Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091515\
Data File : BF081625.D
Acq On : 15 Sep 2015 11:57
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 15 14:04:52 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 14 13:51:56 2015
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	123	0.00
2		1,4-Dioxane	40.000	39.475	1.3	123	0.00
3		Pyridine	40.000	34.811	13.0	93	0.00
4		n-Nitrosodimethylamine	40.000	46.282	-15.7	134	0.00
5	S	2-Fluorophenol	80.000	76.516	4.4	123	0.00
6		Aniline	40.000	38.367	4.1	118	0.00
7	S	Phenol-d6	80.000	79.422	0.7	120	0.00
8		2-Chlorophenol	40.000	39.923	0.2	123	0.00
9		Benzaldehyde	40.000	34.799	13.0	106	0.00
10	C	Phenol	40.000	35.073	12.3	97	0.00
11		bis(2-Chloroethyl)ether	40.000	40.146	-0.4	122	0.00
12		1,3-Dichlorobenzene	40.000	39.010	2.5	122	0.00
13	C	1,4-Dichlorobenzene	40.000	39.013	2.5	120	0.00
14		1,2-Dichlorobenzene	40.000	41.204	-3.0	129	0.00
15		Benzyl Alcohol	40.000	40.585	-1.5	117	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	44.509	-11.3	142	0.00
17		2-Methylphenol	40.000	40.704	-1.8	123	0.00
18		Hexachloroethane	40.000	41.587	-4.0	127	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	43.132	-7.8	138	0.00
20		3+4-Methylphenols	40.000	39.662	0.8	123	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	132	0.00
22		Acetophenone	40.000	38.187	4.5	122	0.00
23	S	Nitrobenzene-d5	80.000	79.854	0.2	129	0.00
24		Nitrobenzene	40.000	39.969	0.1	126	0.00
25		Isophorone	40.000	40.682	-1.7	136	0.00
26	C	2-Nitrophenol	40.000	38.176	4.6	132	0.00
27		2,4-Dimethylphenol	40.000	39.109	2.2	115	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.791	-2.0	120	0.00
29	C	2,4-Dichlorophenol	40.000	39.512	1.2	126	0.00
30		1,2,4-Trichlorobenzene	40.000	40.410	-1.0	127	0.00
31		Naphthalene	40.000	39.115	2.2	129	0.00
32		Benzoic acid	40.000	24.796	38.0#	73	0.00
33		4-Chloroaniline	40.000	39.435	1.4	114	0.00
34	C	Hexachlorobutadiene	40.000	43.162	-7.9	148	0.00
35		Caprolactam	40.000	37.352	6.6	118	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.516	-6.3	140	0.00
37		2-Methylnaphthalene	40.000	39.978	0.1	141	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	135	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.798	0.5	131	0.00
40	P	Hexachlorocyclopentadiene	40.000	37.351	6.6	124	0.00
41	S	2,4,6-Tribromophenol	80.000	80.757	-0.9	130	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.626	0.9	137	0.00
43		2,4,5-Trichlorophenol	40.000	39.711	0.7	131	0.00
44	S	2-Fluorobiphenyl	80.000	80.324	-0.4	130	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091515\
Data File : BF081625.D
Acq On : 15 Sep 2015 11:57
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 15 14:04:52 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 14 13:51:56 2015
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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	38.616	3.5	145	0.00
46		2-Chloronaphthalene	40.000	39.031	2.4	123	0.00
47		2-Nitroaniline	40.000	41.849	-4.6	137	0.00
48		Acenaphthylene	40.000	38.654	3.4	143	0.00
49		Dimethylphthalate	40.000	41.138	-2.8	144	0.00
50		2,6-Dinitrotoluene	40.000	36.877	7.8	118	0.00
51	C	Acenaphthene	40.000	37.923	5.2	143	0.00
52		3-Nitroaniline	40.000	39.769	0.6	132	0.00
53	P	2,4-Dinitrophenol	40.000	33.232	16.9	111	0.00
54		Dibenzofuran	40.000	38.933	2.7	143	0.00
55	P	4-Nitrophenol	40.000	35.029	12.4	101	0.00
56		2,4-Dinitrotoluene	40.000	40.128	-0.3	135	0.00
57		Fluorene	40.000	42.290	-5.7	135	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.743	-4.4	146	0.00
59		Diethylphthalate	40.000	43.597	-9.0	147	0.00
60		4-Chlorophenyl-phenylether	40.000	43.372	-8.4	153	0.00
61		4-Nitroaniline	40.000	36.319	9.2	130	0.00
62		Azobenzene	40.000	42.431	-6.1	135	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	142	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.958	0.1	123	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.029	7.4	131	0.00
66		4-Bromophenyl-phenylether	40.000	40.698	-1.7	162	0.00
67		Hexachlorobenzene	40.000	40.130	-0.3	152	0.00
68		Atrazine	40.000	40.444	-1.1	148	0.00
69	C	Pentachlorophenol	40.000	38.758	3.1	149	0.00
70		Phenanthrene	40.000	38.816	3.0	134	0.00
71		Anthracene	40.000	38.553	3.6	141	0.00
72		Carbazole	40.000	37.204	7.0	141	0.00
73		Di-n-butylphthalate	40.000	45.185	-13.0	170	0.00
74	C	Fluoranthene	40.000	39.262	1.8	149	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	117	0.00
76		Benzidine	40.000	33.239	16.9	111	0.00
77		Pyrene	40.000	45.399	-13.5	137	0.00
78	S	Terphenyl-d14	80.000	90.451	-13.1	160	0.00
79		Butylbenzylphthalate	40.000	46.867	-17.2	150	0.00
80		Benzo(a)anthracene	40.000	42.451	-6.1	128	0.00
81		3,3'-Dichlorobenzidine	40.000	44.379	-10.9	149	0.00
82		Chrysene	40.000	40.700	-1.8	151	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	48.783	-22.0	165	0.00
84	c	Di-n-octyl phthalate	40.000	45.927	-14.8	151	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.116	-0.3	129	0.00
86	I	Perylene-d12	20.000	20.000	0.0	116	0.00
87		Benzo(b)fluoranthene	40.000	42.997	-7.5	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091515\
Data File : BF081625.D
Acq On : 15 Sep 2015 11:57
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 15 14:04:52 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 14 13:51:56 2015
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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	39.724	0.7	139	0.00
89 C	Benzo(a)pyrene	40.000	38.189	4.5	117	0.00
90	Dibenzo(a,h)anthracene	40.000	46.012	-15.0	133	0.00
91	Benzo(g,h,i)perylene	40.000	44.234	-10.6	131	0.00

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