Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
Data File : BF107606.D
Acq On : 24 Jul 2018 00:16
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 24 05:13:50 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 16:29:27 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	56	0.00
2		1,4-Dioxane	0.579	0.555	4.1	54	-0.05
3		Pyridine	1.575	1.426	9.5	50	-0.03
4		n-Nitrosodimethylamine	0.664	0.531	20.0	46#	-0.04
5	S	2-Fluorophenol	1.244	1.182	5.0	53	0.00
6		Aniline	1.926	1.733	10.0	51	0.00
7	S	Phenol-d6	1.494	1.373	8.1	53	0.00
8		2-Chlorophenol	1.333	1.311	1.7	56	0.00
9		Benzaldehyde	0.977	0.914	6.4	57	0.00
10	C	Phenol	1.757	1.546	12.0	51	0.00
11		bis(2-Chloroethyl)ether	1.456	1.192	18.1	45#	0.00
12		1,3-Dichlorobenzene	1.511	1.444	4.4	55	0.00
13	C	1,4-Dichlorobenzene	1.549	1.543	0.4	57	0.00
14		1,2-Dichlorobenzene	1.390	1.381	0.6	58	0.00
15		Benzyl Alcohol	1.018	0.962	5.5	52	0.00
16		2,2'-oxybis(1-Chloropropane	2.457	2.224	9.5	52	0.00
17		2-Methylphenol	1.133	1.005	11.3	50	0.00
18		Hexachloroethane	0.543	0.354	34.8#	36#	0.00
19	P	n-Nitroso-di-n-propylamine	0.965	0.868	10.1	53	0.00
20		3+4-Methylphenols	1.345	1.211	10.0	52	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	52	0.00
22		Acetophenone	0.470	0.469	0.2	55	0.00
23	S	Nitrobenzene-d5	0.296	0.336	-13.5	57	0.00
24		Nitrobenzene	0.339	0.353	-4.1	51	0.00
25		Isophorone	0.633	0.542	14.4	45#	0.00
26	C	2-Nitrophenol	0.143	0.126	11.9	43#	0.00
27		2,4-Dimethylphenol	0.271	0.238	12.2	46#	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.389	8.0	49#	0.00
29	C	2,4-Dichlorophenol	0.277	0.257	7.2	47#	0.00
30		1,2,4-Trichlorobenzene	0.311	0.291	6.4	49#	0.00
31		Naphthalene	0.879	0.822	6.5	51	0.00
32		Benzoic acid	0.166	0.097	41.6#	28#	-0.05
33		4-Chloroaniline	0.384	0.373	2.9	54	0.00
34	C	Hexachlorobutadiene	0.185	0.185	0.0	53	0.00
35		Caprolactam	0.090	0.074	17.8	41#	-0.02
36	C	4-Chloro-3-methylphenol	0.308	0.259	15.9	43#	0.00
37		2-Methylnaphthalene	0.609	0.536	12.0	48#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	44#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.667	0.700	-4.9	47#	0.00
40	P	Hexachlorocyclopentadiene	0.339	0.013#	96.2#	2#	0.00
41	S	2,4,6-Tribromophenol	0.227	0.242	-6.6	45#	0.00
42	C	2,4,6-Trichlorophenol	0.427	0.415	2.8	41#	0.00
43		2,4,5-Trichlorophenol	0.446	0.451	-1.1	42#	0.00
44	S	2-Fluorobiphenyl	1.315	1.507	-14.6	55	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
Data File : BF107606.D
Acq On : 24 Jul 2018 00:16
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 24 05:13:50 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 16:29:27 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		1,1'-Biphenyl	1.742	1.839	-5.6	48#	0.00
46		2-Chloronaphthalene	1.332	1.345	-1.0	46#	0.00
47		2-Nitroaniline	0.388	0.489	-26.0#	52	0.00
48		Acenaphthylene	2.000	2.022	-1.1	46#	0.00
49		Dimethylphthalate	1.507	1.528	-1.4	46#	0.00
50		2,6-Dinitrotoluene	0.299	0.273	8.7	39#	0.00
51	C	Acenaphthene	1.253	1.192	4.9	42#	0.00
52		3-Nitroaniline	0.364	0.452	-24.2	53	0.00
53	P	2,4-Dinitrophenol	0.096	0.005#	94.8#	2#	0.01
54		Dibenzofuran	1.845	1.822	1.2	45#	0.00
55	P	4-Nitrophenol	0.273	0.220	19.4	33#	0.00
56		2,4-Dinitrotoluene	0.361	0.316	12.5	36#	0.00
57		Fluorene	1.316	1.352	-2.7	46#	0.00
58		2,3,4,6-Tetrachlorophenol	0.371	0.379	-2.2	43#	0.00
59		Diethylphthalate	1.574	1.552	1.4	45#	0.00
60		4-Chlorophenyl-phenylether	0.744	0.734	1.3	45#	0.00
61		4-Nitroaniline	0.306	0.484	-58.2#	65	-0.01
62		Azobenzene	1.476	1.344	8.9	40#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	45#	0.00
64		4,6-Dinitro-2-methylphenol	0.084	0.007	91.7#	3#	0.00
65	c	n-Nitrosodiphenylamine	0.679	0.645	5.0	45#	0.00
66		4-Bromophenyl-phenylether	0.235	0.214	8.9	42#	0.00
67		Hexachlorobenzene	0.259	0.243	6.2	44#	0.00
68		Atrazine	0.207	0.182	12.1	40#	0.00
69	C	Pentachlorophenol	0.162	0.126	22.2#	34#	0.00
70		Phenanthrene	0.975	0.960	1.5	47#	0.00
71		Anthracene	0.981	1.019	-3.9	50	0.00
72		Carbazole	1.020	1.082	-6.1	51	0.00
73		Di-n-butylphthalate	1.155	1.226	-6.1	51	0.00
74	C	Fluoranthene	1.046	1.141	-9.1	52	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	46#	0.00
76		Benzidine	0.576	0.890	-54.5#	75	0.00
77		Pyrene	1.368	1.456	-6.4	52	0.00
78	S	Terphenyl-d14	0.781	0.936	-19.8	59	0.00
79		Butylbenzylphthalate	0.588	0.710	-20.7	53	0.00
80		Benzo(a)anthracene	1.172	1.117	4.7	45#	0.00
81		3,3'-Dichlorobenzidine	0.371	0.467	-25.9#	61	0.00
82		Chrysene	1.126	1.088	3.4	48#	0.00
83		Bis(2-ethylhexyl)phthalate	0.830	0.925	-11.4	53	0.00
84	c	Di-n-octyl phthalate	1.290	1.500	-16.3	52	0.00
85		Indeno(1,2,3-cd)pyrene	0.954	0.852	10.7	46#	0.01
86	I	Perylene-d12	1.000	1.000	0.0	42#	0.00
87		Benzo(b)fluoranthene	1.095	0.988	9.8	38#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
Data File : BF107606.D
Acq On : 24 Jul 2018 00:16
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 24 05:13:50 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 20 16:29:27 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(k)fluoranthene	1.072	1.083	-1.0	44#	0.00
89 C	Benzo(a)pyrene	1.001	0.943	5.8	40#	0.00
90	Dibenzo(a,h)anthracene	0.866	0.897	-3.6	46#	0.01
91	Benzo(g,h,i)perylene	0.854	0.873	-2.2	45#	0.01

(#) = Out of Range SPCC's out = 2 CCC's out = 1