Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006183.D
Acq On : 12 Jul 2021 14:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jul 12 15:00:10 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 12 14:59:51 2021
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	103	0.00
2		1,4-Dioxane	0.456	0.433	5.0	100	0.00
3		Pyridine	1.159	1.165	-0.5	102	0.00
4		n-Nitrosodimethylamine	0.456	0.426	6.6	99	0.00
5	S	2-Fluorophenol	1.162	1.154	0.7	102	0.00
6		Aniline	1.622	1.561	3.8	100	0.00
7	S	Phenol-d6	1.537	1.502	2.3	101	0.00
8		2-Chlorophenol	1.361	1.348	1.0	102	0.00
9		Benzaldehyde	0.839	0.788	6.1	97	0.00
10	C	Phenol	1.477	1.452	1.7	102	0.00
11		bis(2-Chloroethyl)ether	1.116	1.094	2.0	101	0.00
12		1,3-Dichlorobenzene	1.459	1.454	0.3	104	0.00
13	C	1,4-Dichlorobenzene	1.490	1.479	0.7	102	0.00
14		1,2-Dichlorobenzene	1.418	1.405	0.9	102	0.00
15		Benzyl Alcohol	0.930	0.898	3.4	97	0.00
16		2,2'-oxybis(1-Chloropropane	1.325	1.225	7.5	98	0.00
17		2-Methylphenol	1.005	0.991	1.4	100	0.00
18		Hexachloroethane	0.496	0.491	1.0	103	0.00
19	P	n-Nitroso-di-n-propylamine	0.838	0.799	4.7	99	0.00
20		3+4-Methylphenols	1.357	1.352	0.4	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.455	0.456	-0.2	99	0.00
23	S	Nitrobenzene-d5	0.317	0.316	0.3	99	0.00
24		Nitrobenzene	0.320	0.319	0.3	100	0.00
25		Isophorone	0.601	0.588	2.2	97	0.00
26	C	2-Nitrophenol	0.178	0.191	-7.3	101	0.00
27		2,4-Dimethylphenol	0.263	0.270	-2.7	98	0.00
28		bis(2-Chloroethoxy)methane	0.333	0.332	0.3	98	0.00
29	C	2,4-Dichlorophenol	0.289	0.306	-5.9	101	0.00
30		1,2,4-Trichlorobenzene	0.307	0.319	-3.9	100	0.00
31		Naphthalene	1.043	1.037	0.6	99	0.00
32		Benzoic acid	0.177	0.200	-13.0	108	0.00
33		4-Chloroaniline	0.402	0.411	-2.2	97	0.00
34	C	Hexachlorobutadiene	0.169	0.181	-7.1	102	0.00
35		Caprolactam	0.091	0.092	-1.1	97	0.00
36	C	4-Chloro-3-methylphenol	0.312	0.309	1.0	97	0.00
37		2-Methylnaphthalene	0.723	0.719	0.6	98	0.00
38		1-Methylnaphthalene	0.685	0.675	1.5	96	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
40		1,2,4,5-Tetrachlorobenzene	0.496	0.561	-13.1	99	0.00
41	P	Hexachlorocyclopentadiene	0.076	0.072	5.3	83	0.00
42	S	2,4,6-Tribromophenol	0.238	0.255	-7.1	89	0.00
43	C	2,4,6-Trichlorophenol	0.359	0.410	-14.2	100	0.00
44		2,4,5-Trichlorophenol	0.385	0.440	-14.3	99	0.00
45	S	2-Fluorobiphenyl	1.353	1.466	-8.4	98	0.00
46		1,1'-Biphenyl	1.502	1.616	-7.6	97	0.00
47		2-Chloronaphthalene	1.155	1.250	-8.2	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006183.D
Acq On : 12 Jul 2021 14:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jul 12 15:00:10 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 12 14:59:51 2021
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)
48		2-Nitroaniline	0.287	0.306	-6.6	95	0.00
49		Acenaphthylene	1.860	1.917	-3.1	93	0.00
50		Dimethylphthalate	1.438	1.498	-4.2	94	0.00
51		2,6-Dinitrotoluene	0.324	0.345	-6.5	95	0.00
52	C	Acenaphthene	1.131	1.115	1.4	90	0.00
53		3-Nitroaniline	0.328	0.349	-6.4	92	0.00
54	P	2,4-Dinitrophenol	0.132	0.151	-14.4	97	0.00
55		Dibenzofuran	1.761	1.724	2.1	88	0.00
56	P	4-Nitrophenol	0.209	0.217	-3.8	91	0.00
57		2,4-Dinitrotoluene	0.437	0.448	-2.5	89	0.00
58		Fluorene	1.415	1.364	3.6	87	0.00
59		2,3,4,6-Tetrachlorophenol	0.301	0.316	-5.0	90	0.00
60		Diethylphthalate	1.475	1.426	3.3	88	0.00
61		4-Chlorophenyl-phenylether	0.620	0.608	1.9	87	0.00
62		4-Nitroaniline	0.302	0.316	-4.6	88	0.00
63		Azobenzene	1.231	1.158	5.9	87	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
65		4,6-Dinitro-2-methylphenol	0.121	0.132	-9.1	89	0.00
66	c	n-Nitrosodiphenylamine	0.651	0.655	-0.6	86	0.00
67		4-Bromophenyl-phenylether	0.197	0.209	-6.1	87	0.00
68		Hexachlorobenzene	0.234	0.249	-6.4	87	0.00
69		Atrazine	0.204	0.213	-4.4	87	0.00
70	C	Pentachlorophenol	0.108	0.124	-14.8	94	0.00
71		Phenanthrene	1.130	1.138	-0.7	87	0.00
72		Anthracene	1.107	1.123	-1.4	86	0.00
73		Carbazole	1.097	1.129	-2.9	88	0.00
74		Di-n-butylphthalate	1.320	1.383	-4.8	90	0.00
75	C	Fluoranthene	1.238	1.273	-2.8	88	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
77		Benzidine	0.593	0.590	0.5	81	0.00
78		Pyrene	1.352	1.324	2.1	86	0.00
79	S	Terphenyl-d14	1.014	1.043	-2.9	88	0.00
80		Butylbenzylphthalate	0.671	0.670	0.1	89	0.00
81		Benzo(a)anthracene	1.265	1.280	-1.2	89	0.00
82		3,3'-Dichlorobenzidine	0.371	0.395	-6.5	91	0.00
83		Chrysene	1.250	1.249	0.1	88	0.00
84		Bis(2-ethylhexyl)phthalate	0.990	1.002	-1.2	92	0.00
85	c	Di-n-octyl phthalate	1.743	1.807	-3.7	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Indeno(1,2,3-cd)pyrene	1.475	1.518	-2.9	94	0.00
88		Benzo(b)fluoranthene	1.213	1.232	-1.6	91	0.00
89		Benzo(k)fluoranthene	1.186	1.171	1.3	93	0.00
90	C	Benzo(a)pyrene	1.138	1.145	-0.6	93	0.00
91		Dibenzo(a,h)anthracene	1.270	1.318	-3.8	95	0.00
92		Benzo(g,h,i)perylene	1.250	1.277	-2.2	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006183.D
Acq On : 12 Jul 2021 14:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jul 12 15:00:10 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 12 14:59:51 2021
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)

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