Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082521\
Data File : BP006882.D
Acq On : 25 Aug 2021 16:43
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Aug 25 17:48:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 25 11:27:41 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	109	0.00
2		1,4-Dioxane	0.567	0.571	-0.7	111	0.04
3		Pyridine	1.666	1.472	11.6	92	0.04
4		n-Nitrosodimethylamine	0.629	0.472	25.0	79	0.03
5	S	2-Fluorophenol	1.302	1.326	-1.8	108	0.02
6		Aniline	2.009	1.753	12.7	87	0.01
7	S	Phenol-d6	1.850	1.650	10.8	91	0.01
8		2-Chlorophenol	1.409	1.374	2.5	101	0.00
9		Benzaldehyde	1.113	0.911	18.1	87	0.00
10	C	Phenol	1.855	1.622	12.6	89	0.01
11		bis(2-Chloroethyl)ether	1.408	1.239	12.0	91	0.01
12		1,3-Dichlorobenzene	1.468	1.481	-0.9	107	0.00
13	C	1,4-Dichlorobenzene	1.490	1.491	-0.1	106	0.00
14		1,2-Dichlorobenzene	1.430	1.405	1.7	102	0.00
15		Benzyl Alcohol	1.387	1.176	15.2	84	0.00
16		2,2'-oxybis(1-Chloropropane	1.676	1.218	27.3#	74	0.00
17		2-Methylphenol	1.171	1.053	10.1	90	0.01
18		Hexachloroethane	0.590	0.584	1.0	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.241	0.871	29.8#	69	0.00
20		3+4-Methylphenols	1.594	1.411	11.5	87	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.530	0.517	2.5	86	0.00
23	S	Nitrobenzene-d5	0.433	0.418	3.5	85	0.00
24		Nitrobenzene	0.450	0.426	5.3	84	0.00
25		Isophorone	0.778	0.728	6.4	81	0.00
26	C	2-Nitrophenol	0.164	0.190	-15.9	100	0.00
27		2,4-Dimethylphenol	0.274	0.272	0.7	87	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.391	6.0	83	0.00
29	C	2,4-Dichlorophenol	0.278	0.297	-6.8	92	0.00
30		1,2,4-Trichlorobenzene	0.299	0.318	-6.4	97	0.00
31		Naphthalene	1.075	1.079	-0.4	90	0.00
32		Benzoic acid	0.212	0.208	1.9	86	0.02
33		4-Chloroaniline	0.435	0.416	4.4	83	0.00
34	C	Hexachlorobutadiene	0.174	0.198	-13.8	105	0.00
35		Caprolactam	0.100	0.093	7.0	75	0.02
36	C	4-Chloro-3-methylphenol	0.353	0.339	4.0	81	0.00
37		2-Methylnaphthalene	0.728	0.714	1.9	86	0.00
38		1-Methylnaphthalene	0.692	0.669	3.3	85	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
40		1,2,4,5-Tetrachlorobenzene	0.523	0.593	-13.4	90	0.00
41	P	Hexachlorocyclopentadiene	0.278	0.262	5.8	73	0.00
42	S	2,4,6-Tribromophenol	0.209	0.230	-10.0	82	0.00
43	C	2,4,6-Trichlorophenol	0.357	0.415	-16.2	88	0.00
44		2,4,5-Trichlorophenol	0.396	0.442	-11.6	85	0.00
45	S	2-Fluorobiphenyl	1.337	1.437	-7.5	85	0.00
46		1,1'-Biphenyl	1.514	1.605	-6.0	83	0.00
47		2-Chloronaphthalene	1.166	1.240	-6.3	83	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.411	0.409	0.5	73	0.00
49		Acenaphthylene	1.880	1.957	-4.1	80	0.00
50		Dimethylphthalate	1.456	1.469	-0.9	77	0.00
51		2,6-Dinitrotoluene	0.325	0.335	-3.1	76	0.00
52	C	Acenaphthene	1.144	1.168	-2.1	78	0.00
53		3-Nitroaniline	0.360	0.361	-0.3	73	0.00
54	P	2,4-Dinitrophenol	0.170	0.170	0.0	76	0.00
55		Dibenzofuran	1.799	1.818	-1.1	78	0.00
56	P	4-Nitrophenol	0.313	0.296	5.4	67	0.01
57		2,4-Dinitrotoluene	0.436	0.455	-4.4	76	0.00
58		Fluorene	1.438	1.440	-0.1	76	0.00
59		2,3,4,6-Tetrachlorophenol	0.332	0.356	-7.2	81	0.00
60		Diethylphthalate	1.529	1.541	-0.8	76	0.00
61		4-Chlorophenyl-phenylether	0.651	0.666	-2.3	79	0.00
62		4-Nitroaniline	0.382	0.368	3.7	69	0.00
63		Azobenzene	1.783	1.682	5.7	70	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
65		4,6-Dinitro-2-methylphenol	0.122	0.131	-7.4	76	0.01
66	c	n-Nitrosodiphenylamine	0.629	0.655	-4.1	74	0.00
67		4-Bromophenyl-phenylether	0.192	0.213	-10.9	80	0.00
68		Hexachlorobenzene	0.218	0.234	-7.3	78	0.00
69		Atrazine	0.196	0.207	-5.6	73	0.00
70	C	Pentachlorophenol	0.139	0.147	-5.8	73	0.00
71		Phenanthrene	1.128	1.143	-1.3	73	0.00
72		Anthracene	1.090	1.130	-3.7	73	0.00
73		Carbazole	1.112	1.102	0.9	69	0.00
74		Di-n-butylphthalate	1.267	1.377	-8.7	76	0.00
75	C	Fluoranthene	1.266	1.260	0.5	70	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	66	0.00
77		Benzidine	0.500	0.524	-4.8	59	0.00
78		Pyrene	1.336	1.444	-8.1	69	0.00
79	S	Terphenyl-d14	0.998	1.085	-8.7	71	0.00
80		Butylbenzylphthalate	0.587	0.698	-18.9	73	0.00
81		Benzo(a)anthracene	1.307	1.345	-2.9	66	0.00
82		3,3'-Dichlorobenzidine	0.343	0.382	-11.4	70	0.00
83		Chrysene	1.267	1.308	-3.2	66	0.01
84		Bis(2-ethylhexyl)phthalate	0.883	0.999	-13.1	71	0.00
85	c	Di-n-octyl phthalate	1.455	1.753	-20.5#	75	0.00
86	I	Perylene-d12	1.000	1.000	0.0	68	0.01
87		Indeno(1,2,3-cd)pyrene	1.450	1.507	-3.9	68	0.02
88		Benzo(b)fluoranthene	1.300	1.326	-2.0	67	0.01
89		Benzo(k)fluoranthene	1.248	1.217	2.5	64	0.01
90	C	Benzo(a)pyrene	1.166	1.205	-3.3	67	0.01
91		Dibenzo(a,h)anthracene	1.254	1.298	-3.5	68	0.02
92		Benzo(g,h,i)perylene	1.202	1.244	-3.5	68	0.02

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 1