Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081619\
Data File : BF116325.D
Acq On : 16 Aug 2019 18:45
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 05:30:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 2019
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	80	0.00
2		1,4-Dioxane	0.604	0.610	-1.0	83	0.00
3		Pyridine	1.686	1.552	7.9	75	0.00
4		n-Nitrosodimethylamine	0.665	0.605	9.0	74	0.00
5	S	2-Fluorophenol	1.195	1.304	-9.1	87	0.00
6		Aniline	2.159	2.125	1.6	78	0.00
7	S	Phenol-d6	1.507	1.564	-3.8	83	0.00
8		2-Chlorophenol	1.321	1.436	-8.7	86	0.00
9		Benzaldehyde	1.040	0.960	7.7	73	0.00
10	C	Phenol	1.775	1.843	-3.8	83	0.00
11		bis(2-Chloroethyl)ether	1.305	1.409	-8.0	87	0.00
12		1,3-Dichlorobenzene	1.527	1.596	-4.5	83	0.00
13	C	1,4-Dichlorobenzene	1.529	1.624	-6.2	84	0.00
14		1,2-Dichlorobenzene	1.408	1.502	-6.7	83	0.00
15		Benzyl Alcohol	1.144	1.143	0.1	78	0.00
16		2,2'-oxybis(1-Chloropropane	1.780	1.869	-5.0	83	0.00
17		2-Methylphenol	1.119	1.150	-2.8	84	0.00
18		Hexachloroethane	0.563	0.600	-6.6	84	0.00
19	P	n-Nitroso-di-n-propylamine	0.934	0.977	-4.6	85	0.00
20		3+4-Methylphenols	1.324	1.460	-10.3	86	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	81	0.00
22		Acetophenone	0.439	0.466	-6.2	85	0.00
23	S	Nitrobenzene-d5	0.346	0.352	-1.7	82	0.00
24		Nitrobenzene	0.374	0.390	-4.3	84	0.00
25		Isophorone	0.663	0.689	-3.9	84	0.00
26	C	2-Nitrophenol	0.176	0.189	-7.4	86	0.00
27		2,4-Dimethylphenol	0.265	0.273	-3.0	82	0.00
28		bis(2-Chloroethoxy)methane	0.411	0.440	-7.1	87	0.00
29	C	2,4-Dichlorophenol	0.280	0.301	-7.5	85	0.00
30		1,2,4-Trichlorobenzene	0.319	0.332	-4.1	83	0.00
31		Naphthalene	0.941	1.017	-8.1	85	0.00
32		Benzoic acid	0.187	0.155	17.1	73	0.00
33		4-Chloroaniline	0.421	0.433	-2.9	82	0.00
34	C	Hexachlorobutadiene	0.184	0.195	-6.0	85	0.00
35		Caprolactam	0.091	0.089	2.2	79	0.00
36	C	4-Chloro-3-methylphenol	0.291	0.307	-5.5	85	0.00
37		2-Methylnaphthalene	0.620	0.658	-6.1	84	0.00
38		1-Methylnaphthalene	0.586	0.620	-5.8	84	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
40		1,2,4,5-Tetrachlorobenzene	0.535	0.578	-8.0	84	0.00
41	P	Hexachlorocyclopentadiene	0.209	0.187	10.5	69	0.00
42	S	2,4,6-Tribromophenol	0.194	0.198	-2.1	80	0.00
43	C	2,4,6-Trichlorophenol	0.368	0.354	3.8	75	0.00
44		2,4,5-Trichlorophenol	0.380	0.357	6.1	73	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081619\
Data File : BF116325.D
Acq On : 16 Aug 2019 18:45
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 05:30:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 2019
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	S	2-Fluorobiphenyl	1.068	1.162	-8.8	84	0.00
46		1,1'-Biphenyl	1.412	1.540	-9.1	85	0.00
47		2-Chloronaphthalene	1.175	1.239	-5.4	83	0.00
48		2-Nitroaniline	0.388	0.400	-3.1	81	0.00
49		Acenaphthylene	1.715	1.844	-7.5	83	0.00
50		Dimethylphthalate	1.362	1.409	-3.5	81	0.00
51		2,6-Dinitrotoluene	0.296	0.309	-4.4	82	0.00
52	C	Acenaphthene	1.052	1.121	-6.6	83	0.00
53		3-Nitroaniline	0.366	0.359	1.9	77	0.00
54	P	2,4-Dinitrophenol	0.126	0.104	17.5	64	0.00
55		Dibenzofuran	1.530	1.587	-3.7	80	0.00
56	P	4-Nitrophenol	0.234	0.221	5.6	74	0.00
57		2,4-Dinitrotoluene	0.394	0.388	1.5	76	0.00
58		Fluorene	1.177	1.304	-10.8	86	0.00
59		2,3,4,6-Tetrachlorophenol	0.320	0.302	5.6	74	0.00
60		Diethylphthalate	1.271	1.351	-6.3	82	0.00
61		4-Chlorophenyl-phenylether	0.596	0.665	-11.6	86	0.00
62		4-Nitroaniline	0.378	0.329	13.0	68	0.00
63		Azobenzene	1.308	1.412	-8.0	84	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
65		4,6-Dinitro-2-methylphenol	0.106	0.098	7.5	68	0.00
66	c	n-Nitrosodiphenylamine	0.602	0.659	-9.5	83	0.00
67		4-Bromophenyl-phenylether	0.214	0.237	-10.7	83	0.00
68		Hexachlorobenzene	0.248	0.267	-7.7	81	0.00
69		Atrazine	0.198	0.201	-1.5	77	0.00
70	C	Pentachlorophenol	0.120	0.107	10.8	66	0.00
71		Phenanthrene	1.022	1.106	-8.2	81	0.00
72		Anthracene	1.035	1.129	-9.1	82	0.00
73		Carbazole	0.939	1.024	-9.1	81	0.00
74		Di-n-butylphthalate	1.095	1.240	-13.2	83	0.00
75	C	Fluoranthene	1.070	1.131	-5.7	80	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	68	0.00
77		Benzidine	0.676	0.603	10.8	57	0.00
78		Pyrene	1.508	1.695	-12.4	78	0.00
79	S	Terphenyl-d14	0.949	1.090	-14.9	80	0.00
80		Butylbenzylphthalate	0.638	0.723	-13.3	76	0.00
81		Benzo(a)anthracene	1.278	1.359	-6.3	72	0.00
82		3,3'-Dichlorobenzidine	0.444	0.449	-1.1	67	0.00
83		Chrysene	1.241	1.334	-7.5	73	0.00
84		Bis(2-ethylhexyl)phthalate	0.829	0.977	-17.9	79	0.00
85	c	Di-n-octyl phthalate	1.316	1.510	-14.7	75	0.00
86		Indeno(1,2,3-cd)pyrene	1.042	1.130	-8.4	77	0.00
87	I	Perylene-d12	1.000	1.000	0.0	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081619\
Data File : BF116325.D
Acq On : 16 Aug 2019 18:45
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 05:30:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 2019
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(b)fluoranthene	1.156	1.175	-1.6	65	0.00
89	Benzo(k)fluoranthene	1.126	1.221	-8.4	75	0.00
90 C	Benzo(a)pyrene	1.034	1.090	-5.4	70	0.00
91	Dibenzo(a,h)anthracene	0.895	1.005	-12.3	78	0.00
92	Benzo(g,h,i)perylene	0.879	0.974	-10.8	79	0.00

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