Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011216\
Data File : BF084412.D
Acq On : 12 Jan 2016 13:12
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 13 00:47:39 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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	88	0.01
2		1,4-Dioxane	0.586	0.619	-5.6	89	0.01
3		Pyridine	1.633	1.861	-14.0	93	0.00
4		n-Nitrosodimethylamine	0.838	0.743	11.3	74	0.01
5	S	2-Fluorophenol	1.236	1.248	-1.0	95	0.01
6		Aniline	2.272	2.144	5.6	81	0.01
7	S	Phenol-d6	1.553	1.484	4.4	84	0.01
8		2-Chlorophenol	1.403	1.408	-0.4	90	0.01
9		Benzaldehyde	1.011	0.897	11.3	79	0.00
10	C	Phenol	1.766	1.526	13.6	71	0.00
11		bis(2-Chloroethyl)ether	1.368	1.365	0.2	92	0.00
12		1,3-Dichlorobenzene	1.447	1.490	-3.0	93	0.01
13	C	1,4-Dichlorobenzene	1.451	1.487	-2.5	86	0.01
14		1,2-Dichlorobenzene	1.390	1.300	6.5	88	0.01
15		Benzyl Alcohol	1.306	1.098	15.9	71	0.01
16		2,2'-oxybis(1-Chloropropane	2.491	2.196	11.8	80	0.01
17		2-Methylphenol	1.176	1.187	-0.9	83	0.01
18		Hexachloroethane	0.580	0.566	2.4	83	0.01
19	P	n-Nitroso-di-n-propylamine	1.051	0.891	15.2	77	0.01
20		3+4-Methylphenols	1.382	1.215	12.1	83	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	90	0.01
22		Acetophenone	0.481	0.454	5.6	79	0.01
23	S	Nitrobenzene-d5	0.359	0.320	10.9	81	0.01
24		Nitrobenzene	0.364	0.377	-3.6	83	0.01
25		Isophorone	0.687	0.658	4.2	85	0.01
26	C	2-Nitrophenol	0.166	0.163	1.8	89	0.00
27		2,4-Dimethylphenol	0.336	0.292	13.1	73	0.01
28		bis(2-Chloroethoxy)methane	0.420	0.382	9.0	88	0.00
29	C	2,4-Dichlorophenol	0.280	0.288	-2.9	94	0.01
30		1,2,4-Trichlorobenzene	0.289	0.295	-2.1	88	0.01
31		Naphthalene	0.907	0.932	-2.8	93	0.01
32		Benzoic acid	0.198	0.138	30.3#	59	0.01
33		4-Chloroaniline	0.416	0.385	7.5	85	0.00
34	C	Hexachlorobutadiene	0.166	0.150	9.6	81	0.01
35		Caprolactam	0.094	0.094	0.0	79	0.01
36	C	4-Chloro-3-methylphenol	0.313	0.284	9.3	85	0.01
37		2-Methylnaphthalene	0.651	0.565	13.2	81	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	83	0.01
39		1,2,4,5-Tetrachlorobenzene	0.575	0.621	-8.0	90	0.01
40	P	Hexachlorocyclopentadiene	0.347	0.317	8.6	74	0.01
41	S	2,4,6-Tribromophenol	0.202	0.206	-2.0	80	0.01
42	C	2,4,6-Trichlorophenol	0.430	0.426	0.9	85	0.01
43		2,4,5-Trichlorophenol	0.412	0.429	-4.1	84	0.01
44	S	2-Fluorobiphenyl	1.317	1.084	17.7	77	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011216\
Data File : BF084412.D
Acq On : 12 Jan 2016 13:12
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 13 00:47:39 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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		1,1'-Biphenyl	1.590	1.591	-0.1	89	0.01
46		2-Chloronaphthalene	1.249	1.181	5.4	87	0.01
47		2-Nitroaniline	0.412	0.463	-12.4	97	0.01
48		Acenaphthylene	1.845	1.803	2.3	78	0.01
49		Dimethylphthalate	1.430	1.504	-5.2	84	0.01
50		2,6-Dinitrotoluene	0.314	0.324	-3.2	78	0.01
51	C	Acenaphthene	1.237	1.201	2.9	75	0.01
52		3-Nitroaniline	0.389	0.432	-11.1	85	0.01
53	P	2,4-Dinitrophenol	0.093	0.153	-64.5#	96	0.01
54		Dibenzofuran	1.812	1.648	9.1	79	0.01
55	P	4-Nitrophenol	0.282	0.323	-14.5	79	0.01
56		2,4-Dinitrotoluene	0.397	0.396	0.3	71	0.01
57		Fluorene	1.400	1.176	16.0	74	0.01
58		2,3,4,6-Tetrachlorophenol	0.352	0.348	1.1	78	0.01
59		Diethylphthalate	1.410	1.423	-0.9	82	0.01
60		4-Chlorophenyl-phenylether	0.564	0.613	-8.7	91	0.01
61		4-Nitroaniline	0.346	0.409	-18.2	101	0.02
62		Azobenzene	1.546	1.618	-4.7	86	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.01
64		4,6-Dinitro-2-methylphenol	0.109	0.131	-20.2	98	0.01
65	c	n-Nitrosodiphenylamine	0.639	0.621	2.8	82	0.01
66		4-Bromophenyl-phenylether	0.215	0.229	-6.5	85	0.01
67		Hexachlorobenzene	0.242	0.217	10.3	81	0.01
68		Atrazine	0.209	0.228	-9.1	83	0.01
69	C	Pentachlorophenol	0.126	0.105	16.7	63	0.01
70		Phenanthrene	1.046	1.053	-0.7	79	0.01
71		Anthracene	1.026	1.012	1.4	88	0.01
72		Carbazole	1.003	1.022	-1.9	83	0.01
73		Di-n-butylphthalate	1.111	1.132	-1.9	85	0.01
74	C	Fluoranthene	1.093	0.952	12.9	77	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	80	0.01
76		Benzidine	0.717	0.604	15.8	67	0.01
77		Pyrene	1.569	1.363	13.1	79	0.01
78	S	Terphenyl-d14	0.896	0.726	19.0	76	0.01
79		Butylbenzylphthalate	0.679	0.644	5.2	75	0.01
80		Benzo(a)anthracene	1.174	1.037	11.7	75	0.01
81		3,3'-Dichlorobenzidine	0.410	0.375	8.5	72	0.01
82		Chrysene	1.128	0.946	16.1	68	0.01
83		Bis(2-ethylhexyl)phthalate	0.858	0.724	15.6	71	0.01
84	c	Di-n-octyl phthalate	1.449	1.169	19.3	62	0.01
85		Indeno(1,2,3-cd)pyrene	0.895	0.971	-8.5	89	0.03
86	I	Perylene-d12	1.000	1.000	0.0	80	0.01
87		Benzo(b)fluoranthene	1.308	1.176	10.1	69	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011216\
Data File : BF084412.D
Acq On : 12 Jan 2016 13:12
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 13 00:47:39 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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(k)fluoranthene	1.070	1.093	-2.1	82	0.01
89 C	Benzo(a)pyrene	1.110	1.099	1.0	76	0.01
90	Dibenzo(a,h)anthracene	0.955	1.003	-5.0	88	0.03
91	Benzo(g,h,i)perylene	0.989	1.075	-8.7	93	0.02

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