Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030216\
Data File : BF085128.D
Acq On : 2 Mar 2016 19:23
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 03 04:17:46 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 13:25:39 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	57	0.00
2		1,4-Dioxane	0.611	0.589	3.6	55	0.00
3		Pyridine	1.567	1.661	-6.0	57	0.00
4		n-Nitrosodimethylamine	0.713	0.652	8.6	50	-0.02
5	S	2-Fluorophenol	1.232	1.176	4.5	59	0.00
6		Aniline	2.292	2.124	7.3	52	0.00
7	S	Phenol-d6	1.615	1.594	1.3	58	0.00
8		2-Chlorophenol	1.420	1.465	-3.2	60	0.00
9		Benzaldehyde	0.842	0.771	8.4	58	0.00
10	C	Phenol	1.807	2.013	-11.4	65	0.00
11		bis(2-Chloroethyl)ether	1.432	1.242	13.3	47#	0.00
12		1,3-Dichlorobenzene	1.525	1.501	1.6	56	0.00
13	C	1,4-Dichlorobenzene	1.555	1.436	7.7	56	0.00
14		1,2-Dichlorobenzene	1.369	1.504	-9.9	64	0.00
15		Benzyl Alcohol	1.177	1.240	-5.4	63	0.00
16		2,2'-oxybis(1-Chloropropane	2.111	2.005	5.0	56	0.00
17		2-Methylphenol	1.190	1.123	5.6	52	0.00
18		Hexachloroethane	0.575	0.564	1.9	57	0.00
19	P	n-Nitroso-di-n-propylamine	1.064	1.055	0.8	54	0.00
20		3+4-Methylphenols	1.422	1.463	-2.9	64	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	55	0.00
22		Acetophenone	0.503	0.532	-5.8	60	0.00
23	S	Nitrobenzene-d5	0.376	0.366	2.7	51	0.00
24		Nitrobenzene	0.390	0.434	-11.3	64	0.00
25		Isophorone	0.719	0.735	-2.2	56	0.00
26	C	2-Nitrophenol	0.174	0.173	0.6	50#	0.00
27		2,4-Dimethylphenol	0.338	0.324	4.1	50	0.00
28		bis(2-Chloroethoxy)methane	0.397	0.448	-12.8	65	0.00
29	C	2,4-Dichlorophenol	0.278	0.273	1.8	52	0.00
30		1,2,4-Trichlorobenzene	0.303	0.317	-4.6	60	0.00
31		Naphthalene	1.011	0.968	4.3	55	0.00
32		Benzoic acid	0.155	0.237	-52.9#	74	0.00
33		4-Chloroaniline	0.400	0.407	-1.7	59	0.00
34	C	Hexachlorobutadiene	0.176	0.188	-6.8	56	0.00
35		Caprolactam	0.087	0.096	-10.3	59	0.00
36	C	4-Chloro-3-methylphenol	0.307	0.337	-9.8	60	0.00
37		2-Methylnaphthalene	0.618	0.655	-6.0	58	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	59	0.00
39		1,2,4,5-Tetrachlorobenzene	0.616	0.593	3.7	59	0.00
40	P	Hexachlorocyclopentadiene	0.351	0.268	23.6	42#	0.00
41	S	2,4,6-Tribromophenol	0.201	0.230	-14.4	63	0.00
42	C	2,4,6-Trichlorophenol	0.415	0.423	-1.9	56	0.00
43		2,4,5-Trichlorophenol	0.397	0.425	-7.1	59	0.00
44	S	2-Fluorobiphenyl	1.358	1.174	13.5	54	0.00

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030216\
Data File : BF085128.D
Acq On : 2 Mar 2016 19:23
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 03 04:17:46 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 13:25:39 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	1.775	1.607	9.5	53	0.00
46		2-Chloronaphthalene	1.295	1.219	5.9	54	0.00
47		2-Nitroaniline	0.406	0.347	14.5	46#	0.00
48		Acenaphthylene	1.924	2.041	-6.1	65	0.00
49		Dimethylphthalate	1.483	1.361	8.2	53	0.00
50		2,6-Dinitrotoluene	0.308	0.329	-6.8	58	0.00
51	C	Acenaphthene	1.208	1.249	-3.4	64	0.00
52		3-Nitroaniline	0.365	0.319	12.6	43#	0.00
53	P	2,4-Dinitrophenol	0.095	0.084	11.6	46#	0.00
54		Dibenzofuran	1.568	1.658	-5.7	64	0.00
55	P	4-Nitrophenol	0.280	0.229	18.2	41#	0.00
56		2,4-Dinitrotoluene	0.414	0.445	-7.5	63	0.00
57		Fluorene	1.233	1.359	-10.2	66	0.00
58		2,3,4,6-Tetrachlorophenol	0.301	0.332	-10.3	57	0.00
59		Diethylphthalate	1.421	1.385	2.5	54	0.00
60		4-Chlorophenyl-phenylether	0.636	0.663	-4.2	66	0.00
61		4-Nitroaniline	0.342	0.355	-3.8	52	0.00
62		Azobenzene	1.436	1.404	2.2	55	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	53	0.00
64		4,6-Dinitro-2-methylphenol	0.100	0.085	15.0	36#	0.00
65	c	n-Nitrosodiphenylamine	0.630	0.611	3.0	51	0.00
66		4-Bromophenyl-phenylether	0.210	0.246	-17.1	65	0.00
67		Hexachlorobenzene	0.237	0.259	-9.3	57	0.00
68		Atrazine	0.192	0.193	-0.5	57	0.00
69	C	Pentachlorophenol	0.139	0.127	8.6	48#	0.00
70		Phenanthrene	0.995	1.045	-5.0	62	0.00
71		Anthracene	1.117	1.069	4.3	53	0.00
72		Carbazole	0.972	0.973	-0.1	54	0.00
73		Di-n-butylphthalate	1.157	1.205	-4.1	54	0.00
74	C	Fluoranthene	1.092	1.019	6.7	52	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	42#	0.00
76		Benzidine	0.605	0.684	-13.1	51	0.00
77		Pyrene	1.365	1.564	-14.6	51	0.00
78	S	Terphenyl-d14	0.853	0.878	-2.9	43#	0.00
79		Butylbenzylphthalate	0.578	0.627	-8.5	44#	0.00
80		Benzo(a)anthracene	1.161	1.188	-2.3	44#	0.00
81		3,3'-Dichlorobenzidine	0.363	0.451	-24.2	51	0.00
82		Chrysene	1.060	1.210	-14.2	51	0.00
83		Bis(2-ethylhexyl)phthalate	0.668	0.804	-20.4	49#	0.00
84	c	Di-n-octyl phthalate	0.991	1.358	-37.0#	53	0.00
85		Indeno(1,2,3-cd)pyrene	0.897	1.437	-60.2#	68	0.01
86	I	Perylene-d12	1.000	1.000	0.0	64	0.01
87		Benzo(b)fluoranthene	1.218	1.049	13.9	54	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030216\
Data File : BF085128.D
Acq On : 2 Mar 2016 19:23
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 03 04:17:46 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 13:25:39 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	1.055	1.037	1.7	62	0.01
89 C	Benzo(a)pyrene	1.024	1.041	-1.7	64	0.00
90	Dibenzo(a,h)anthracene	0.923	1.053	-14.1	69	0.01
91	Benzo(g,h,i)perylene	0.938	0.989	-5.4	64	0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1