Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121918\
Data File : BE098530.D
Acq On : 19 Dec 2018 17:15
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 19 23:44:05 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	97	0.00
2		1,4-Dioxane	0.400	0.323	19.3	80	-0.01
3		n-Nitrosodimethylamine	0.400	0.347	13.3	91	0.00
4	S	2-Fluorophenol	0.400	0.430	-7.5	107	0.01
5	S	Phenol-d6	0.400	0.465	-16.3	112	0.00
6		bis(2-Chloroethyl)ether	0.400	0.352	12.0	88	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	111	0.00
8	S	Nitrobenzene-d5	0.400	0.359	10.3	109	0.01
9		Naphthalene	0.400	0.393	1.8	112	0.01
10		Hexachlorobutadiene	0.400	0.385	3.8	110	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.403	-0.8	113	0.00
12		2-Methylnaphthalene	0.400	0.395	1.3	112	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14	S	2,4,6-Tribromophenol	0.400	0.407	-1.7	113	0.00
15	S	2-Fluorobiphenyl	0.400	0.394	1.5	114	0.00
16		Acenaphthylene	0.400	0.365	8.8	99	0.00
17		Acenaphthene	0.400	0.394	1.5	105	0.00
18		Fluorene	0.400	0.372	7.0	99	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	91	-0.01
20		4-Bromophenyl-phenylether	0.400	0.393	1.8	93	0.00
21		Hexachlorobenzene	0.400	0.415	-3.7	96	0.00
22		Pentachlorophenol	0.400	0.443	-10.7	104	0.00
23		Phenanthrene	0.400	0.390	2.5	90	0.00
24		Anthracene	0.400	0.368	8.0	84	0.00
25	SURR	Fluoranthene-d10	0.400	0.373	6.8	81	0.00
26		Fluoranthene	0.400	0.370	7.5	80	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	95	-0.01
28		Pyrene	0.400	0.361	9.8	84	-0.01
29	S	Terphenyl-d14	0.400	0.358	10.5	86	-0.01
30		Benzo(a)anthracene	0.400	0.368	8.0	89	-0.01
31		Chrysene	0.400	0.388	3.0	93	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.363	9.3	87	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.261	34.8#	65	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	94	-0.01
35		Benzo(b)fluoranthene	0.400	0.401	-0.3	97	-0.01
36		Benzo(k)fluoranthene	0.400	0.456	-14.0	111	-0.02
37	C	Benzo(a)pyrene	0.400	0.401	-0.3	96	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.312	22.0	75	-0.02
39		Benzo(g,h,i)perylene	0.400	0.245	38.8#	59	-0.02

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

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