Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE121418\
Data File : BECCC52.D
Acq On : 15 Dec 2018 1:50
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId :

Quant Time: Dec 15 02:46:47 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 10 14:59:55 2018
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	91	-0.01
2		1,4-Dioxane	0.400	0.281	29.8#	65	0.00
3		n-Nitrosodimethylamine	0.400	0.321	19.8	79	0.00
4	S	2-Fluorophenol	0.400	0.484	-21.0	114	0.00
5	S	Phenol-d6	0.400	0.462	-15.5	105	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.331	17.3	78	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	89	-0.01
8	S	Nitrobenzene-d5	0.400	0.408	-2.0	100	0.00
9		Naphthalene	0.400	0.390	2.5	89	0.00
10		Hexachlorobutadiene	0.400	0.407	-1.7	94	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.382	4.5	86	-0.01
12		2-Methylnaphthalene	0.400	0.381	4.8	87	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	88	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.491	-22.7	114	0.00
15	S	2-Fluorobiphenyl	0.400	0.370	7.5	89	0.00
16		Acenaphthylene	0.400	0.402	-0.5	91	0.00
17		Acenaphthene	0.400	0.373	6.8	83	0.00
18		Fluorene	0.400	0.379	5.3	84	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	83	-0.01
20		4-Bromophenyl-phenylether	0.400	0.370	7.5	80	0.00
21		Hexachlorobenzene	0.400	0.382	4.5	81	0.00
22		Pentachlorophenol	0.400	0.644	-61.0#	149	-0.01
23		Phenanthrene	0.400	0.380	5.0	80	0.00
24		Anthracene	0.400	0.368	8.0	77	0.00
25	SURR	Fluoranthene-d10	0.400	0.419	-4.7	84	0.00
26		Fluoranthene	0.400	0.413	-3.2	82	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	99	-0.01
28		Pyrene	0.400	0.350	12.5	85	0.00
29	S	Terphenyl-d14	0.400	0.323	19.3	81	0.00
30		Benzo(a)anthracene	0.400	0.369	7.8	94	-0.01
31		Chrysene	0.400	0.384	4.0	96	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.408	-2.0	102	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	105	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	106	-0.01
35		Benzo(b)fluoranthene	0.400	0.448	-12.0	122	0.04
36		Benzo(k)fluoranthene	0.400	0.394	1.5	108	-0.01
37	C	Benzo(a)pyrene	0.400	0.400	0.0	108	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.386	3.5	105	-0.02
39		Benzo(g,h,i)perylene	0.400	0.390	2.5	106	-0.02

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

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