Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030315\
Data File : BF077562.D
Acq On : 3 Mar 2015 23:33
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 01:30:44 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 03 23:57:38 2015
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	91	0.00
2		1,4-Dioxane	0.508	0.508	0.0	92	0.00
3		Pyridine	1.454	1.422	2.2	83	0.00
4		n-Nitrosodimethylamine	0.632	0.644	-1.9	95	0.00
5	S	2-Fluorophenol	1.198	1.217	-1.6	91	0.00
6		Aniline	2.129	2.009	5.6	83	0.00
7	S	Phenol-d6	1.540	1.469	4.6	84	0.00
8		2-Chlorophenol	1.413	1.395	1.3	88	0.00
9		Benzaldehyde	0.935	0.990	-5.9	97	0.00
10	C	Phenol	1.828	1.799	1.6	84	0.00
11		bis(2-Chloroethyl)ether	1.313	1.160	11.7	80	0.00
12		1,3-Dichlorobenzene	1.539	1.422	7.6	82	0.00
13	C	1,4-Dichlorobenzene	1.566	1.490	4.9	84	0.00
14		1,2-Dichlorobenzene	1.489	1.377	7.5	81	0.00
15		Benzyl Alcohol	1.171	1.082	7.6	82	0.00
16		2,2'-oxybis(1-Chloropropane	1.877	1.705	9.2	80	0.00
17		2-Methylphenol	1.105	1.088	1.5	83	0.00
18		Hexachloroethane	0.523	0.506	3.3	86	0.00
19	P	n-Nitroso-di-n-propylamine	0.978	0.928	5.1	82	0.00
20		3+4-Methylphenols	1.455	1.384	4.9	82	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
22		Acetophenone	0.470	0.480	-2.1	89	0.00
23	S	Nitrobenzene-d5	0.322	0.317	1.6	81	0.00
24		Nitrobenzene	0.338	0.337	0.3	83	0.00
25		Isophorone	0.638	0.581	8.9	73	0.00
26	C	2-Nitrophenol	0.139	0.165	-18.7	91	0.00
27		2,4-Dimethylphenol	0.310	0.300	3.2	81	0.00
28		bis(2-Chloroethoxy)methane	0.403	0.390	3.2	81	0.00
29	C	2,4-Dichlorophenol	0.291	0.290	0.3	82	0.00
30		1,2,4-Trichlorobenzene	0.320	0.307	4.1	80	0.00
31		Naphthalene	1.000	0.989	1.1	84	0.00
32		Benzoic acid	0.151	0.171	-13.2	89	0.00
33		4-Chloroaniline	0.445	0.413	7.2	75	0.00
34	C	Hexachlorobutadiene	0.183	0.163	10.9	74	0.00
35		Caprolactam	0.089	0.088	1.1	80	0.00
36	C	4-Chloro-3-methylphenol	0.293	0.284	3.1	78	0.00
37		2-Methylnaphthalene	0.710	0.693	2.4	79	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	73	0.00
39		1,2,4,5-Tetrachlorobenzene	0.574	0.626	-9.1	76	0.00
40	P	Hexachlorocyclopentadiene	0.308	0.328	-6.5	70	0.00
41	S	2,4,6-Tribromophenol	0.193	0.186	3.6	66	0.00
42	C	2,4,6-Trichlorophenol	0.380	0.429	-12.9	76	0.00
43		2,4,5-Trichlorophenol	0.406	0.463	-14.0	78	0.00
44	S	2-Fluorobiphenyl	1.283	1.379	-7.5	77	0.00

Data Path : Z:\HPCHEM1\BNA F\Data\BF030315\
Data File : BF077562.D
Acq On : 3 Mar 2015 23:33
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 01:30:44 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 03 23:57:38 2015
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)
45		1,1'-Biphenyl	1.515	1.652	-9.0	76	0.00
46		2-Chloronaphthalene	1.188	1.335	-12.4	81	0.00
47		2-Nitroaniline	0.293	0.366	-24.9	86	0.00
48		Acenaphthylene	1.881	2.124	-12.9	81	0.00
49		Dimethylphthalate	1.406	1.408	-0.1	71	0.00
50		2,6-Dinitrotoluene	0.282	0.318	-12.8	73	0.00
51	C	Acenaphthene	1.123	1.091	2.8	68	0.00
52		3-Nitroaniline	0.325	0.385	-18.5	79	0.00
53	P	2,4-Dinitrophenol	0.086	0.086	0.0	69	0.00
54		Dibenzofuran	1.733	1.603	7.5	65	0.00
55	P	4-Nitrophenol	0.261	0.277	-6.1	69	0.00
56		2,4-Dinitrotoluene	0.381	0.388	-1.8	65	0.00
57		Fluorene	1.386	1.297	6.4	66	0.00
58		2,3,4,6-Tetrachlorophenol	0.327	0.322	1.5	66	0.00
59		Diethylphthalate	1.386	1.217	12.2	62	0.00
60		4-Chlorophenyl-phenylether	0.668	0.585	12.4	61	0.00
61		4-Nitroaniline	0.312	0.365	-17.0	80	0.00
62		Azobenzene	1.315	1.182	10.1	62	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
64		4,6-Dinitro-2-methylphenol	0.075	0.078	-4.0	69	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.616	7.2	62	0.00
66		4-Bromophenyl-phenylether	0.234	0.197	15.8	59	0.00
67		Hexachlorobenzene	0.251	0.219	12.7	62	0.00
68		Atrazine	0.216	0.177	18.1	61	0.00
69	C	Pentachlorophenol	0.132	0.121	8.3	60	0.00
70		Phenanthrene	1.159	1.101	5.0	67	0.00
71		Anthracene	1.150	1.078	6.3	67	0.00
72		Carbazole	1.094	1.037	5.2	63	0.00
73		Di-n-butylphthalate	1.284	1.054	17.9	54	0.00
74	C	Fluoranthene	1.266	1.220	3.6	66	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
76		Benzidine	0.676	0.659	2.5	83	0.00
77		Pyrene	1.223	0.984	19.5	64	0.00
78	S	Terphenyl-d14	0.856	0.651	23.9	60	0.00
79		Butylbenzylphthalate	0.503	0.407	19.1	60	0.00
80		Benzo(a)anthracene	1.172	1.099	6.2	75	0.00
81		3,3'-Dichlorobenzidine	0.430	0.411	4.4	74	0.00
82		Chrysene	1.101	0.994	9.7	73	0.00
83		Bis(2-ethylhexyl)phthalate	0.807	0.556	31.1#	53	0.00
84	c	Di-n-octyl phthalate	1.348	0.824	38.9#	46#	0.01
85		Indeno(1,2,3-cd)pyrene	1.255	1.108	11.7	69	0.03
86	I	Perylene-d12	1.000	1.000	0.0	75	0.03
87		Benzo(b)fluoranthene	1.163	1.220	-4.9	74	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030315\
Data File : BF077562.D
Acq On : 3 Mar 2015 23:33
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 01:30:44 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 03 23:57:38 2015
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)
88	Benzo(k)fluoranthene	1.140	0.888	22.1	62	0.02
89 C	Benzo(a)pyrene	1.108	1.063	4.1	70	0.03
90	Dibenzo(a,h)anthracene	1.056	1.004	4.9	69	0.03
91	Benzo(g,h,i)perylene	1.066	1.028	3.6	71	0.03

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