

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080122\
 Data File : BF129476.D
 Acq On : 01 Aug 2022 16:37
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 02 01:16:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 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	81	0.00
2	1,4-Dioxane	40.000	38.657	3.4	79	0.02
3	Pyridine	40.000	39.043	2.4	78	0.02
4	n-Nitrosodimethylamine	40.000	39.162	2.1	79	0.02
5 S	2-Fluorophenol	80.000	78.124	2.3	82	0.01
6	Aniline	40.000	33.909	15.2	70	0.00
7 S	Phenol-d6	80.000	77.287	3.4	81	0.00
8	2-Chlorophenol	40.000	39.397	1.5	81	0.00
9	Benzaldehyde	40.000	35.061	12.3	76	0.00
10 C	Phenol	40.000	39.375	1.6	82	0.01
11	bis(2-Chloroethyl)ether	40.000	39.189	2.0	81	0.00
12	1,3-Dichlorobenzene	40.000	39.565	1.1	82	0.00
13 C	1,4-Dichlorobenzene	40.000	39.517	1.2	82	0.00
14	1,2-Dichlorobenzene	40.000	39.484	1.3	81	0.00
15	Benzyl Alcohol	40.000	40.061	-0.2	81	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	34.875	12.8	72	0.00
17	2-Methylphenol	40.000	38.103	4.7	79	0.00
18	Hexachloroethane	40.000	40.116	-0.3	83	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.229	4.4	81	0.00
20	3+4-Methylphenols	40.000	38.073	4.8	82	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	79	0.00
22	Acetophenone	40.000	40.621	-1.6	85	0.00
23 S	Nitrobenzene-d5	80.000	82.112	-2.6	83	0.00
24	Nitrobenzene	40.000	40.202	-0.5	80	0.00
25	Isophorone	40.000	38.360	4.1	78	0.00
26 C	2-Nitrophenol	40.000	40.772	-1.9	81	0.00
27	2,4-Dimethylphenol	40.000	40.133	-0.3	81	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.591	3.5	79	0.00
29 C	2,4-Dichlorophenol	40.000	39.619	1.0	79	0.00
30	1,2,4-Trichlorobenzene	40.000	40.328	-0.8	82	0.00
31	Naphthalene	40.000	39.722	0.7	81	0.00
32	Benzoic acid	40.000	23.535	41.2#	45	0.01
33	4-Chloroaniline	40.000	37.461	6.3	76	0.00
34 C	Hexachlorobutadiene	40.000	42.086	-5.2	86	0.00
35	Caprolactam	40.000	37.984	5.0	76	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.025	-0.1	81	0.00
37	2-Methylnaphthalene	40.000	39.336	1.7	80	0.00
38	1-Methylnaphthalene	40.000	39.559	1.1	82	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	78	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	41.084	-2.7	82	0.00
41 P	Hexachlorocyclopentadiene	40.000	42.604	-6.5	77	0.00
42 S	2,4,6-Tribromophenol	80.000	79.738	0.3	79	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.588	-1.5	79	0.00
44	2,4,5-Trichlorophenol	40.000	40.134	-0.3	79	0.00
45 S	2-Fluorobiphenyl	80.000	75.288	5.9	83	0.00
46	1,1'-Biphenyl	40.000	40.412	-1.0	80	0.00
47	2-Chloronaphthalene	40.000	40.196	-0.5	80	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080122\
 Data File : BF129476.D
 Acq On : 01 Aug 2022 16:37
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 02 01:16:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 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	40.846	-2.1	79	0.00
49	Acenaphthylene	40.000	40.086	-0.2	80	0.00
50	Dimethylphthalate	40.000	39.366	1.6	79	0.00
51	2,6-Dinitrotoluene	40.000	40.445	-1.1	79	0.00
52 C	Acenaphthene	40.000	39.657	0.9	78	0.00
53	3-Nitroaniline	40.000	38.050	4.9	74	0.00
54 P	2,4-Dinitrophenol	40.000	39.002	2.5	72	0.00
55	Dibenzofuran	40.000	39.752	0.6	79	0.00
56 P	4-Nitrophenol	40.000	36.212	9.5	70	0.01
57	2,4-Dinitrotoluene	40.000	40.086	-0.2	76	0.00
58	Fluorene	40.000	39.926	0.2	81	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	38.758	3.1	75	0.00
60	Diethylphthalate	40.000	38.998	2.5	78	0.00
61	4-Chlorophenyl-phenylether	40.000	39.765	0.6	80	0.00
62	4-Nitroaniline	40.000	38.313	4.2	75	0.00
63	Azobenzene	40.000	38.187	4.5	76	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	76	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.326	-3.3	74	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.884	0.3	77	0.00
67	4-Bromophenyl-phenylether	40.000	40.648	-1.6	78	0.00
68	Hexachlorobenzene	40.000	40.777	-1.9	79	0.00
69	Atrazine	40.000	38.810	3.0	75	0.00
70 C	Pentachlorophenol	40.000	33.610	16.0	61	0.00
71	Phenanthrene	40.000	39.767	0.6	78	0.00
72	Anthracene	40.000	39.760	0.6	79	0.00
73	Carbazole	40.000	39.242	1.9	77	0.00
74	Di-n-butylphthalate	40.000	39.407	1.5	77	0.00
75 C	Fluoranthene	40.000	40.035	-0.1	78	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	75	0.00
77	Benzydine	40.000	35.012	12.5	63	0.00
78	Pyrene	40.000	41.704	-4.3	79	0.00
79 S	Terphenyl-d14	80.000	83.690	-4.6	81	0.00
80	Butylbenzylphthalate	40.000	40.943	-2.4	75	0.00
81	Benzo(a)anthracene	40.000	39.869	0.3	76	0.00
82	3,3'-Dichlorobenzidine	40.000	37.602	6.0	71	0.00
83	Chrysene	40.000	39.878	0.3	77	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.854	2.9	73	0.00
85 c	Di-n-octyl phthalate	40.000	39.067	2.3	75	0.00
86 I	Perylene-d12	20.000	20.000	0.0	78	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	41.594	-4.0	76	0.00
88	Benzo(b)fluoranthene	40.000	39.436	1.4	81	0.00
89	Benzo(k)fluoranthene	40.000	38.469	3.8	75	0.00
90 C	Benzo(a)pyrene	40.000	39.220	2.0	77	0.00
91	Dibenzo(a,h)anthracene	40.000	41.328	-3.3	75	0.00
92	Benzo(g,h,i)perylene	40.000	41.616	-4.0	74	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080122\
Data File : BF129476.D
Acq On : 01 Aug 2022 16:37
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 02 01:16:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 02 01:11:25 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