

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070122\
 Data File : BF129092.D
 Acq On : 01 Jul 2022 15:01
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 02 01:25:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 29 17:05:31 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	82	0.00
2	1,4-Dioxane	40.000	39.082	2.3	81	-0.04
3	Pyridine	40.000	42.130	-5.3	85	-0.06
4	n-Nitrosodimethylamine	40.000	37.866	5.3	78	-0.04
5 S	2-Fluorophenol	80.000	78.644	1.7	82	-0.01
6	Aniline	40.000	40.256	-0.6	84	0.00
7 S	Phenol-d6	80.000	79.357	0.8	83	0.00
8	2-Chlorophenol	40.000	40.126	-0.3	84	0.00
9	Benzaldehyde	40.000	44.012	-10.0	93	0.00
10 C	Phenol	40.000	39.128	2.2	82	0.00
11	bis(2-Chloroethyl)ether	40.000	39.118	2.2	82	0.00
12	1,3-Dichlorobenzene	40.000	39.250	1.9	83	0.00
13 C	1,4-Dichlorobenzene	40.000	39.636	0.9	83	0.00
14	1,2-Dichlorobenzene	40.000	39.794	0.5	83	0.00
15	Benzyl Alcohol	40.000	39.420	1.4	81	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	38.902	2.7	82	0.00
17	2-Methylphenol	40.000	39.259	1.9	82	0.00
18	Hexachloroethane	40.000	39.605	1.0	83	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.891	2.8	82	0.00
20	3+4-Methylphenols	40.000	40.239	-0.6	83	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	83	0.00
22	Acetophenone	40.000	38.911	2.7	82	0.00
23 S	Nitrobenzene-d5	80.000	83.099	-3.9	85	0.00
24	Nitrobenzene	40.000	40.607	-1.5	84	0.00
25	Isophorone	40.000	39.476	1.3	83	-0.01
26 C	2-Nitrophenol	40.000	46.905	-17.3	92	0.00
27	2,4-Dimethylphenol	40.000	40.343	-0.9	84	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.435	1.4	82	0.00
29 C	2,4-Dichlorophenol	40.000	40.289	-0.7	83	0.00
30	1,2,4-Trichlorobenzene	40.000	39.240	1.9	83	0.00
31	Naphthalene	40.000	40.214	-0.5	83	0.00
32	Benzoic acid	40.000	41.268	-3.2	86	-0.01
33	4-Chloroaniline	40.000	40.739	-1.8	85	0.00
34 C	Hexachlorobutadiene	40.000	39.140	2.1	83	0.00
35	Caprolactam	40.000	40.941	-2.4	86	-0.02
36 C	4-Chloro-3-methylphenol	40.000	40.292	-0.7	84	0.00
37	2-Methylnaphthalene	40.000	40.193	-0.5	84	0.00
38	1-Methylnaphthalene	40.000	40.029	-0.1	83	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	84	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.855	0.4	84	0.00
41 P	Hexachlorocyclopentadiene	40.000	40.780	-2.0	83	0.00
42 S	2,4,6-Tribromophenol	80.000	81.992	-2.5	86	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.330	-0.8	83	0.00
44	2,4,5-Trichlorophenol	40.000	39.946	0.1	83	0.00
45 S	2-Fluorobiphenyl	80.000	75.199	6.0	86	0.00
46	1,1'-Biphenyl	40.000	40.236	-0.6	85	0.00
47	2-Chloronaphthalene	40.000	39.255	1.9	83	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070122\
 Data File : BF129092.D
 Acq On : 01 Jul 2022 15:01
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 02 01:25:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 29 17:05:31 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	44.325	-10.8	89	0.00
49	Acenaphthylene	40.000	40.297	-0.7	84	0.00
50	Dimethylphthalate	40.000	39.798	0.5	84	0.00
51	2,6-Dinitrotoluene	40.000	42.698	-6.7	87	0.00
52 C	Acenaphthene	40.000	40.416	-1.0	85	0.00
53	3-Nitroaniline	40.000	42.813	-7.0	87	0.00
54 P	2,4-Dinitrophenol	40.000	47.469	-18.7	113	0.00
55	Dibenzofuran	40.000	40.020	-0.1	84	0.00
56 P	4-Nitrophenol	40.000	41.483	-3.7	83	0.00
57	2,4-Dinitrotoluene	40.000	45.834	-14.6	91	0.00
58	Fluorene	40.000	39.491	1.3	84	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.302	-0.8	84	0.00
60	Diethylphthalate	40.000	39.953	0.1	85	0.00
61	4-Chlorophenyl-phenylether	40.000	39.617	1.0	84	0.00
62	4-Nitroaniline	40.000	42.186	-5.5	86	0.00
63	Azobenzene	40.000	39.358	1.6	82	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	85	0.00
65	4,6-Dinitro-2-methylphenol	40.000	51.716	-29.3#	107	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.800	0.5	84	0.00
67	4-Bromophenyl-phenylether	40.000	40.113	-0.3	84	0.00
68	Hexachlorobenzene	40.000	39.309	1.7	85	0.00
69	Atrazine	40.000	41.275	-3.2	86	0.00
70 C	Pentachlorophenol	40.000	41.123	-2.8	83	0.00
71	Phenanthrene	40.000	39.433	1.4	84	0.00
72	Anthracene	40.000	39.998	0.0	85	0.00
73	Carbazole	40.000	39.848	0.4	86	0.00
74	Di-n-butylphthalate	40.000	41.666	-4.2	88	0.00
75 C	Fluoranthene	40.000	40.426	-1.1	87	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	87	0.00
77	Benzydine	40.000	50.121	-25.3#	103	0.00
78	Pyrene	40.000	40.244	-0.6	87	0.00
79 S	Terphenyl-d14	80.000	79.603	0.5	89	0.00
80	Butylbenzylphthalate	40.000	42.597	-6.5	91	0.00
81	Benzo(a)anthracene	40.000	39.745	0.6	88	0.00
82	3,3'-Dichlorobenzidine	40.000	43.198	-8.0	98	0.00
83	Chrysene	40.000	39.358	1.6	87	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	42.176	-5.4	91	0.00
85 c	Di-n-octyl phthalate	40.000	42.794	-7.0	94	-0.01
86 I	Perylene-d12	20.000	20.000	0.0	89	-0.01
87	Indeno(1,2,3-cd)pyrene	40.000	38.847	2.9	91	-0.01
88	Benzo(b)fluoranthene	40.000	41.069	-2.7	92	-0.01
89	Benzo(k)fluoranthene	40.000	38.991	2.5	87	0.00
90 C	Benzo(a)pyrene	40.000	39.542	1.1	89	-0.01
91	Dibenzo(a,h)anthracene	40.000	39.419	1.5	91	-0.01
92	Benzo(g,h,i)perylene	40.000	37.952	5.1	89	-0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070122\
Data File : BF129092.D
Acq On : 01 Jul 2022 15:01
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jul 02 01:25:11 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 29 17:05:31 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