Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF042618\
Data File : BF104857.D
Acq On : 27 Apr 2018 2:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 27 12:54:42 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 25 14:01:50 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	95	0.00
2		1,4-Dioxane	0.609	0.516	15.3	80	-0.05
3		Pyridine	1.714	1.468	14.4	81	-0.05
4		n-Nitrosodimethylamine	0.599	0.492	17.9	77	-0.04
5	S	2-Fluorophenol	1.308	1.185	9.4	87	0.00
6		Aniline	2.233	1.945	12.9	83	0.00
7	S	Phenol-d6	1.607	1.470	8.5	89	0.00
8		2-Chlorophenol	1.381	1.326	4.0	91	0.00
9		Benzaldehyde	0.802	0.742	7.5	85	0.00
10	C	Phenol	1.705	1.604	5.9	91	0.00
11		bis(2-Chloroethyl)ether	1.361	1.258	7.6	88	0.00
12		1,3-Dichlorobenzene	1.564	1.485	5.1	91	0.00
13	C	1,4-Dichlorobenzene	1.559	1.488	4.6	91	0.00
14		1,2-Dichlorobenzene	1.445	1.389	3.9	91	0.00
15		Benzyl Alcohol	1.221	1.077	11.8	85	0.00
16		2,2'-oxybis(1-Chloropropane	1.827	1.435	21.5	75	0.00
17		2-Methylphenol	1.200	1.140	5.0	91	0.00
18		Hexachloroethane	0.556	0.385	30.8#	66	0.00
19	P	n-Nitroso-di-n-propylamine	1.050	0.854	18.7	81	0.00
20		3+4-Methylphenols	1.488	1.376	7.5	89	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.492	0.440	10.6	89	0.00
23	S	Nitrobenzene-d5	0.306	0.320	-4.6	100	0.00
24		Nitrobenzene	0.338	0.338	0.0	94	0.00
25		Isophorone	0.648	0.559	13.7	85	0.00
26	C	2-Nitrophenol	0.138	0.158	-14.5	107	0.00
27		2,4-Dimethylphenol	0.286	0.245	14.3	83	0.00
28		bis(2-Chloroethoxy)methane	0.396	0.347	12.4	87	0.00
29	C	2,4-Dichlorophenol	0.273	0.257	5.9	92	0.00
30		1,2,4-Trichlorobenzene	0.299	0.284	5.0	93	0.00
31		Naphthalene	0.996	0.918	7.8	91	0.00
32		Benzoic acid	0.228	0.183	19.7	74	0.00
33		4-Chloroaniline	0.427	0.400	6.3	93	0.00
34	C	Hexachlorobutadiene	0.164	0.155	5.5	94	0.00
35		Caprolactam	0.095	0.089	6.3	90	0.00
36	C	4-Chloro-3-methylphenol	0.292	0.277	5.1	93	0.00
37		2-Methylnaphthalene	0.644	0.591	8.2	91	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
39		1,2,4,5-Tetrachlorobenzene	0.573	0.580	-1.2	96	0.00
40	P	Hexachlorocyclopentadiene	0.321	0.027#	91.6#	8#	0.00
41	S	2,4,6-Tribromophenol	0.207	0.184	11.1	83	0.00
42	C	2,4,6-Trichlorophenol	0.397	0.384	3.3	88	0.00
43		2,4,5-Trichlorophenol	0.445	0.440	1.1	93	0.00
44	S	2-Fluorobiphenyl	1.266	1.215	4.0	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF042618\
Data File : BF104857.D
Acq On : 27 Apr 2018 2:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 27 12:54:42 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 25 14:01:50 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	1.680	1.607	4.3	91	0.00
46		2-Chloronaphthalene	1.327	1.272	4.1	91	0.00
47		2-Nitroaniline	0.328	0.396	-20.7	105	0.00
48		Acenaphthylene	2.082	1.870	10.2	84	0.00
49		Dimethylphthalate	1.577	1.530	3.0	92	0.00
50		2,6-Dinitrotoluene	0.292	0.319	-9.2	95	0.00
51	C	Acenaphthene	1.270	1.098	13.5	82	0.00
52		3-Nitroaniline	0.342	0.396	-15.8	101	0.00
53	P	2,4-Dinitrophenol	0.089	0.033#	62.9#	36#	0.00
54		Dibenzofuran	1.770	1.527	13.7	81	0.00
55	P	4-Nitrophenol	0.268	0.222	17.2	71	0.00
56		2,4-Dinitrotoluene	0.383	0.373	2.6	87	0.00
57		Fluorene	1.289	1.248	3.2	90	0.00
58		2,3,4,6-Tetrachlorophenol	0.332	0.287	13.6	81	0.00
59		Diethylphthalate	1.571	1.418	9.7	86	0.00
60		4-Chlorophenyl-phenylether	0.574	0.586	-2.1	95	0.00
61		4-Nitroaniline	0.334	0.390	-16.8	100	0.00
62		Azobenzene	1.501	1.228	18.2	77	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	82	0.00
64		4,6-Dinitro-2-methylphenol	0.087	0.047	46.0#	42#	0.00
65	c	n-Nitrosodiphenylamine	0.693	0.689	0.6	83	0.00
66		4-Bromophenyl-phenylether	0.213	0.214	-0.5	85	0.00
67		Hexachlorobenzene	0.239	0.236	1.3	83	0.00
68		Atrazine	0.209	0.190	9.1	76	0.00
69	C	Pentachlorophenol	0.148	0.107	27.7#	57	0.00
70		Phenanthrene	1.114	1.012	9.2	76	0.00
71		Anthracene	1.132	1.026	9.4	76	0.00
72		Carbazole	1.106	0.976	11.8	75	0.00
73		Di-n-butylphthalate	1.351	1.253	7.3	78	0.00
74	C	Fluoranthene	1.164	0.952	18.2	69	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	73	0.00
76		Benzidine	0.692	0.732	-5.8	77	0.00
77		Pyrene	1.482	1.369	7.6	68	0.00
78	S	Terphenyl-d14	0.877	0.818	6.7	71	0.00
79		Butylbenzylphthalate	0.698	0.636	8.9	66	0.00
80		Benzo(a)anthracene	1.195	1.189	0.5	74	0.00
81		3,3'-Dichlorobenzidine	0.445	0.449	-0.9	72	0.00
82		Chrysene	1.227	1.141	7.0	68	0.00
83		Bis(2-ethylhexyl)phthalate	0.898	0.853	5.0	70	0.00
84	c	Di-n-octyl phthalate	1.501	1.342	10.6	64	0.00
85		Indeno(1,2,3-cd)pyrene	1.043	0.944	9.5	69	0.00
86	I	Perylene-d12	1.000	1.000	0.0	69	0.00
87		Benzo(b)fluoranthene	1.263	1.199	5.1	65	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF042618\
Data File : BF104857.D
Acq On : 27 Apr 2018 2:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 27 12:54:42 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 25 14:01:50 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	1.193	1.198	-0.4	70	0.00
89 C	Benzo(a)pyrene	1.130	1.090	3.5	67	0.00
90	Dibenzo(a,h)anthracene	0.928	0.898	3.2	69	0.00
91	Benzo(g,h,i)perylene	0.899	0.851	5.3	70	0.00

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