Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011520\
Data File : BE101068.D
Acq On : 15 Jan 2020 21:33
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 16 04:43:56 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 08 18:36:25 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	110	-0.01
2		1,4-Dioxane	0.400	0.377	5.8	108	0.00
3		n-Nitrosodimethylamine	0.400	0.456	-14.0	146	-0.01
4	S	2-Fluorophenol	0.400	0.508	-27.0#	142	-0.01
5	S	Phenol-d6	0.400	0.477	-19.2	148	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.390	2.5	118	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	110	0.00
8	S	Nitrobenzene-d5	0.400	0.431	-7.7	132	0.00
9		Naphthalene	0.400	0.430	-7.5	116	0.00
10		Hexachlorobutadiene	0.400	0.435	-8.7	115	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.366	8.5	100	-0.01
12		2-Methylnaphthalene	0.400	0.405	-1.3	113	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	100	0.00
14	S	2,4,6-Tribromophenol	0.400	0.398	0.5	116	-0.01
15	S	2-Fluorobiphenyl	0.400	0.407	-1.7	101	-0.01
16		Acenaphthylene	0.400	0.448	-12.0	125	-0.01
17		Acenaphthene	0.400	0.429	-7.2	111	-0.01
18		Fluorene	0.400	0.416	-4.0	107	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	95	0.00
20		4-Bromophenyl-phenylether	0.400	0.404	-1.0	101	-0.01
21		Hexachlorobenzene	0.400	0.426	-6.5	101	-0.01
22		Pentachlorophenol	0.400	0.543	-35.8#	130	-0.01
23		Phenanthrene	0.400	0.405	-1.3	98	0.00
24		Anthracene	0.400	0.436	-9.0	110	0.00
25	SURR	Fluoranthene-d10	0.400	0.387	3.3	95	-0.01
26		Fluoranthene	0.400	0.411	-2.7	104	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	104	-0.01
28		Pyrene	0.400	0.414	-3.5	105	0.00
29	S	Terphenyl-d14	0.400	0.375	6.3	96	-0.01
30		Benzo(a)anthracene	0.400	0.440	-10.0	119	0.00
31		Chrysene	0.400	0.454	-13.5	115	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.584	-46.0#	175	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.592	-48.0#	158	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	143	-0.01
35		Benzo(b)fluoranthene	0.400	0.359	10.3	137	-0.01
36		Benzo(k)fluoranthene	0.400	0.333	16.8	119	0.00
37	C	Benzo(a)pyrene	0.400	0.420	-5.0	155	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.445	-11.2	172	-0.01
39		Benzo(g,h,i)perylene	0.400	0.414	-3.5	144	-0.02

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

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