Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_F\DATA\BF021518\
Data File : BF103036.D
Acq On : 16 Feb 2018 00:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 16 04:50:06 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 15 14:08:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	79	0.00
2		1,4-Dioxane	40.000	40.368	-0.9	78	-0.04
3		Pyridine	40.000	40.832	-2.1	78	-0.02
4		n-Nitrosodimethylamine	40.000	37.960	5.1	73	-0.04
5	S	2-Fluorophenol	80.000	79.695	0.4	79	0.00
6		Aniline	40.000	38.907	2.7	77	0.00
7	S	Phenol-d6	80.000	79.308	0.9	79	0.00
8		2-Chlorophenol	40.000	40.355	-0.9	79	0.00
9		Benzaldehyde	40.000	34.819	13.0	68	0.00
10	C	Phenol	40.000	43.317	-8.3	87	0.00
11		bis(2-Chloroethyl)ether	40.000	39.503	1.2	78	0.00
12		1,3-Dichlorobenzene	40.000	38.982	2.5	77	0.00
13	C	1,4-Dichlorobenzene	40.000	39.817	0.5	79	0.00
14		1,2-Dichlorobenzene	40.000	39.395	1.5	78	0.00
15		Benzyl Alcohol	40.000	39.548	1.1	76	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.029	9.9	70	0.00
17		2-Methylphenol	40.000	38.930	2.7	77	0.00
18		Hexachloroethane	40.000	32.718	18.2	63	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.427	8.9	74	0.00
20		3+4-Methylphenols	40.000	37.165	7.1	72	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	79	0.00
22		Acetophenone	40.000	36.394	9.0	74	0.00
23	S	Nitrobenzene-d5	80.000	90.431	-13.0	85	0.00
24		Nitrobenzene	40.000	42.597	-6.5	82	0.00
25		Isophorone	40.000	37.302	6.7	75	0.00
26	C	2-Nitrophenol	40.000	36.015	10.0	74	0.00
27		2,4-Dimethylphenol	40.000	36.893	7.8	72	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.092	2.3	78	0.00
29	C	2,4-Dichlorophenol	40.000	38.902	2.7	74	0.00
30		1,2,4-Trichlorobenzene	40.000	38.338	4.2	77	0.00
31		Naphthalene	40.000	38.137	4.7	76	0.00
32		Benzoic acid	40.000	30.706	23.2	56	-0.04
33		4-Chloroaniline	40.000	38.667	3.3	77	0.00
34	C	Hexachlorobutadiene	40.000	38.056	4.9	75	0.00
35		Caprolactam	40.000	38.631	3.4	74	-0.01
36	C	4-Chloro-3-methylphenol	40.000	38.249	4.4	74	0.00
37		2-Methylnaphthalene	40.000	37.811	5.5	76	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	75	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	41.084	-2.7	78	0.00
40	P	Hexachlorocyclopentadiene	40.000	2.898	92.8#	5	0.00
41	S	2,4,6-Tribromophenol	80.000	72.657	9.2	64	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.557	3.6	69	0.00
43		2,4,5-Trichlorophenol	40.000	37.651	5.9	66	0.00
44	S	2-Fluorobiphenyl	80.000	80.214	-0.3	74	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_F\DATA\BF021518\
Data File : BF103036.D
Acq On : 16 Feb 2018 00:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 16 04:50:06 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 15 14:08:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	39.367	1.6	75	0.00
46		2-Chloronaphthalene	40.000	39.824	0.4	75	0.00
47		2-Nitroaniline	40.000	49.134	-22.8	94	0.00
48		Acenaphthylene	40.000	38.322	4.2	73	0.00
49		Dimethylphthalate	40.000	39.951	0.1	76	0.00
50		2,6-Dinitrotoluene	40.000	39.927	0.2	71	0.00
51	C	Acenaphthene	40.000	36.919	7.7	71	0.00
52		3-Nitroaniline	40.000	49.599	-24.0	89	0.00
53	P	2,4-Dinitrophenol	40.000	11.606	71.0#	7	0.00
54		Dibenzofuran	40.000	38.382	4.0	73	0.00
55	P	4-Nitrophenol	40.000	28.712	28.2#	52	0.00
56		2,4-Dinitrotoluene	40.000	37.001	7.5	69	0.00
57		Fluorene	40.000	39.255	1.9	73	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	32.846	17.9	59	0.00
59		Diethylphthalate	40.000	38.605	3.5	73	0.00
60		4-Chlorophenyl-phenylether	40.000	41.283	-3.2	77	0.00
61		4-Nitroaniline	40.000	48.202	-20.5	88	0.00
62		Azobenzene	40.000	37.267	6.8	71	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	66	0.00
64		4,6-Dinitro-2-methylphenol	40.000	13.749	65.6#	10	0.00
65	c	n-Nitrosodiphenylamine	40.000	43.263	-8.2	72	0.00
66		4-Bromophenyl-phenylether	40.000	42.668	-6.7	71	0.00
67		Hexachlorobenzene	40.000	40.207	-0.5	67	0.00
68		Atrazine	40.000	41.000	-2.5	66	0.00
69	C	Pentachlorophenol	40.000	25.477	36.3#	40	0.00
70		Phenanthrene	40.000	37.932	5.2	64	0.00
71		Anthracene	40.000	38.621	3.4	65	0.00
72		Carbazole	40.000	37.559	6.1	63	0.00
73		Di-n-butylphthalate	40.000	43.469	-8.7	71	0.00
74	C	Fluoranthene	40.000	34.596	13.5	58	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	64	0.00
76		Benzidine	40.000	30.224	24.4	45	0.00
77		Pyrene	40.000	36.257	9.4	58	0.00
78	S	Terphenyl-d14	80.000	75.639	5.5	62	0.00
79		Butylbenzylphthalate	40.000	41.669	-4.2	66	0.00
80		Benzo(a)anthracene	40.000	39.546	1.1	65	0.00
81		3,3'-Dichlorobenzidine	40.000	40.241	-0.6	62	0.00
82		Chrysene	40.000	37.878	5.3	62	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	46.374	-15.9	73	0.00
84	c	Di-n-octyl phthalate	40.000	48.576	-21.4#	75	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	35.108	12.2	61	0.00
86	I	Perylene-d12	20.000	20.000	0.0	65	0.00
87		Benzo(b)fluoranthene	40.000	41.047	-2.6	66	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_F\DATA\BF021518\
Data File : BF103036.D
Acq On : 16 Feb 2018 00:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 16 04:50:06 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 15 14:08:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	38.076	4.8	61	0.00
89 C	Benzo(a)pyrene	40.000	38.743	3.1	63	0.00
90	Dibenzo(a,h)anthracene	40.000	37.191	7.0	61	0.00
91	Benzo(g,h,i)perylene	40.000	36.626	8.4	61	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 2