Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
Data File : BF111528.D
Acq On : 17 Dec 2018 9:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 18 04:17:58 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 14 13:26:17 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	83	0.00
2		1,4-Dioxane	0.580	0.550	5.2	77	0.00
3		Pyridine	1.458	1.452	0.4	80	0.00
4		n-Nitrosodimethylamine	0.662	0.651	1.7	77	0.00
5	S	2-Fluorophenol	1.202	1.117	7.1	77	0.00
6		Aniline	1.983	1.768	10.8	76	0.00
7	S	Phenol-d6	1.599	1.579	1.3	85	0.00
8		2-Chlorophenol	1.436	1.438	-0.1	84	0.00
9		Benzaldehyde	0.970	0.894	7.8	83	0.00
10	C	Phenol	1.781	1.614	9.4	77	0.00
11		bis(2-Chloroethyl)ether	1.396	1.247	10.7	76	0.00
12		1,3-Dichlorobenzene	1.581	1.607	-1.6	87	0.00
13	C	1,4-Dichlorobenzene	1.605	1.635	-1.9	87	0.00
14		1,2-Dichlorobenzene	1.509	1.421	5.8	80	0.00
15		Benzyl Alcohol	1.216	1.118	8.1	78	0.00
16		2,2'-oxybis(1-Chloropropane	2.726	2.638	3.2	82	0.00
17		2-Methylphenol	1.156	1.117	3.4	83	0.00
18		Hexachloroethane	0.597	0.574	3.9	84	0.00
19	P	n-Nitroso-di-n-propylamine	1.055	1.022	3.1	84	0.00
20		3+4-Methylphenols	1.450	1.428	1.5	85	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	67	0.00
22		Acetophenone	0.446	0.571	-28.0#	86	0.00
23	S	Nitrobenzene-d5	0.336	0.417	-24.1	83	0.00
24		Nitrobenzene	0.343	0.376	-9.6	74	0.00
25		Isophorone	0.568	0.696	-22.5	84	0.00
26	C	2-Nitrophenol	0.178	0.193	-8.4	72	0.00
27		2,4-Dimethylphenol	0.258	0.272	-5.4	71	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.395	-0.8	69	0.00
29	C	2,4-Dichlorophenol	0.275	0.291	-5.8	72	0.00
30		1,2,4-Trichlorobenzene	0.288	0.296	-2.8	72	0.00
31		Naphthalene	0.935	0.958	-2.5	71	0.00
32		Benzoic acid	0.223	0.196	12.1	58	0.00
33		4-Chloroaniline	0.416	0.420	-1.0	71	0.00
34	C	Hexachlorobutadiene	0.159	0.172	-8.2	73	0.00
35		Caprolactam	0.102	0.104	-2.0	70	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.310	-2.3	72	0.00
37		2-Methylnaphthalene	0.633	0.630	0.5	72	0.00
38		1-Methylnaphthalene	0.612	0.606	1.0	72	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	69	0.00
40		1,2,4,5-Tetrachlorobenzene	0.548	0.566	-3.3	74	0.00
41	P	Hexachlorocyclopentadiene	0.282	0.278	1.4	67	0.00
42	S	2,4,6-Tribromophenol	0.179	0.268	-49.7#	106	0.00
43	C	2,4,6-Trichlorophenol	0.388	0.390	-0.5	72	0.00
44		2,4,5-Trichlorophenol	0.413	0.425	-2.9	73	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
Data File : BF111528.D
Acq On : 17 Dec 2018 9:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 18 04:17:58 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 14 13:26:17 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.295	1.349	-4.2	76	0.00
46		1,1'-Biphenyl	1.695	1.733	-2.2	74	0.00
47		2-Chloronaphthalene	1.288	1.335	-3.6	75	0.00
48		2-Nitroaniline	0.418	0.419	-0.2	72	0.00
49		Acenaphthylene	1.912	1.990	-4.1	75	0.00
50		Dimethylphthalate	1.440	1.484	-3.1	74	0.00
51		2,6-Dinitrotoluene	0.324	0.338	-4.3	75	0.00
52	C	Acenaphthene	1.208	1.273	-5.4	76	0.00
53		3-Nitroaniline	0.394	0.410	-4.1	75	0.00
54	P	2,4-Dinitrophenol	0.123	0.133	-8.1	77	0.00
55		Dibenzofuran	1.754	1.881	-7.2	78	0.00
56	P	4-Nitrophenol	0.273	0.289	-5.9	72	0.00
57		2,4-Dinitrotoluene	0.426	0.460	-8.0	76	0.00
58		Fluorene	1.272	1.839	-44.6#	106	0.00
59		2,3,4,6-Tetrachlorophenol	0.323	0.390	-20.7	87	0.00
60		Diethylphthalate	1.459	1.721	-18.0	85	0.00
61		4-Chlorophenyl-phenylether	0.623	0.869	-39.5#	100	0.00
62		4-Nitroaniline	0.390	0.592	-51.8#	107	0.00
63		Azobenzene	1.438	2.113	-46.9#	107	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
65		4,6-Dinitro-2-methylphenol	0.113	0.113	0.0	106	0.00
66	c	n-Nitrosodiphenylamine	0.690	0.690	0.0	108	0.00
67		4-Bromophenyl-phenylether	0.216	0.215	0.5	106	0.00
68		Hexachlorobenzene	0.222	0.222	0.0	107	0.00
69		Atrazine	0.199	0.199	0.0	104	0.00
70	C	Pentachlorophenol	0.136	0.144	-5.9	110	0.00
71		Phenanthrene	1.031	1.022	0.9	107	0.00
72		Anthracene	1.054	1.040	1.3	106	0.00
73		Carbazole	1.090	1.097	-0.6	108	0.00
74		Di-n-butylphthalate	1.214	1.247	-2.7	108	0.00
75	C	Fluoranthene	1.008	1.043	-3.5	108	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	108	0.00
77		Benzidine	0.735	0.732	0.4	118	0.00
78		Pyrene	1.410	1.394	1.1	109	0.00
79	S	Terphenyl-d14	0.852	0.804	5.6	106	0.00
80		Butylbenzylphthalate	0.687	0.700	-1.9	113	0.00
81		Benzo(a)anthracene	1.087	1.067	1.8	110	0.00
82		3,3'-Dichlorobenzidine	0.442	0.439	0.7	112	0.00
83		Chrysene	1.146	1.119	2.4	112	0.00
84		Bis(2-ethylhexyl)phthalate	0.878	0.889	-1.3	113	0.00
85	c	Di-n-octyl phthalate	1.481	1.519	-2.6	114	0.00
86		Indeno(1,2,3-cd)pyrene	0.827	0.712	13.9	112	0.00
87	I	Perylene-d12	1.000	1.000	0.0	110	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
Data File : BF111528.D
Acq On : 17 Dec 2018 9:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 18 04:17:58 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 14 13:26:17 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.166	1.188	-1.9	110	0.00
89	Benzo(k)fluoranthene	1.114	1.184	-6.3	124	0.00
90 C	Benzo(a)pyrene	1.051	1.069	-1.7	116	0.00
91	Dibenzo(a,h)anthracene	0.829	0.757	8.7	111	0.00
92	Benzo(g,h,i)perylene	0.814	0.745	8.5	114	0.00

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