

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020421\
 Data File : BF123107.D
 Acq On : 4 Feb 2021 9:51
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 04 10:31:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 29 12:29:00 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	97	0.00
2	1,4-Dioxane	0.645	0.652	-1.1	98	0.01
3	Pyridine	1.725	1.701	1.4	94	0.01
4	n-Nitrosodimethylamine	0.726	0.699	3.7	90	0.00
5 S	2-Fluorophenol	1.297	1.305	-0.6	97	0.00
6	Aniline	2.163	2.105	2.7	92	0.00
7 S	Phenol-d6	1.757	1.729	1.6	94	0.00
8	2-Chlorophenol	1.405	1.434	-2.1	97	0.00
9	Benzaldehyde	0.803	0.840	-4.6	99	0.00
10 C	Phenol	1.743	1.849	-6.1	99	0.00
11	bis(2-Chloroethyl)ether	1.450	1.411	2.7	92	0.00
12	1,3-Dichlorobenzene	1.639	1.652	-0.8	96	0.00
13 C	1,4-Dichlorobenzene	1.707	1.670	2.2	97	0.00
14	1,2-Dichlorobenzene	1.597	1.548	3.1	97	0.00
15	Benzyl Alcohol	1.232	1.220	1.0	91	0.00
16	2,2'-oxybis(1-Chloropropane	2.237	2.139	4.4	91	0.00
17	2-Methylphenol	1.196	1.195	0.1	94	0.00
18	Hexachloroethane	0.598	0.611	-2.2	97	0.00
19 P	n-Nitroso-di-n-propylamine	1.057	1.007	4.7	90	0.00
20	3+4-Methylphenols	1.409	1.461	-3.7	94	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22	Acetophenone	0.521	0.514	1.3	94	0.00
23 S	Nitrobenzene-d5	0.397	0.414	-4.3	97	0.00
24	Nitrobenzene	0.409	0.416	-1.7	94	0.00
25	Isophorone	0.728	0.711	2.3	92	0.00
26 C	2-Nitrophenol	0.169	0.197	-16.6	99	0.00
27	2,4-Dimethylphenol	0.306	0.304	0.7	93	0.00
28	bis(2-Chloroethoxy)methane	0.497	0.485	2.4	92	0.00
29 C	2,4-Dichlorophenol	0.319	0.330	-3.4	96	0.00
30	1,2,4-Trichlorobenzene	0.360	0.369	-2.5	97	0.00
31	Naphthalene	1.097	1.100	-0.3	96	0.00
32	Benzoic acid	0.224	0.243	-8.5	95	0.00
33	4-Chloroaniline	0.487	0.479	1.6	93	0.00
34 C	Hexachlorobutadiene	0.208	0.220	-5.8	99	0.00
35	Caprolactam	0.108	0.107	0.9	91	0.00
36 C	4-Chloro-3-methylphenol	0.332	0.347	-4.5	96	0.00
37	2-Methylnaphthalene	0.728	0.723	0.7	94	0.00
38	1-Methylnaphthalene	0.690	0.686	0.6	94	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
40	1,2,4,5-Tetrachlorobenzene	0.622	0.657	-5.6	98	0.00
41 P	Hexachlorocyclopentadiene	0.295	0.312	-5.8	92	0.00
42 S	2,4,6-Tribromophenol	0.217	0.241	-11.1	100	0.00
43 C	2,4,6-Trichlorophenol	0.427	0.459	-7.5	95	0.00
44	2,4,5-Trichlorophenol	0.446	0.469	-5.2	94	0.00
45 S	2-Fluorobiphenyl	1.345	1.369	-1.8	96	0.00
46	1,1'-Biphenyl	1.640	1.675	-2.1	94	0.00
47	2-Chloronaphthalene	1.373	1.399	-1.9	94	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020421\
 Data File : BF123107.D
 Acq On : 4 Feb 2021 9:51
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 04 10:31:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 29 12:29:00 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	0.416	0.454	-9.1	95	0.00
49	Acenaphthylene	2.010	2.025	-0.7	93	0.00
50	Dimethylphthalate	1.569	1.592	-1.5	95	0.00
51	2,6-Dinitrotoluene	0.334	0.358	-7.2	94	0.00
52 C	Acenaphthene	1.291	1.299	-0.6	93	0.00
53	3-Nitroaniline	0.395	0.412	-4.3	93	0.00
54 P	2,4-Dinitrophenol	0.126	0.153	-21.4	103	0.00
55	Dibenzofuran	1.879	1.827	2.8	93	0.00
56 P	4-Nitrophenol	0.286	0.298	-4.2	92	0.00
57	2,4-Dinitrotoluene	0.435	0.474	-9.0	95	0.00
58	Fluorene	1.501	1.376	8.3	94	0.00
59	2,3,4,6-Tetrachlorophenol	0.378	0.405	-7.1	97	0.00
60	Diethylphthalate	1.543	1.562	-1.2	94	0.00
61	4-Chlorophenyl-phenylether	0.705	0.704	0.1	97	0.00
62	4-Nitroaniline	0.397	0.406	-2.3	91	0.00
63	Azobenzene	1.510	1.472	2.5	90	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
65	4,6-Dinitro-2-methylphenol	0.101	0.116	-14.9	101	0.00
66 c	n-Nitrosodiphenylamine	0.721	0.710	1.5	92	0.00
67	4-Bromophenyl-phenylether	0.255	0.263	-3.1	95	0.00
68	Hexachlorobenzene	0.273	0.288	-5.5	97	0.00
69	Atrazine	0.199	0.207	-4.0	98	0.00
70 C	Pentachlorophenol	0.148	0.164	-10.8	97	0.00
71	Phenanthrene	1.242	1.186	4.5	94	0.00
72	Anthracene	1.192	1.178	1.2	94	0.00
73	Carbazole	1.106	1.079	2.4	92	0.00
74	Di-n-butylphthalate	1.311	1.311	0.0	94	0.00
75 C	Fluoranthene	1.241	1.209	2.6	94	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	88	0.00
77	Benzydine	0.453	0.464	-2.4	74	0.00
78	Pyrene	1.510	1.606	-6.4	92	0.00
79 S	Terphenyl-d14	1.018	1.148	-12.8	96	0.00
80	Butylbenzylphthalate	0.650	0.739	-13.7	94	0.00
81	Benzo(a)anthracene	1.354	1.403	-3.6	91	0.00
82	3,3'-Dichlorobenzidine	0.426	0.468	-9.9	90	0.00
83	Chrysene	1.355	1.372	-1.3	89	0.00
84	Bis(2-ethylhexyl)phthalate	0.863	0.971	-12.5	95	0.00
85 c	Di-n-octyl phthalate	1.450	1.585	-9.3	89	0.00
86 I	Perylene-d12	1.000	1.000	0.0	79	0.00
87	Indeno(1,2,3-cd)pyrene	1.332	1.367	-2.6	77	0.00
88	Benzo(b)fluoranthene	1.444	1.452	-0.6	76	0.00
89	Benzo(k)fluoranthene	1.341	1.494	-11.4	89	0.00
90 C	Benzo(a)pyrene	1.268	1.336	-5.4	80	0.00
91	Dibenzo(a,h)anthracene	1.096	1.133	-3.4	78	0.00
92	Benzo(g,h,i)perylene	1.062	1.091	-2.7	78	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020421\
Data File : BF123107.D
Acq On : 4 Feb 2021 9:51
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 04 10:31:40 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 29 12:29:00 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 = 0