Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003741.D
Acq On : 26 Oct 2020 15:19
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Oct 26 16:57:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 26 16:22:21 2020
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	90	0.00
2		1,4-Dioxane	0.511	0.456	10.8	85	0.00
3		Pyridine	1.445	1.298	10.2	83	0.00
4		n-Nitrosodimethylamine	0.617	0.507	17.8	76	0.00
5	S	2-Fluorophenol	1.196	1.079	9.8	83	0.00
6		Aniline	1.876	1.705	9.1	83	0.00
7	S	Phenol-d6	1.690	1.544	8.6	84	0.00
8		2-Chlorophenol	1.191	1.093	8.2	84	0.00
9		Benzaldehyde	0.891	0.719	19.3	79	0.00
10	C	Phenol	1.609	1.466	8.9	84	0.00
11		bis(2-Chloroethyl)ether	1.277	1.138	10.9	82	0.00
12		1,3-Dichlorobenzene	1.536	1.380	10.2	84	0.00
13	C	1,4-Dichlorobenzene	1.551	1.398	9.9	84	0.00
14		1,2-Dichlorobenzene	1.469	1.329	9.5	84	0.00
15		Benzyl Alcohol	1.292	1.144	11.5	80	0.00
16		2,2'-oxybis(1-Chloropropane	1.266	1.124	11.2	82	0.00
17		2-Methylphenol	1.066	0.976	8.4	83	0.00
18		Hexachloroethane	0.598	0.526	12.0	81	0.00
19	P	n-Nitroso-di-n-propylamine	1.033	0.912	11.7	80	0.00
20		3+4-Methylphenols	1.435	1.313	8.5	84	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
22		Acetophenone	0.585	0.509	13.0	82	0.00
23	S	Nitrobenzene-d5	0.473	0.406	14.2	80	0.00
24		Nitrobenzene	0.455	0.395	13.2	82	0.00
25		Isophorone	0.762	0.665	12.7	81	0.00
26	C	2-Nitrophenol	0.170	0.157	7.6	84	0.00
27		2,4-Dimethylphenol	0.267	0.241	9.7	84	0.00
28		bis(2-Chloroethoxy)methane	0.486	0.430	11.5	84	0.00
29	C	2,4-Dichlorophenol	0.350	0.314	10.3	83	0.00
30		1,2,4-Trichlorobenzene	0.449	0.400	10.9	85	0.00
31		Naphthalene	1.078	0.959	11.0	85	0.00
32		Benzoic acid	0.158	0.151	4.4	82	0.00
33		4-Chloroaniline	0.444	0.409	7.9	85	0.00
34	C	Hexachlorobutadiene	0.339	0.293	13.6	82	0.00
35		Caprolactam	0.103	0.096	6.8	84	0.00
36	C	4-Chloro-3-methylphenol	0.380	0.347	8.7	85	0.00
37		2-Methylnaphthalene	0.782	0.704	10.0	85	0.00
38		1-Methylnaphthalene	0.740	0.664	10.3	85	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
40		1,2,4,5-Tetrachlorobenzene	0.792	0.690	12.9	85	0.00
41	P	Hexachlorocyclopentadiene	0.459	0.386	15.9	79	0.00
42	S	2,4,6-Tribromophenol	0.312	0.294	5.8	88	0.00
43	C	2,4,6-Trichlorophenol	0.471	0.428	9.1	85	0.00
44		2,4,5-Trichlorophenol	0.490	0.442	9.8	85	0.00
45	S	2-Fluorobiphenyl	1.547	1.348	12.9	85	0.00
46		1,1'-Biphenyl	1.560	1.358	12.9	85	0.00
47		2-Chloronaphthalene	1.304	1.147	12.0	86	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003741.D
Acq On : 26 Oct 2020 15:19
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Oct 26 16:57:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 26 16:22:21 2020
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	0.385	0.338	12.2	81	0.00
49		Acenaphthylene	1.862	1.648	11.5	85	0.00
50		Dimethylphthalate	1.643	1.445	12.1	85	0.00
51		2,6-Dinitrotoluene	0.334	0.303	9.3	86	0.00
52	C	Acenaphthene	1.257	1.115	11.3	85	0.00
53		3-Nitroaniline	0.328	0.306	6.7	87	0.00
54	P	2,4-Dinitrophenol	0.153	0.142	7.2	81	0.00
55		Dibenzofuran	1.900	1.680	11.6	86	0.00
56	P	4-Nitrophenol	0.246	0.231	6.1	87	0.00
57		2,4-Dinitrotoluene	0.452	0.418	7.5	86	0.00
58		Fluorene	1.533	1.360	11.3	86	0.00
59		2,3,4,6-Tetrachlorophenol	0.470	0.436	7.2	87	0.00
60		Diethylphthalate	1.610	1.409	12.5	84	0.00
61		4-Chlorophenyl-phenylether	0.927	0.821	11.4	87	0.00
62		4-Nitroaniline	0.364	0.338	7.1	86	0.00
63		Azobenzene	1.513	1.289	14.8	82	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
65		4,6-Dinitro-2-methylphenol	0.105	0.094	10.5	88	0.00
66	c	n-Nitrosodiphenylamine	0.595	0.512	13.9	86	0.00
67		4-Bromophenyl-phenylether	0.268	0.233	13.1	87	0.00
68		Hexachlorobenzene	0.308	0.272	11.7	88	0.00
69		Atrazine	0.232	0.200	13.8	84	0.00
70	C	Pentachlorophenol	0.152	0.132	13.2	85	0.00
71		Phenanthrene	1.111	0.969	12.8	88	0.00
72		Anthracene	1.073	0.943	12.1	88	0.00
73		Carbazole	0.955	0.855	10.5	89	0.00
74		Di-n-butylphthalate	1.166	1.023	12.3	85	0.00
75	C	Fluoranthene	1.365	1.228	10.0	89	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
77		Benzidine	0.479	0.431	10.0	87	0.00
78		Pyrene	1.340	1.159	13.5	89	0.00
79	S	Terphenyl-d14	1.137	0.978	14.0	89	0.00
80		Butylbenzylphthalate	0.491	0.425	13.4	86	0.00
81		Benzo(a)anthracene	1.333	1.181	11.4	91	0.00
82		3,3'-Dichlorobenzidine	0.467	0.430	7.9	92	0.00
83		Chrysene	1.264	1.118	11.6	91	0.00
84		Bis(2-ethylhexyl)phthalate	0.702	0.608	13.4	86	0.00
85	c	Di-n-octyl phthalate	1.140	1.008	11.6	87	0.00
86	I	Perylene-d12	1.000	1.000	0.0	105	0.00
87		Indeno(1,2,3-cd)pyrene	1.435	1.361	5.2	104	0.00
88		Benzo(b)fluoranthene	1.396	1.206	13.6	92	0.00
89		Benzo(k)fluoranthene	1.332	1.192	10.5	97	0.00
90	C	Benzo(a)pyrene	1.244	1.113	10.5	96	0.00
91		Dibenzo(a,h)anthracene	1.200	1.143	4.7	104	0.00
92		Benzo(g,h,i)perylene	1.112	1.076	3.2	107	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003741.D
Acq On : 26 Oct 2020 15:19
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Oct 26 16:57:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 26 16:22:21 2020
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