Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100841.D
Acq On : 11 Dec 2019 10:20
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 11 11:44:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 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	117	-0.01
2		1,4-Dioxane	0.729	0.642	11.9	103	-0.01
3		n-Nitrosodimethylamine	0.594	0.579	2.5	130	-0.01
4	S	2-Fluorophenol	0.949	1.076	-13.4	141	-0.01
5	S	Phenol-d6	1.433	1.718	-19.9	151#	0.00
6		bis(2-Chloroethyl)ether	1.385	1.372	0.9	122	-0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	131	0.00
8	S	Nitrobenzene-d5	0.190	0.246	-29.5#	194#	0.00
9		Naphthalene	4.250	4.059	4.5	130	0.00
10		Hexachlorobutadiene	0.166	0.156	6.0	127	0.00
11	SURR	2-Methylnaphthalene-d10	0.580	0.568	2.1	136	0.00
12		2-Methylnaphthalene	0.664	0.655	1.4	137	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	152#	0.00
14	S	2,4,6-Tribromophenol	0.044	0.126	-186.4#	493#	0.00
15	S	2-Fluorobiphenyl	1.587	1.327	16.4	133	-0.01
16		Acenaphthylene	1.655	1.557	5.9	166#	0.00
17		Acenaphthene	1.140	1.109	2.7	163#	0.00
18		Fluorene	1.471	1.789	-21.6	208#	-0.01
19	I	Phenanthrene-d10	1.000	1.000	0.0	241#	-0.01
20		4-Bromophenyl-phenylether	0.196	0.172	12.2	233#	-0.01
21		Hexachlorobenzene	0.213	0.194	8.9	233#	0.00
22		Pentachlorophenol	0.015	0.025	-66.7#	2091#	0.00
23		Phenanthrene	1.212	1.149	5.2	243#	0.00
24		Anthracene	1.013	0.968	4.4	268#	0.00
25	SURR	Fluoranthene-d10	4.728	4.224	10.7	239#	0.00
26		Fluoranthene	1.446	1.314	9.1	243#	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	224#	0.00
28		Pyrene	1.300	1.321	-1.6	243#	0.00
29	S	Terphenyl-d14	0.958	0.940	1.9	230#	0.00
30		Benzo(a)anthracene	1.179	1.242	-5.3	265#	-0.01
31		Chrysene	1.385	1.304	5.8	218#	0.00
32		Bis(2-ethylhexyl)phthalate	1.069	1.379	-29.0#	381#	-0.01
33		Indeno(1,2,3-cd)pyrene	1.207	1.377	-14.1	289#	-0.03
34	I	Perylene-d12	1.000	1.000	0.0	270#	-0.01
35		Benzo(b)fluoranthene	1.095	0.951	13.2	247#	-0.01
36		Benzo(k)fluoranthene	1.228	1.100	10.4	255#	-0.01
37	C	Benzo(a)pyrene	0.948	0.838	11.6	262#	-0.01
38		Dibenzo(a,h)anthracene	0.943	0.897	4.9	289#	-0.03
39		Benzo(g,h,i)perylene	1.009	0.932	7.6	265#	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100841.D
Acq On : 11 Dec 2019 10:20
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 11 11:44:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 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)

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