Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE031819\
Data File : BE099013.D
Acq On : 18 Mar 2019 23:18
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 19 02:49:38 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 15 05:31:33 2019
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	114	0.00
2		1,4-Dioxane	0.400	0.370	7.5	109	0.00
3		n-Nitrosodimethylamine	0.400	0.182	54.5#	51	0.06
4	S	2-Fluorophenol	0.400	0.285	28.8#	78	0.02
5	S	Phenol-d6	0.400	0.353	11.8	100	0.01
6		bis(2-Chloroethyl)ether	0.400	0.317	20.8	93	0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	128	0.00
8	S	Nitrobenzene-d5	0.400	0.431	-7.7	153	0.01
9		Naphthalene	0.400	0.396	1.0	130	-0.01
10		Hexachlorobutadiene	0.400	0.364	9.0	120	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.373	6.8	121	0.00
12		2-Methylnaphthalene	0.400	0.372	7.0	122	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	128	0.00
14	S	2,4,6-Tribromophenol	0.400	0.285	28.8#	92	0.00
15	S	2-Fluorobiphenyl	0.400	0.570	-42.5#	186	0.00
16		Acenaphthylene	0.400	0.390	2.5	131	0.00
17		Acenaphthene	0.400	0.396	1.0	128	-0.01
18		Fluorene	0.400	0.375	6.3	124	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
20		4-Bromophenyl-phenylether	0.400	0.449	-12.2	126	0.00
21		Hexachlorobenzene	0.400	0.437	-9.2	127	0.00
22		Pentachlorophenol	0.400	0.375	6.3	69	0.01
23		Phenanthrene	0.400	0.401	-0.3	113	0.00
24		Anthracene	0.400	0.344	14.0	100	0.00
25	SURR	Fluoranthene-d10	0.400	0.362	9.5	103	0.00
26		Fluoranthene	0.400	0.375	6.3	107	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	99	-0.01
28		Pyrene	0.400	0.433	-8.2	107	0.00
29	S	Terphenyl-d14	0.400	1.207	-201.8#	301	0.00
30		Benzo(a)anthracene	0.400	0.370	7.5	94	0.00
31		Chrysene	0.400	0.388	3.0	98	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.302	24.5	81	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.334	16.5	84	0.00
34	I	Perylene-d12	0.400	0.400	0.0	99	-0.01
35		Benzo(b)fluoranthene	0.400	0.405	-1.3	104	0.00
36		Benzo(k)fluoranthene	0.400	0.447	-11.7	117	-0.01
37	C	Benzo(a)pyrene	0.400	0.422	-5.5	108	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.256	36.0#	65	0.00
39		Benzo(g,h,i)perylene	0.400	0.359	10.3	91	0.00

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

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