Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100518\
Data File : BE097757.D
Acq On : 5 Oct 2018 16:25
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 06 04:18:55 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 03 12:17:19 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	60	-0.01
2		1,4-Dioxane	0.400	0.394	1.5	64	0.00
3		n-Nitrosodimethylamine	0.400	0.341	14.8	54	0.00
4	S	2-Fluorophenol	0.400	0.376	6.0	60	-0.01
5	S	Phenol-d6	0.400	0.365	8.8	56	0.00
6		bis(2-Chloroethyl)ether	0.400	0.382	4.5	60	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	63	0.00
8	S	Nitrobenzene-d5	0.400	0.604	-51.0#	101	0.00
9		Naphthalene	0.400	0.385	3.8	62	0.00
10		Hexachlorobutadiene	0.400	0.386	3.5	63	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.380	5.0	61	0.00
12		2-Methylnaphthalene	0.400	0.378	5.5	62	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	64	0.00
14	S	2,4,6-Tribromophenol	0.400	0.432	-8.0	87	-0.01
15	S	2-Fluorobiphenyl	0.400	0.378	5.5	65	-0.02
16		Acenaphthylene	0.400	0.353	11.8	60	0.00
17		Acenaphthene	0.400	0.378	5.5	64	0.00
18		Fluorene	0.400	0.378	5.5	64	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	63	-0.01
20		4-Bromophenyl-phenylether	0.400	0.380	5.0	64	-0.01
21		Hexachlorobenzene	0.400	0.364	9.0	60	0.00
22		Pentachlorophenol	0.400	0.435	-8.7	78	-0.01
23		Phenanthrene	0.400	0.385	3.8	64	0.00
24		Anthracene	0.400	0.358	10.5	62	0.00
25	SURR	Fluoranthene-d10	0.400	0.373	6.8	63	0.00
26		Fluoranthene	0.400	0.370	7.5	62	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	57	-0.01
28		Pyrene	0.400	0.397	0.8	63	0.00
29	S	Terphenyl-d14	0.400	0.394	1.5	63	0.00
30		Benzo(a)anthracene	0.400	0.351	12.3	56	-0.01
31		Chrysene	0.400	0.405	-1.3	63	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.385	3.8	64	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.331	17.3	52	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	55	-0.01
35		Benzo(b)fluoranthene	0.400	0.373	6.8	57	0.00
36		Benzo(k)fluoranthene	0.400	0.401	-0.3	61	0.00
37	C	Benzo(a)pyrene	0.400	0.359	10.3	55	0.00
38		Dibenzo(a,h)anthracene	0.400	0.347	13.3	53	-0.02
39		Benzo(g,h,i)perylene	0.400	0.367	8.3	55	-0.01

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

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