 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	160	0.00
2		1,4-Dioxane	0.400	0.383	4.3	149	0.00
3		n-Nitrosodimethylamine	0.400	0.289	27.8#	123	0.00
4	S	2-Fluorophenol	0.400	0.442	-10.5	179	0.00
5	S	Phenol-d6	0.400	0.433	-8.2	178	0.00
6		bis(2-Chloroethyl)ether	0.400	0.421	-5.2	166	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	161	-0.01
8	S	Nitrobenzene-d5	0.400	0.437	-9.2	184	0.00
9		Naphthalene	0.400	0.410	-2.5	164	-0.01
10		Hexachlorobutadiene	0.400	0.418	-4.5	169	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.451	-12.7	185	0.00
12		2-Methylnaphthalene	0.400	0.400	0.0	164	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	156	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.392	2.0	171	0.00
15	S	2-Fluorobiphenyl	0.400	0.412	-3.0	159	0.00
16		Acenaphthylene	0.400	0.374	6.5	157	0.00
17		Acenaphthene	0.400	0.410	-2.5	161	0.00
18		Fluorene	0.400	0.393	1.8	158	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	154	0.00
20		4-Bromophenyl-phenylether	0.400	0.381	4.8	150	0.00
21		Hexachlorobenzene	0.400	0.414	-3.5	160	0.00
22		Pentachlorophenol	0.400	0.382	4.5	172	0.00
23		Phenanthrene	0.400	0.392	2.0	154	0.00
24		Anthracene	0.400	0.373	6.8	153	0.00
25	SURR	Fluoranthene-d10	0.400	0.430	-7.5	174	0.00
26		Fluoranthene	0.400	0.388	3.0	156	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	143	0.00
28		Pyrene	0.400	0.446	-11.5	161	0.00
29	S	Terphenyl-d14	0.400	0.424	-6.0	151	0.00
30		Benzo(a)anthracene	0.400	0.393	1.8	147	0.00
31		Chrysene	0.400	0.416	-4.0	148	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.366	8.5	146	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.451	-12.7	149	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	141	-0.01
35		Benzo(b)fluoranthene	0.400	0.380	5.0	147	-0.01
36		Benzo(k)fluoranthene	0.400	0.420	-5.0	152	-0.02
37	C	Benzo(a)pyrene	0.400	0.387	3.3	150	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.419	-4.7	147	-0.02
39		Benzo(g,h,i)perylene	0.400	0.397	0.8	148	-0.02

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

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