Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080823\
Data File : BP016529.D
Acq On : 08 Aug 2023 09:52
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Aug 08 16:03:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 08 16:03:02 2023
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	52	0.00
2		1,4-Dioxane	0.478	0.441	7.7	50#	0.00
3		Pyridine	1.429	1.561	-9.2	58	0.00
4		n-Nitrosodimethylamine	0.539	0.540	-0.2	53	0.00
5	S	2-Fluorophenol	1.226	1.204	1.8	51	0.00
6		Aniline	2.076	2.315	-11.5	57	0.00
7	S	Phenol-d6	1.680	1.872	-11.4	57	0.00
8		2-Chlorophenol	1.285	1.332	-3.7	53	0.00
9		Benzaldehyde	0.866	0.884	-2.1	57	0.00
10	C	Phenol	1.632	1.815	-11.2	58	0.00
11		bis(2-Chloroethyl)ether	1.326	1.398	-5.4	55	0.00
12		1,3-Dichlorobenzene	1.380	1.328	3.8	50	0.00
13	C	1,4-Dichlorobenzene	1.400	1.355	3.2	50	0.00
14		1,2-Dichlorobenzene	1.348	1.333	1.1	51	0.00
15		Benzyl Alcohol	1.148	1.317	-14.7	58	0.00
16		2,2'-oxybis(1-Chloropropane	1.504	1.635	-8.7	57	0.00
17		2-Methylphenol	1.176	1.328	-12.9	58	0.00
18		Hexachloroethane	0.499	0.475	4.8	50#	0.00
19	P	n-Nitroso-di-n-propylamine	1.053	1.200	-14.0	59	0.00
20		3+4-Methylphenols	1.597	1.838	-15.1	58	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	60	0.00
22		Acetophenone	0.505	0.495	2.0	58	0.00
23	S	Nitrobenzene-d5	0.357	0.359	-0.6	59	0.00
24		Nitrobenzene	0.367	0.365	0.5	59	0.00
25		Isophorone	0.721	0.724	-0.4	59	0.00
26	C	2-Nitrophenol	0.141	0.155	-9.9	61	0.00
27		2,4-Dimethylphenol	0.323	0.318	1.5	58	0.00
28		bis(2-Chloroethoxy)methane	0.442	0.438	0.9	59	0.00
29	C	2,4-Dichlorophenol	0.302	0.298	1.3	56	0.00
30		1,2,4-Trichlorobenzene	0.327	0.295	9.8	52	0.00
31		Naphthalene	1.046	1.001	4.3	57	0.00
32		Benzoic acid	0.191	0.236	-23.6	72	0.00
33		4-Chloroaniline	0.472	0.488	-3.4	61	0.00
34	C	Hexachlorobutadiene	0.194	0.164	15.5	48#	0.00
35		Caprolactam	0.110	0.131	-19.1	69	0.00
36	C	4-Chloro-3-methylphenol	0.336	0.368	-9.5	63	0.00
37		2-Methylnaphthalene	0.760	0.757	0.4	58	0.00
38		1-Methylnaphthalene	0.713	0.716	-0.4	59	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	65	0.00
40		1,2,4,5-Tetrachlorobenzene	0.608	0.523	14.0	54	0.00
41	P	Hexachlorocyclopentadiene	0.305	0.241	21.0	47#	0.00
42	S	2,4,6-Tribromophenol	0.294	0.275	6.5	56	0.00
43	C	2,4,6-Trichlorophenol	0.390	0.375	3.8	59	0.00
44		2,4,5-Trichlorophenol	0.473	0.457	3.4	59	0.00
45	S	2-Fluorobiphenyl	1.395	1.258	9.8	58	0.00
46		1,1'-Biphenyl	1.525	1.406	7.8	60	0.00
47		2-Chloronaphthalene	1.150	1.069	7.0	60	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080823\
Data File : BP016529.D
Acq On : 08 Aug 2023 09:52
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Aug 08 16:03:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 08 16:03:02 2023
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	0.324	0.359	-10.8	69	0.00
49		Acenaphthylene	1.908	1.838	3.7	62	0.00
50		Dimethylphthalate	1.559	1.512	3.0	63	0.00
51		2,6-Dinitrotoluene	0.319	0.332	-4.1	65	0.00
52	C	Acenaphthene	1.135	1.092	3.8	62	0.00
53		3-Nitroaniline	0.361	0.396	-9.7	69	0.00
54	P	2,4-Dinitrophenol	0.121	0.151	-24.8	81	0.00
55		Dibenzofuran	1.861	1.795	3.5	62	0.00
56	P	4-Nitrophenol	0.283	0.338	-19.4	75	0.00
57		2,4-Dinitrotoluene	0.416	0.459	-10.3	67	0.00
58		Fluorene	1.462	1.455	0.5	64	0.00
59		2,3,4,6-Tetrachlorophenol	0.389	0.391	-0.5	61	0.00
60		Diethylphthalate	1.549	1.528	1.4	64	0.00
61		4-Chlorophenyl-phenylether	0.738	0.702	4.9	60	0.00
62		4-Nitroaniline	0.358	0.415	-15.9	72	0.00
63		Azobenzene	1.462	1.487	-1.7	68	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
65		4,6-Dinitro-2-methylphenol	0.084	0.095	-13.1	78	0.00
66	c	n-Nitrosodiphenylamine	0.589	0.535	9.2	64	0.00
67		4-Bromophenyl-phenylether	0.212	0.183	13.7	58	0.00
68		Hexachlorobenzene	0.236	0.204	13.6	58	0.00
69		Atrazine	0.175	0.165	5.7	65	0.00
70	C	Pentachlorophenol	0.163	0.159	2.5	66	0.00
71		Phenanthrene	1.062	1.008	5.1	67	0.00
72		Anthracene	1.068	1.031	3.5	68	0.00
73		Carbazole	1.001	1.013	-1.2	72	0.00
74		Di-n-butylphthalate	1.158	1.176	-1.6	70	0.00
75	C	Fluoranthene	1.245	1.272	-2.2	71	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	80	0.00
77		Benzidine	0.347	0.402	-15.9	84	0.00
78		Pyrene	1.341	1.240	7.5	73	0.00
79	S	Terphenyl-d14	0.947	0.865	8.7	69	0.00
80		Butylbenzylphthalate	0.527	0.541	-2.7	80	0.00
81		Benzo(a)anthracene	1.342	1.299	3.2	76	0.00
82		3,3'-Dichlorobenzidine	0.441	0.443	-0.5	75	0.00
83		Chrysene	1.281	1.241	3.1	77	0.00
84		Bis(2-ethylhexyl)phthalate	0.791	0.804	-1.6	79	0.00
85	c	Di-n-octyl phthalate	1.330	1.458	-9.6	86	0.00
86	I	Perylene-d12	1.000	1.000	0.0	90	0.00
87		Indeno(1,2,3-cd)pyrene	1.348	1.351	-0.2	90	0.00
88		Benzo(b)fluoranthene	1.159	1.095	5.5	81	0.00
89		Benzo(k)fluoranthene	1.179	1.126	4.5	83	0.00
90	C	Benzo(a)pyrene	1.123	1.100	2.0	85	0.00
91		Dibenzo(a,h)anthracene	1.124	1.122	0.2	89	0.00
92		Benzo(g,h,i)perylene	1.084	1.107	-2.1	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080823\
Data File : BP016529.D
Acq On : 08 Aug 2023 09:52
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Aug 08 16:03:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 08 16:03:02 2023
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