Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE071918\
Data File : BE097000.D
Acq On : 19 Jul 2018 12:04
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 19 13:31:13 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 17 15:02:56 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	63	0.00
2		1,4-Dioxane	0.400	0.428	-7.0	70	0.00
3		n-Nitrosodimethylamine	0.400	0.416	-4.0	69	0.00
4	S	2-Fluorophenol	0.400	0.406	-1.5	72	0.00
5	S	Phenol-d6	0.400	0.375	6.3	64	0.00
6		bis(2-Chloroethyl)ether	0.400	0.387	3.3	63	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	55	0.00
8	S	Nitrobenzene-d5	0.400	0.424	-6.0	63	0.00
9		Naphthalene	0.400	0.385	3.8	56	0.00
10		Hexachlorobutadiene	0.400	0.428	-7.0	62	0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.374	6.5	55	0.00
12		2-Methylnaphthalene	0.400	0.381	4.8	55	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	65	0.00
14	S	2,4,6-Tribromophenol	0.400	0.463	-15.8	81	0.00
15	S	2-Fluorobiphenyl	0.400	0.355	11.3	59	0.01
16		Acenaphthylene	0.400	0.363	9.3	61	0.01
17		Acenaphthene	0.400	0.352	12.0	60	0.00
18		Fluorene	0.400	0.403	-0.8	69	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	79	0.01
20		4-Bromophenyl-phenylether	0.400	0.372	7.0	78	0.01
21		Hexachlorobenzene	0.400	0.368	8.0	75	0.00
22		Pentachlorophenol	0.400	0.582	-45.5#	122	0.01
23		Phenanthrene	0.400	0.381	4.8	79	0.00
24		Anthracene	0.400	0.373	6.8	79	0.00
25	SURR	Fluoranthene-d10	0.400	0.425	-6.2	91	0.00
26		Fluoranthene	0.400	0.401	-0.3	84	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28		Pyrene	0.400	0.341	14.8	88	0.00
29	S	Terphenyl-d14	0.400	0.368	8.0	96	0.00
30		Benzo(a)anthracene	0.400	0.386	3.5	105	0.00
31		Chrysene	0.400	0.364	9.0	93	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.372	7.0	106	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.497	-24.2	139	0.00
34	I	Perylene-d12	0.400	0.400	0.0	108	0.00
35		Benzo(b)fluoranthene	0.400	0.358	10.5	104	0.00
36		Benzo(k)fluoranthene	0.400	0.344	14.0	97	0.00
37	C	Benzo(a)pyrene	0.400	0.390	2.5	110	0.00
38		Dibenzo(a,h)anthracene	0.400	0.477	-19.2	148	0.00
39		Benzo(g,h,i)perylene	0.400	0.454	-13.5	136	0.00

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

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