Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110518\
Data File : BE098163.D
Acq On : 5 Nov 2018 16:47
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 05 17:32:33 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE110118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 01 13:00:52 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	93	0.00
2		1,4-Dioxane	0.400	0.416	-4.0	97	0.00
3		n-Nitrosodimethylamine	0.400	0.429	-7.2	107	0.00
4	S	2-Fluorophenol	0.400	0.402	-0.5	92	0.00
5	S	Phenol-d6	0.400	0.401	-0.3	93	0.00
6		bis(2-Chloroethyl)ether	0.400	0.403	-0.8	97	0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	91	0.01
8	S	Nitrobenzene-d5	0.400	0.405	-1.3	95	0.00
9		Naphthalene	0.400	0.414	-3.5	95	0.01
10		Hexachlorobutadiene	0.400	0.415	-3.7	96	0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.416	-4.0	94	0.02
12		2-Methylnaphthalene	0.400	0.409	-2.2	95	0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	85	0.00
14	S	2,4,6-Tribromophenol	0.400	0.387	3.3	97	0.00
15	S	2-Fluorobiphenyl	0.400	0.409	-2.2	96	0.00
16		Acenaphthylene	0.400	0.420	-5.0	96	0.00
17		Acenaphthene	0.400	0.421	-5.2	93	0.00
18		Fluorene	0.400	0.408	-2.0	93	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	88	0.01
20		4-Bromophenyl-phenylether	0.400	0.415	-3.7	94	0.00
21		Hexachlorobenzene	0.400	0.420	-5.0	97	0.00
22		Pentachlorophenol	0.400	0.500	-25.0	122	0.01
23		Phenanthrene	0.400	0.406	-1.5	93	0.00
24		Anthracene	0.400	0.413	-3.2	97	0.00
25	SURR	Fluoranthene-d10	0.400	0.401	-0.3	93	0.00
26		Fluoranthene	0.400	0.403	-0.8	95	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	81	0.01
28		Pyrene	0.400	0.425	-6.2	94	0.00
29	S	Terphenyl-d14	0.400	0.423	-5.7	94	0.00
30		Benzo(a)anthracene	0.400	0.414	-3.5	92	0.00
31		Chrysene	0.400	0.427	-6.7	94	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.288	28.0#	70	0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.400	0.0	88	0.00
34	I	Perylene-d12	0.400	0.400	0.0	83	0.00
35		Benzo(b)fluoranthene	0.400	0.413	-3.2	91	0.00
36		Benzo(k)fluoranthene	0.400	0.414	-3.5	92	0.00
37	C	Benzo(a)pyrene	0.400	0.399	0.3	88	0.00
38		Dibenzo(a,h)anthracene	0.400	0.391	2.3	85	0.00
39		Benzo(g,h,i)perylene	0.400	0.407	-1.7	89	0.00

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

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