Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082620\
Data File : BN011653.D
Acq On : 26 Aug 2020 19:57
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 27 00:29:49 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 25 03:31:11 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	118	-0.02
2		1,4-Dioxane	0.400	0.439	-9.7	133	-0.02
3		n-Nitrosodimethylamine	0.400	0.451	-12.7	123	0.00
4	S	2-Fluorophenol	0.400	0.397	0.8	115	0.00
5	S	Phenol-d6	0.400	0.383	4.3	107	0.00
6		bis(2-Chloroethyl)ether	0.400	0.419	-4.7	122	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	108	-0.01
8	S	Nitrobenzene-d5	0.400	0.444	-11.0	124	-0.03
9		Naphthalene	0.400	0.401	-0.3	116	-0.01
10		Hexachlorobutadiene	0.400	0.409	-2.2	118	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.378	5.5	112	-0.01
12		2-Methylnaphthalene	0.400	0.381	4.8	110	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	97	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.351	12.3	86	0.00
15	S	2-Fluorobiphenyl	0.400	0.414	-3.5	105	-0.03
16		Acenaphthylene	0.400	0.356	11.0	94	-0.01
17		Acenaphthene	0.400	0.391	2.3	101	-0.01
18		Fluorene	0.400	0.379	5.3	99	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	92	-0.01
20		4,6-Dinitro-2-methylphenol	0.400	0.519	-29.8#	143	-0.01
21		4-Bromophenyl-phenylether	0.400	0.382	4.5	95	-0.01
22		Hexachlorobenzene	0.400	0.412	-3.0	99	-0.01
23		Atrazine	0.400	0.260	35.0#	84	-0.01
24		Pentachlorophenol	0.400	0.403	-0.8	97	-0.01
25		Phenanthrene	0.400	0.398	0.5	97	0.00
26		Anthracene	0.400	0.357	10.8	87	-0.01
27	SURR	Fluoranthene-d10	0.400	0.381	4.8	97	0.00
28		Fluoranthene	0.400	0.390	2.5	95	-0.01
29	I	Chrysene-d12	0.400	0.400	0.0	91	-0.01
30		Pyrene	0.400	0.389	2.8	95	-0.01
31	S	Terphenyl-d14	0.400	0.396	1.0	94	-0.01
32		Benzo(a)anthracene	0.400	0.404	-1.0	97	0.00
33		Chrysene	0.400	0.422	-5.5	106	0.00
34		Bis(2-ethylhexyl)phthalate	0.400	0.252	37.0#	68	-0.01
35		Indeno(1,2,3-cd)pyrene	0.400	0.460	-15.0	111	-0.01
36	I	Perylene-d12	0.400	0.400	0.0	100	0.00
37		Benzo(b)fluoranthene	0.400	0.401	-0.3	108	0.00
38		Benzo(k)fluoranthene	0.400	0.415	-3.7	110	0.00
39	C	Benzo(a)pyrene	0.400	0.395	1.3	105	0.00
40		Dibenzo(a,h)anthracene	0.400	0.416	-4.0	110	-0.01
41		Benzo(g,h,i)perylene	0.400	0.417	-4.2	110	-0.01

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

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