

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021622\
 Data File : BF126920.D
 Acq On : 16 Feb 2022 19:02
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 17 01:09:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 16 14:55:48 2022
 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	103	0.00
2	1,4-Dioxane	0.592	0.569	3.9	102	0.00
3	Pyridine	1.553	1.491	4.0	100	0.00
4	n-Nitrosodimethylamine	0.761	0.731	3.9	98	0.00
5 S	2-Fluorophenol	1.294	1.218	5.9	99	0.00
6	Aniline	2.069	2.027	2.0	100	0.00
7 S	Phenol-d6	1.746	1.674	4.1	100	0.00
8	2-Chlorophenol	1.388	1.331	4.1	100	0.00
9	Benzaldehyde	1.063	0.942	11.4	102	0.00
10 C	Phenol	1.764	1.683	4.6	101	0.00
11	bis(2-Chloroethyl)ether	1.384	1.337	3.4	101	0.00
12	1,3-Dichlorobenzene	1.498	1.425	4.9	101	0.00
13 C	1,4-Dichlorobenzene	1.518	1.448	4.6	101	0.00
14	1,2-Dichlorobenzene	1.448	1.375	5.0	100	0.00
15	Benzyl Alcohol	1.471	1.436	2.4	100	0.00
16	2,2'-oxybis(1-Chloropropane	2.179	2.065	5.2	98	0.00
17	2-Methylphenol	1.201	1.169	2.7	101	0.00
18	Hexachloroethane	0.582	0.560	3.8	101	0.00
19 P	n-Nitroso-di-n-propylamine	1.125	1.078	4.2	100	0.00
20	3+4-Methylphenols	1.560	1.513	3.0	100	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22	Acetophenone	0.527	0.498	5.5	100	0.00
23 S	Nitrobenzene-d5	0.443	0.426	3.8	101	0.00
24	Nitrobenzene	0.419	0.401	4.3	100	0.00
25	Isophorone	0.733	0.702	4.2	99	0.00
26 C	2-Nitrophenol	0.178	0.178	0.0	101	0.00
27	2,4-Dimethylphenol	0.293	0.281	4.1	101	0.00
28	bis(2-Chloroethoxy)methane	0.445	0.428	3.8	101	0.00
29 C	2,4-Dichlorophenol	0.275	0.269	2.2	101	0.00
30	1,2,4-Trichlorobenzene	0.308	0.292	5.2	100	0.00
31	Naphthalene	1.044	1.000	4.2	101	0.00
32	Benzoic acid	0.208	0.213	-2.4	98	0.00
33	4-Chloroaniline	0.411	0.393	4.4	100	0.00
34 C	Hexachlorobutadiene	0.180	0.174	3.3	101	0.00
35	Caprolactam	0.096	0.094	2.1	100	0.00
36 C	4-Chloro-3-methylphenol	0.352	0.342	2.8	99	0.00
37	2-Methylnaphthalene	0.677	0.650	4.0	100	0.00
38	1-Methylnaphthalene	0.649	0.628	3.2	102	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
40	1,2,4,5-Tetrachlorobenzene	0.526	0.516	1.9	103	0.00
41 P	Hexachlorocyclopentadiene	0.286	0.295	-3.1	101	0.00
42 S	2,4,6-Tribromophenol	0.230	0.228	0.9	99	0.00
43 C	2,4,6-Trichlorophenol	0.386	0.377	2.3	100	0.00
44	2,4,5-Trichlorophenol	0.406	0.391	3.7	98	0.00
45 S	2-Fluorobiphenyl	1.359	1.288	5.2	100	0.00
46	1,1'-Biphenyl	1.585	1.529	3.5	101	0.00
47	2-Chloronaphthalene	1.175	1.126	4.2	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021622\
 Data File : BF126920.D
 Acq On : 16 Feb 2022 19:02
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 17 01:09:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 16 14:55:48 2022
 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	0.464	0.455	1.9	100	0.00
49	Acenaphthylene	1.909	1.860	2.6	100	0.00
50	Dimethylphthalate	1.430	1.373	4.0	99	0.00
51	2,6-Dinitrotoluene	0.291	0.292	-0.3	101	0.00
52 C	Acenaphthene	1.241	1.207	2.7	99	0.00
53	3-Nitroaniline	0.356	0.351	1.4	99	0.00
54 P	2,4-Dinitrophenol	0.154	0.163	-5.8	100	0.00
55	Dibenzofuran	1.707	1.664	2.5	101	0.00
56 P	4-Nitrophenol	0.263	0.259	1.5	97	0.00
57	2,4-Dinitrotoluene	0.393	0.394	-0.3	100	0.00
58	Fluorene	1.330	1.271	4.4	99	0.00
59	2,3,4,6-Tetrachlorophenol	0.322	0.319	0.9	101	0.00
60	Diethylphthalate	1.494	1.448	3.1	99	0.00
61	4-Chlorophenyl-phenylether	0.627	0.603	3.8	99	0.00
62	4-Nitroaniline	0.351	0.338	3.7	96	0.00
63	Azobenzene	1.663	1.600	3.8	99	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
65	4,6-Dinitro-2-methylphenol	0.118	0.122	-3.4	98	0.00
66 c	n-Nitrosodiphenylamine	0.656	0.629	4.1	98	0.00
67	4-Bromophenyl-phenylether	0.223	0.217	2.7	100	0.00
68	Hexachlorobenzene	0.241	0.231	4.1	100	0.00
69	Atrazine	0.203	0.197	3.0	100	0.00
70 C	Pentachlorophenol	0.131	0.134	-2.3	97	0.00
71	Phenanthrene	1.119	1.062	5.1	100	0.00
72	Anthracene	1.114	1.065	4.4	99	0.00
73	Carbazole	1.022	0.967	5.4	98	0.00
74	Di-n-butylphthalate	1.314	1.256	4.4	98	0.00
75 C	Fluoranthene	1.063	1.003	5.6	98	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
77	Benzydine	0.544	0.560	-2.9	98	0.00
78	Pyrene	1.738	1.743	-0.3	98	0.00
79 S	Terphenyl-d14	1.361	1.357	0.3	100	0.00
80	Butylbenzylphthalate	0.823	0.827	-0.5	98	0.00
81	Benzo(a)anthracene	1.348	1.329	1.4	98	0.00
82	3,3'-Dichlorobenzidine	0.425	0.422	0.7	97	0.00
83	Chrysene	1.268	1.244	1.9	98	0.00
84	Bis(2-ethylhexyl)phthalate	1.080	1.073	0.6	98	0.00
85 c	Di-n-octyl phthalate	1.555	1.520	2.3	99	0.00
86 I	Perylene-d12	1.000	1.000	0.0	100	0.00
87	Indeno(1,2,3-cd)pyrene	1.311	1.366	-4.2	100	0.00
88	Benzo(b)fluoranthene	1.343	1.322	1.6	102	0.00
89	Benzo(k)fluoranthene	1.296	1.263	2.5	98	0.00
90 C	Benzo(a)pyrene	1.049	1.041	0.8	100	0.00
91	Dibenzo(a,h)anthracene	1.085	1.128	-4.0	100	0.00
92	Benzo(g,h,i)perylene	1.069	1.122	-5.0	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021622\
Data File : BF126920.D
Acq On : 16 Feb 2022 19:02
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 17 01:09:00 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 16 14:55:48 2022
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