

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072222\
 Data File : BF129306.D
 Acq On : 22 Jul 2022 09:14
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 22 10:56:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 20 13:42:18 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	60	0.00
2	1,4-Dioxane	40.000	38.048	4.9	59	0.00
3	Pyridine	40.000	36.469	8.8	54	-0.01
4	n-Nitrosodimethylamine	40.000	43.103	-7.8	65	0.00
5 S	2-Fluorophenol	80.000	79.483	0.6	63	0.00
6	Aniline	40.000	39.019	2.5	61	0.00
7 S	Phenol-d6	80.000	78.637	1.7	62	0.00
8	2-Chlorophenol	40.000	40.162	-0.4	63	0.00
9	Benzaldehyde	40.000	38.869	2.8	61	0.00
10 C	Phenol	40.000	39.691	0.8	62	0.00
11	bis(2-Chloroethyl)ether	40.000	38.407	4.0	60	0.00
12	1,3-Dichlorobenzene	40.000	40.385	-1.0	63	0.00
13 C	1,4-Dichlorobenzene	40.000	40.263	-0.7	63	0.00
14	1,2-Dichlorobenzene	40.000	41.117	-2.8	65	0.00
15	Benzyl Alcohol	40.000	43.055	-7.6	66	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	37.735	5.7	59	0.00
17	2-Methylphenol	40.000	39.758	0.6	62	0.00
18	Hexachloroethane	40.000	41.592	-4.0	65	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	40.994	-2.5	66	0.00
20	3+4-Methylphenols	40.000	39.250	1.9	64	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	62	0.00
22	Acetophenone	40.000	40.721	-1.8	66	0.00
23 S	Nitrobenzene-d5	80.000	83.181	-4.0	66	0.00
24	Nitrobenzene	40.000	40.761	-1.9	65	0.00
25	Isophorone	40.000	39.788	0.5	64	0.00
26 C	2-Nitrophenol	40.000	39.875	0.3	62	0.00
27	2,4-Dimethylphenol	40.000	39.430	1.4	62	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.747	3.1	62	0.00
29 C	2,4-Dichlorophenol	40.000	39.813	0.5	63	0.00
30	1,2,4-Trichlorobenzene	40.000	39.945	0.1	64	0.00
31	Naphthalene	40.000	40.286	-0.7	65	0.00
32	Benzoic acid	40.000	39.741	0.6	60	0.00
33	4-Chloroaniline	40.000	39.097	2.3	62	0.00
34 C	Hexachlorobutadiene	40.000	43.538	-8.8	69	0.00
35	Caprolactam	40.000	40.416	-1.0	64	0.00
36 C	4-Chloro-3-methylphenol	40.000	41.011	-2.5	65	0.00
37	2-Methylnaphthalene	40.000	40.080	-0.2	64	0.00
38	1-Methylnaphthalene	40.000	40.099	-0.2	64	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	62	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	41.331	-3.3	66	0.00
41 P	Hexachlorocyclopentadiene	40.000	47.312	-18.3	69	0.00
42 S	2,4,6-Tribromophenol	80.000	87.830	-9.8	71	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.444	-1.1	65	0.00
44	2,4,5-Trichlorophenol	40.000	41.124	-2.8	65	0.00
45 S	2-Fluorobiphenyl	80.000	78.250	2.2	70	0.00
46	1,1'-Biphenyl	40.000	40.075	-0.2	64	0.00
47	2-Chloronaphthalene	40.000	40.337	-0.8	65	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072222\
 Data File : BF129306.D
 Acq On : 22 Jul 2022 09:14
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 22 10:56:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 20 13:42:18 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	41.545	-3.9	64	0.00
49	Acenaphthylene	40.000	40.609	-1.5	65	0.00
50	Dimethylphthalate	40.000	40.470	-1.2	65	0.00
51	2,6-Dinitrotoluene	40.000	40.939	-2.3	65	0.00
52 C	Acenaphthene	40.000	41.033	-2.6	66	0.00
53	3-Nitroaniline	40.000	38.962	2.6	61	0.00
54 P	2,4-Dinitrophenol	40.000	42.219	-5.5	65	0.00
55	Dibenzofuran	40.000	41.294	-3.2	66	0.00
56 P	4-Nitrophenol	40.000	41.171	-2.9	63	0.00
57	2,4-Dinitrotoluene	40.000	42.384	-6.0	66	0.00
58	Fluorene	40.000	41.400	-3.5	67	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	41.537	-3.8	66	0.00
60	Diethylphthalate	40.000	42.258	-5.6	68	0.00
61	4-Chlorophenyl-phenylether	40.000	41.268	-3.2	67	0.00
62	4-Nitroaniline	40.000	40.197	-0.5	63	0.00
63	Azobenzene	40.000	40.976	-2.4	66	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	64	0.00
65	4,6-Dinitro-2-methylphenol	40.000	43.552	-8.9	65	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.865	0.3	65	0.00
67	4-Bromophenyl-phenylether	40.000	41.105	-2.8	67	0.00
68	Hexachlorobenzene	40.000	41.715	-4.3	69	0.00
69	Atrazine	40.000	41.333	-3.3	67	0.00
70 C	Pentachlorophenol	40.000	42.199	-5.5	65	0.00
71	Phenanthrene	40.000	40.217	-0.5	66	0.00
72	Anthracene	40.000	40.597	-1.5	67	0.00
73	Carbazole	40.000	40.108	-0.3	66	0.00
74	Di-n-butylphthalate	40.000	42.179	-5.4	69	0.00
75 C	Fluoranthene	40.000	40.695	-1.7	67	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	67	0.00
77	Benzydine	40.000	41.028	-2.6	66	0.00
78	Pyrene	40.000	39.459	1.4	66	0.00
79 S	Terphenyl-d14	80.000	83.517	-4.4	73	0.00
80	Butylbenzylphthalate	40.000	40.550	-1.4	67	0.00
81	Benzo(a)anthracene	40.000	39.291	1.8	67	0.00
82	3,3'-Dichlorobenzidine	40.000	39.495	1.3	66	0.00
83	Chrysene	40.000	39.207	2.0	68	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.722	-1.8	69	0.00
85 c	Di-n-octyl phthalate	40.000	39.866	0.3	67	0.00
86 I	Perylene-d12	20.000	20.000	0.0	63	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	42.945	-7.4	66	-0.01
88	Benzo(b)fluoranthene	40.000	42.248	-5.6	67	0.00
89	Benzo(k)fluoranthene	40.000	38.923	2.7	65	-0.01
90 C	Benzo(a)pyrene	40.000	39.459	1.4	64	0.00
91	Dibenzo(a,h)anthracene	40.000	43.146	-7.9	66	-0.01
92	Benzo(g,h,i)perylene	40.000	42.942	-7.4	65	-0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072222\
Data File : BF129306.D
Acq On : 22 Jul 2022 09:14
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jul 22 10:56:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 20 13:42:18 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