Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070819\
Data File : BN006748.D
Acq On : 08 Jul 2019 09:10
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 08 10:03:20 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN070219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 05 05:30:28 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	80	-0.02
2		1,4-Dioxane	0.400	0.438	-9.5	88	-0.04
3		n-Nitrosodimethylamine	0.400	0.403	-0.8	87	-0.02
4	S	2-Fluorophenol	0.400	0.362	9.5	75	0.00
5	S	Phenol-d6	0.400	0.382	4.5	80	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.458	-14.5	103	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	86	-0.01
8	S	Nitrobenzene-d5	0.400	0.376	6.0	85	-0.01
9		Naphthalene	0.400	0.433	-8.2	90	-0.03
10		Hexachlorobutadiene	0.400	0.451	-12.7	89	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.447	-11.7	93	-0.02
12		2-Methylnaphthalene	0.400	0.401	-0.3	87	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	96	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.415	-3.7	107	-0.02
15	S	2-Fluorobiphenyl	0.400	0.411	-2.7	96	-0.02
16		Acenaphthylene	0.400	0.409	-2.2	97	-0.02
17		Acenaphthene	0.400	0.443	-10.7	100	-0.02
18		Fluorene	0.400	0.440	-10.0	101	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	104	-0.01
20		4-Bromophenyl-phenylether	0.400	0.413	-3.2	107	-0.02
21		Hexachlorobenzene	0.400	0.452	-13.0	111	-0.01
22		Pentachlorophenol	0.400	0.598	-49.5#	180	-0.04
23		Phenanthrene	0.400	0.414	-3.5	105	-0.01
24		Anthracene	0.400	0.390	2.5	101	-0.01
25	SURR	Fluoranthene-d10	0.400	0.450	-12.5	111	-0.01
26		Fluoranthene	0.400	0.438	-9.5	108	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	112	-0.01
28		Pyrene	0.400	0.400	0.0	107	-0.01
29	S	Terphenyl-d14	0.400	0.428	-7.0	115	-0.01
30		Benzo(a)anthracene	0.400	0.401	-0.3	104	-0.02
31		Chrysene	0.400	0.436	-9.0	117	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.066	83.5#	29	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.451	-12.7	116	-0.04
34	I	Perylene-d12	0.400	0.400	0.0	115	-0.03
35		Benzo(b)fluoranthene	0.400	0.418	-4.5	112	-0.02
36		Benzo(k)fluoranthene	0.400	0.434	-8.5	117	-0.02
37	C	Benzo(a)pyrene	0.400	0.422	-5.5	111	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.419	-4.7	112	-0.04
39		Benzo(g,h,i)perylene	0.400	0.442	-10.5	119	-0.04

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

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