Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
Data File : BP006843.D
Acq On : 24 Aug 2021 12:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Aug 24 13:41:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 24 13:40:38 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	87	0.00
2		1,4-Dioxane	0.567	0.543	4.2	84	0.00
3		Pyridine	1.666	1.553	6.8	77	0.00
4		n-Nitrosodimethylamine	0.629	0.524	16.7	69	0.00
5	S	2-Fluorophenol	1.302	1.351	-3.8	87	0.00
6		Aniline	2.009	1.986	1.1	78	0.00
7	S	Phenol-d6	1.850	1.846	0.2	81	0.01
8		2-Chlorophenol	1.409	1.495	-6.1	87	0.00
9		Benzaldehyde	1.113	1.029	7.5	78	0.00
10	C	Phenol	1.855	1.832	1.2	80	0.00
11		bis(2-Chloroethyl)ether	1.408	1.368	2.8	80	0.00
12		1,3-Dichlorobenzene	1.468	1.559	-6.2	89	0.00
13	C	1,4-Dichlorobenzene	1.490	1.582	-6.2	89	0.00
14		1,2-Dichlorobenzene	1.430	1.519	-6.2	88	0.00
15		Benzyl Alcohol	1.387	1.381	0.4	78	0.00
16		2,2'-oxybis(1-Chloropropane	1.676	1.376	17.9	66	0.00
17		2-Methylphenol	1.171	1.227	-4.8	83	0.01
18		Hexachloroethane	0.590	0.616	-4.4	88	0.00
19	P	n-Nitroso-di-n-propylamine	1.241	1.171	5.6	73	0.00
20		3+4-Methylphenols	1.594	1.670	-4.8	82	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	80	0.01
22		Acetophenone	0.530	0.554	-4.5	80	0.00
23	S	Nitrobenzene-d5	0.433	0.447	-3.2	79	0.00
24		Nitrobenzene	0.450	0.453	-0.7	78	0.00
25		Isophorone	0.778	0.811	-4.2	78	0.01
26	C	2-Nitrophenol	0.164	0.205	-25.0#	94	0.00
27		2,4-Dimethylphenol	0.274	0.299	-9.1	83	0.01
28		bis(2-Chloroethoxy)methane	0.416	0.425	-2.2	79	0.00
29	C	2,4-Dichlorophenol	0.278	0.323	-16.2	87	0.01
30		1,2,4-Trichlorobenzene	0.299	0.333	-11.4	88	0.01
31		Naphthalene	1.075	1.148	-6.8	83	0.01
32		Benzoic acid	0.212	0.266	-25.5#	96	0.02
33		4-Chloroaniline	0.435	0.467	-7.4	81	0.00
34	C	Hexachlorobutadiene	0.174	0.204	-17.2	94	0.01
35		Caprolactam	0.100	0.114	-14.0	80	0.02
36	C	4-Chloro-3-methylphenol	0.353	0.388	-9.9	81	0.02
37		2-Methylnaphthalene	0.728	0.784	-7.7	82	0.01
38		1-Methylnaphthalene	0.692	0.743	-7.4	82	0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	76	0.01
40		1,2,4,5-Tetrachlorobenzene	0.523	0.587	-12.2	86	0.01
41	P	Hexachlorocyclopentadiene	0.278	0.265	4.7	71	0.01
42	S	2,4,6-Tribromophenol	0.209	0.251	-20.1	86	0.02
43	C	2,4,6-Trichlorophenol	0.357	0.428	-19.9	88	0.01
44		2,4,5-Trichlorophenol	0.396	0.461	-16.4	85	0.01
45	S	2-Fluorobiphenyl	1.337	1.452	-8.6	82	0.01
46		1,1'-Biphenyl	1.514	1.630	-7.7	81	0.01
47		2-Chloronaphthalene	1.166	1.265	-8.5	82	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
Data File : BP006843.D
Acq On : 24 Aug 2021 12:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Aug 24 13:41:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 24 13:40:38 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	0.411	0.441	-7.3	75	0.01
49		Acenaphthylene	1.880	2.061	-9.6	81	0.01
50		Dimethylphthalate	1.456	1.576	-8.2	80	0.02
51		2,6-Dinitrotoluene	0.325	0.367	-12.9	81	0.01
52	C	Acenaphthene	1.144	1.219	-6.6	79	0.01
53		3-Nitroaniline	0.360	0.401	-11.4	78	0.01
54	P	2,4-Dinitrophenol	0.170	0.200	-17.6	86	0.01
55		Dibenzofuran	1.799	1.925	-7.0	79	0.01
56	P	4-Nitrophenol	0.313	0.346	-10.5	76	0.02
57		2,4-Dinitrotoluene	0.436	0.506	-16.1	82	0.02
58		Fluorene	1.438	1.546	-7.5	79	0.02
59		2,3,4,6-Tetrachlorophenol	0.332	0.380	-14.5	83	0.02
60		Diethylphthalate	1.529	1.691	-10.6	81	0.01
61		4-Chlorophenyl-phenylether	0.651	0.713	-9.5	81	0.01
62		4-Nitroaniline	0.382	0.402	-5.2	72	0.02
63		Azobenzene	1.783	1.819	-2.0	73	0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	75	0.01
65		4,6-Dinitro-2-methylphenol	0.122	0.142	-16.4	84	0.02
66	c	n-Nitrosodiphenylamine	0.629	0.676	-7.5	78	0.02
67		4-Bromophenyl-phenylether	0.192	0.217	-13.0	84	0.01
68		Hexachlorobenzene	0.218	0.244	-11.9	84	0.01
69		Atrazine	0.196	0.208	-6.1	75	0.01
70	C	Pentachlorophenol	0.139	0.159	-14.4	80	0.02
71		Phenanthrene	1.128	1.202	-6.6	78	0.02
72		Anthracene	1.090	1.189	-9.1	79	0.01
73		Carbazole	1.112	1.195	-7.5	76	0.02
74		Di-n-butylphthalate	1.267	1.467	-15.8	83	0.02
75	C	Fluoranthene	1.266	1.398	-10.4	79	0.02
76	I	Chrysene-d12	1.000	1.000	0.0	75	0.02
77		Benzidine	0.500	0.597	-19.4	77	0.01
78		Pyrene	1.336	1.442	-7.9	79	0.02
79	S	Terphenyl-d14	0.998	1.100	-10.2	82	0.02
80		Butylbenzylphthalate	0.587	0.716	-22.0	86	0.02
81		Benzo(a)anthracene	1.307	1.408	-7.7	79	0.02
82		3,3'-Dichlorobenzidine	0.343	0.410	-19.5	85	0.02
83		Chrysene	1.267	1.359	-7.3	79	0.02
84		Bis(2-ethylhexyl)phthalate	0.883	1.037	-17.4	85	0.02
85	c	Di-n-octyl phthalate	1.455	1.825	-25.4#	90	0.02
86	I	Perylene-d12	1.000	1.000	0.0	78	0.03
87		Indeno(1,2,3-cd)pyrene	1.450	1.589	-9.6	82	0.05
88		Benzo(b)fluoranthene	1.300	1.421	-9.3	82	0.03
89		Benzo(k)fluoranthene	1.248	1.278	-2.4	76	0.03
90	C	Benzo(a)pyrene	1.166	1.275	-9.3	81	0.03
91		Dibenzo(a,h)anthracene	1.254	1.364	-8.8	82	0.05
92		Benzo(g,h,i)perylene	1.202	1.318	-9.7	82	0.05

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
Data File : BP006843.D
Acq On : 24 Aug 2021 12:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Aug 24 13:41:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 24 13:40:38 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 = 2