Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
Data File : BE100276.D
Acq On : 1 Jul 2019 21:30
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 02 02:07:11 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 14:06:53 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	252#	0.00
2		1,4-Dioxane	0.621	0.401	35.4#	157#	0.01
3		n-Nitrosodimethylamine	0.289	0.393	-36.0#	416#	-0.02
4	S	2-Fluorophenol	1.020	0.952	6.7	234#	0.00
5	S	Phenol-d6	1.342	1.766	-31.6#	333#	0.00
6		bis(2-Chloroethyl)ether	1.102	1.642	-49.0#	339#	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	343#	-0.01
8	S	Nitrobenzene-d5	0.397	0.424	-6.8	372#	-0.03
9		Naphthalene	4.411	4.476	-1.5	342#	0.00
10		Hexachlorobutadiene	0.425	0.432	-1.6	338#	0.00
11	SURR	2-Methylnaphthalene-d10	0.752	0.822	-9.3	369#	-0.01
12		2-Methylnaphthalene	0.774	0.804	-3.9	373#	-0.01
13	I	Acenaphthene-d10	1.000	1.000	0.0	365#	0.00
14	S	2,4,6-Tribromophenol	0.280	0.309	-10.4	409#	-0.01
15	S	2-Fluorobiphenyl	1.559	1.610	-3.3	366#	-0.01
16		Acenaphthylene	1.920	2.000	-4.2	367#	0.00
17		Acenaphthene	1.088	1.116	-2.6	354#	0.00
18		Fluorene	1.609	1.618	-0.6	354#	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	318#	0.00
20		4-Bromophenyl-phenylether	0.255	0.280	-9.8	335#	-0.01
21		Hexachlorobenzene	0.265	0.279	-5.3	321#	0.01
22		Pentachlorophenol	0.080	0.101	-26.3#	452#	-0.04
23		Phenanthrene	0.969	1.025	-5.8	335#	-0.01
24		Anthracene	0.873	0.971	-11.2	358#	-0.01
25	SURR	Fluoranthene-d10	5.805	5.711	1.6	309#	0.00
26		Fluoranthene	1.544	1.510	2.2	303#	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	261#	0.00
28		Pyrene	1.059	1.278	-20.7	308#	0.00
29	S	Terphenyl-d14	0.829	1.009	-21.7	314#	0.00
30		Benzo(a)anthracene	1.315	1.462	-11.2	284#	0.01
31		Chrysene	1.321	1.439	-8.9	277#	0.00
32		Bis(2-ethylhexyl)phthalate	2.255	2.229	1.2	265#	0.00
33		Indeno(1,2,3-cd)pyrene	1.747	1.749	-0.1	241#	0.01
34	I	Perylene-d12	1.000	1.000	0.0	240#	0.00
35		Benzo(b)fluoranthene	1.084	1.163	-7.3	254#	0.00
36		Benzo(k)fluoranthene	1.211	1.216	-0.4	232#	0.00
37	C	Benzo(a)pyrene	1.087	1.134	-4.3	241#	0.00
38		Dibenzo(a,h)anthracene	1.135	1.202	-5.9	252#	0.01
39		Benzo(g,h,i)perylene	1.166	1.222	-4.8	242#	0.01

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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