Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032219\
Data File : BN004902.D
Acq On : 23 Mar 2019 01:30
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 23 02:55:48 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 20 16:06:20 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	195	0.00
2		1,4-Dioxane	0.400	0.417	-4.2	201	0.00
3		n-Nitrosodimethylamine	0.400	0.337	15.8	177	0.00
4	S	2-Fluorophenol	0.400	0.406	-1.5	194	0.00
5	S	Phenol-d6	0.400	0.402	-0.5	197	0.00
6		bis(2-Chloroethyl)ether	0.400	0.411	-2.7	205	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	195	-0.01
8	S	Nitrobenzene-d5	0.400	0.402	-0.5	198	-0.01
9		Naphthalene	0.400	0.411	-2.7	198	-0.01
10		Hexachlorobutadiene	0.400	0.392	2.0	189	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.403	-0.8	197	0.00
12		2-Methylnaphthalene	0.400	0.401	-0.3	194	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	184	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.325	18.8	159	-0.01
15	S	2-Fluorobiphenyl	0.400	0.417	-4.2	188	-0.01
16		Acenaphthylene	0.400	0.401	-0.3	193	-0.01
17		Acenaphthene	0.400	0.407	-1.7	185	-0.01
18		Fluorene	0.400	0.395	1.3	178	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	175	-0.01
20		4-Bromophenyl-phenylether	0.400	0.390	2.5	171	-0.01
21		Hexachlorobenzene	0.400	0.383	4.3	165	-0.01
22		Pentachlorophenol	0.400	0.443	-10.7	224	-0.01
23		Phenanthrene	0.400	0.396	1.0	171	-0.01
24		Anthracene	0.400	0.381	4.8	170	-0.01
25	SURR	Fluoranthene-d10	0.400	0.387	3.3	170	-0.01
26		Fluoranthene	0.400	0.375	6.3	163	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	160	-0.01
28		Pyrene	0.400	0.411	-2.7	161	-0.01
29	S	Terphenyl-d14	0.400	0.414	-3.5	163	-0.01
30		Benzo(a)anthracene	0.400	0.394	1.5	160	-0.01
31		Chrysene	0.400	0.400	0.0	155	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.364	9.0	158	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.330	17.5	127	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	152	-0.01
35		Benzo(b)fluoranthene	0.400	0.396	1.0	148	-0.01
36		Benzo(k)fluoranthene	0.400	0.380	5.0	146	-0.01
37	C	Benzo(a)pyrene	0.400	0.386	3.5	147	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.340	15.0	128	-0.01
39		Benzo(g,h,i)perylene	0.400	0.322	19.5	120	-0.02

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

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