Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012420\
Data File : BN009442.D
Acq On : 24 Jan 2020 15:02
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jan 24 15:53:11 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN011420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 14 09:33:55 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	131	-0.02
2		1,4-Dioxane	0.400	0.393	1.8	122	0.00
3		n-Nitrosodimethylamine	0.400	0.420	-5.0	148	-0.02
4	S	2-Fluorophenol	0.400	0.398	0.5	127	-0.02
5	S	Phenol-d6	0.400	0.423	-5.7	136	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.402	-0.5	143	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	143	-0.01
8	S	Nitrobenzene-d5	0.400	0.396	1.0	149	-0.03
9		Naphthalene	0.400	0.389	2.8	138	-0.01
10		Hexachlorobutadiene	0.400	0.397	0.8	139	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	142	-0.02
12		2-Methylnaphthalene	0.400	0.392	2.0	142	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	150	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.364	9.0	146	-0.02
15	S	2-Fluorobiphenyl	0.400	0.415	-3.7	155	-0.01
16		Acenaphthylene	0.400	0.375	6.3	147	-0.01
17		Acenaphthene	0.400	0.386	3.5	144	-0.01
18		Fluorene	0.400	0.377	5.8	140	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	146	-0.02
20		4-Bromophenyl-phenylether	0.400	0.383	4.3	143	-0.01
21		Hexachlorobenzene	0.400	0.383	4.3	138	-0.01
22		Pentachlorophenol	0.400	0.413	-3.2	151	-0.02
23		Phenanthrene	0.400	0.379	5.3	137	-0.01
24		Anthracene	0.400	0.376	6.0	142	-0.02
25	SURR	Fluoranthene-d10	0.400	0.360	10.0	134	-0.01
26		Fluoranthene	0.400	0.365	8.8	136	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	139	-0.01
28		Pyrene	0.400	0.396	1.0	135	-0.01
29	S	Terphenyl-d14	0.400	0.408	-2.0	140	-0.01
30		Benzo(a)anthracene	0.400	0.385	3.8	139	-0.01
31		Chrysene	0.400	0.388	3.0	134	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.404	-1.0	151	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.414	-3.5	140	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	145	-0.01
35		Benzo(b)fluoranthene	0.400	0.354	11.5	134	-0.01
36		Benzo(k)fluoranthene	0.400	0.390	2.5	135	-0.01
37	C	Benzo(a)pyrene	0.400	0.385	3.8	142	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.381	4.8	141	-0.02
39		Benzo(g,h,i)perylene	0.400	0.376	6.0	136	-0.02

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

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