

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060222\
 Data File : BF128646.D
 Acq On : 02 Jun 2022 19:08
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 03 03:27:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 02 17:07:23 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	87	0.00
2	1,4-Dioxane	40.000	39.389	1.5	81	-0.04
3	Pyridine	40.000	38.526	3.7	80	-0.04
4	n-Nitrosodimethylamine	40.000	38.739	3.2	80	-0.04
5 S	2-Fluorophenol	80.000	77.179	3.5	82	0.00
6	Aniline	40.000	37.950	5.1	83	0.00
7 S	Phenol-d6	80.000	76.659	4.2	83	0.00
8	2-Chlorophenol	40.000	38.264	4.3	83	0.00
9	Benzaldehyde	40.000	33.654	15.9	90	0.00
10 C	Phenol	40.000	38.406	4.0	82	0.00
11	bis(2-Chloroethyl)ether	40.000	38.518	3.7	83	0.00
12	1,3-Dichlorobenzene	40.000	39.136	2.2	84	0.00
13 C	1,4-Dichlorobenzene	40.000	38.916	2.7	85	0.00
14	1,2-Dichlorobenzene	40.000	39.325	1.7	87	0.00
15	Benzyl Alcohol	40.000	39.265	1.8	86	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	39.635	0.9	88	0.00
17	2-Methylphenol	40.000	39.587	1.0	87	0.00
18	Hexachloroethane	40.000	41.551	-3.9	89	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	39.973	0.1	86	0.00
20	3+4-Methylphenols	40.000	39.075	2.3	84	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	87	0.00
22	Acetophenone	40.000	40.148	-0.4	85	0.00
23 S	Nitrobenzene-d5	80.000	80.527	-0.7	86	0.00
24	Nitrobenzene	40.000	40.328	-0.8	85	0.00
25	Isophorone	40.000	39.398	1.5	84	0.00
26 C	2-Nitrophenol	40.000	40.726	-1.8	83	0.00
27	2,4-Dimethylphenol	40.000	39.600	1.0	81	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.712	0.7	83	0.00
29 C	2,4-Dichlorophenol	40.000	40.411	-1.0	84	0.00
30	1,2,4-Trichlorobenzene	40.000	40.160	-0.4	84	0.00
31	Naphthalene	40.000	39.861	0.3	87	0.00
32	Benzoic acid	40.000	41.905	-4.8	85	-0.02
33	4-Chloroaniline	40.000	38.899	2.8	86	0.00
34 C	Hexachlorobutadiene	40.000	40.302	-0.8	89	0.00
35	Caprolactam	40.000	39.076	2.3	86	-0.01
36 C	4-Chloro-3-methylphenol	40.000	40.331	-0.8	89	0.00
37	2-Methylnaphthalene	40.000	39.877	0.3	88	0.00
38	1-Methylnaphthalene	40.000	39.486	1.3	89	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	87	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	41.991	-5.0	88	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.471	-3.7	85	0.00
42 S	2,4,6-Tribromophenol	80.000	84.252	-5.3	86	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.877	-2.2	87	0.00
44	2,4,5-Trichlorophenol	40.000	41.036	-2.6	87	0.00
45 S	2-Fluorobiphenyl	80.000	83.322	-4.2	89	0.00
46	1,1'-Biphenyl	40.000	40.822	-2.1	88	0.00
47	2-Chloronaphthalene	40.000	41.344	-3.4	89	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060222\
 Data File : BF128646.D
 Acq On : 02 Jun 2022 19:08
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 03 03:27:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 02 17:07:23 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	43.064	-7.7	91	0.00
49	Acenaphthylene	40.000	41.047	-2.6	88	0.00
50	Dimethylphthalate	40.000	41.011	-2.5	87	0.00
51	2,6-Dinitrotoluene	40.000	42.379	-5.9	88	0.00
52 C	Acenaphthene	40.000	41.673	-4.2	87	0.00
53	3-Nitroaniline	40.000	41.240	-3.1	85	0.00
54 P	2,4-Dinitrophenol	40.000	41.095	-2.7	84	0.00
55	Dibenzofuran	40.000	41.177	-2.9	88	0.00
56 P	4-Nitrophenol	40.000	40.777	-1.9	88	0.00
57	2,4-Dinitrotoluene	40.000	42.216	-5.5	86	0.00
58	Fluorene	40.000	40.957	-2.4	86	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	42.014	-5.0	87	0.00
60	Diethylphthalate	40.000	42.236	-5.6	90	0.00
61	4-Chlorophenyl-phenylether	40.000	42.235	-5.6	89	0.00
62	4-Nitroaniline	40.000	40.284	-0.7	84	0.00
63	Azobenzene	40.000	41.786	-4.5	88	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	85	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.480	-3.7	84	0.00
66 c	n-Nitrosodiphenylamine	40.000	41.289	-3.2	85	0.00
67	4-Bromophenyl-phenylether	40.000	42.134	-5.3	87	0.00
68	Hexachlorobenzene	40.000	41.280	-3.2	85	0.00
69	Atrazine	40.000	39.868	0.3	81	0.00
70 C	Pentachlorophenol	40.000	40.677	-1.7	82	0.00
71	Phenanthrene	40.000	40.274	-0.7	85	0.00
72	Anthracene	40.000	40.509	-1.3	85	0.00
73	Carbazole	40.000	39.422	1.4	83	0.00
74	Di-n-butylphthalate	40.000	41.094	-2.7	84	0.00
75 C	Fluoranthene	40.000	40.268	-0.7	82	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	76	0.00
77	Benzydine	40.000	46.806	-17.0	96	0.00
78	Pyrene	40.000	41.074	-2.7	81	0.00
79 S	Terphenyl-d14	80.000	83.992	-5.0	79	0.00
80	Butylbenzylphthalate	40.000	41.683	-4.2	77	0.00
81	Benzo(a)anthracene	40.000	40.914	-2.3	77	0.00
82	3,3'-Dichlorobenzidine	40.000	43.082	-7.7	82	0.00
83	Chrysene	40.000	41.329	-3.3	79	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	42.432	-6.1	79	0.00
85 c	Di-n-octyl phthalate	40.000	44.430	-11.1	82	0.00
86 I	Perylene-d12	20.000	20.000	0.0	88	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	45.978	-14.9	112	0.00
88	Benzo(b)fluoranthene	40.000	41.283	-3.2	88	0.00
89	Benzo(k)fluoranthene	40.000	38.653	3.4	83	0.00
90 C	Benzo(a)pyrene	40.000	41.588	-4.0	90	0.00
91	Dibenzo(a,h)anthracene	40.000	45.896	-14.7	109	0.00
92	Benzo(g,h,i)perylene	40.000	46.307	-15.8	113	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060222\
Data File : BF128646.D
Acq On : 02 Jun 2022 19:08
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 03 03:27:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 02 17:07:23 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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

(#) = Out of Range

SPCC's out = 0 CCC's out = 0