

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060122\
 Data File : BP010644.D
 Acq On : 01 Jun 2022 10:57
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 02 05:00:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 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	107	0.00
2	1,4-Dioxane	40.000	34.283	14.3	93	0.00
3	Pyridine	40.000	38.345	4.1	101	0.00
4	n-Nitrosodimethylamine	40.000	34.911	12.7	96	0.00
5 S	2-Fluorophenol	80.000	80.617	-0.8	107	0.00
6	Aniline	40.000	42.330	-5.8	115	0.00
7 S	Phenol-d6	80.000	83.998	-5.0	113	0.00
8	2-Chlorophenol	40.000	41.354	-3.4	112	0.00
9	Benzaldehyde	40.000	34.673	13.3	105	0.00
10 C	Phenol	40.000	41.515	-3.8	113	0.00
11	bis(2-Chloroethyl)ether	40.000	40.712	-1.8	109	0.00
12	1,3-Dichlorobenzene	40.000	39.618	1.0	108	0.00
13 C	1,4-Dichlorobenzene	40.000	39.728	0.7	108	0.00
14	1,2-Dichlorobenzene	40.000	39.499	1.3	108	0.00
15	Benzyl Alcohol	40.000	43.074	-7.7	115	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	37.752	5.6	104	0.00
17	2-Methylphenol	40.000	42.917	-7.3	116	0.00
18	Hexachloroethane	40.000	38.879	2.8	106	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	42.583	-6.5	119	0.00
20	3+4-Methylphenols	40.000	43.694	-9.2	116	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	115	0.00
22	Acetophenone	40.000	39.998	0.0	116	0.00
23 S	Nitrobenzene-d5	80.000	78.948	1.3	114	0.00
24	Nitrobenzene	40.000	39.279	1.8	112	0.00
25	Isophorone	40.000	42.735	-6.8	123	0.00
26 C	2-Nitrophenol	40.000	44.376	-10.9	125	0.00
27	2,4-Dimethylphenol	40.000	42.590	-6.5	121	0.00
28	bis(2-Chloroethoxy)methane	40.000	40.670	-1.7	117	0.00
29 C	2,4-Dichlorophenol	40.000	42.660	-6.6	119	0.00
30	1,2,4-Trichlorobenzene	40.000	39.441	1.4	113	0.00
31	Naphthalene	40.000	39.852	0.4	115	0.00
32	Benzoic acid	40.000	40.983	-2.5	131	0.02
33	4-Chloroaniline	40.000	43.650	-9.1	125	0.00
34 C	Hexachlorobutadiene	40.000	38.176	4.6	112	0.00
35	Caprolactam	40.000	54.280	-35.7#	149	0.01
36 C	4-Chloro-3-methylphenol	40.000	45.307	-13.3	130	0.00
37	2-Methylnaphthalene	40.000	41.437	-3.6	121	0.00
38	1-Methylnaphthalene	40.000	41.555	-3.9	121	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	128	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	37.520	6.2	120	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.700	-4.3	129	0.00
42 S	2,4,6-Tribromophenol	80.000	90.453	-13.1	142	0.00
43 C	2,4,6-Trichlorophenol	40.000	41.714	-4.3	128	0.00
44	2,4,5-Trichlorophenol	40.000	40.738	-1.8	128	0.00
45 S	2-Fluorobiphenyl	80.000	76.381	4.5	124	0.00
46	1,1'-Biphenyl	40.000	38.293	4.3	124	0.00
47	2-Chloronaphthalene	40.000	38.295	4.3	124	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060122\
 Data File : BP010644.D
 Acq On : 01 Jun 2022 10:57
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 02 05:00:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 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	42.477	-6.2	132	0.00
49	Acenaphthylene	40.000	40.465	-1.2	130	0.00
50	Dimethylphthalate	40.000	41.923	-4.8	135	0.00
51	2,6-Dinitrotoluene	40.000	43.735	-9.3	137	0.00
52 C	Acenaphthene	40.000	39.446	1.4	128	0.00
53	3-Nitroaniline	40.000	47.118	-17.8	144	0.00
54 P	2,4-Dinitrophenol	40.000	45.205	-13.0	156	0.00
55	Dibenzofuran	40.000	40.630	-1.6	131	0.00
56 P	4-Nitrophenol	40.000	43.717	-9.3	141	0.00
57	2,4-Dinitrotoluene	40.000	46.987	-17.5	143	0.00
58	Fluorene	40.000	41.491	-3.7	133	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	43.826	-9.6	138	0.00
60	Diethylphthalate	40.000	44.000	-10.0	139	0.00
61	4-Chlorophenyl-phenylether	40.000	41.242	-3.1	134	0.00
62	4-Nitroaniline	40.000	47.455	-18.6	151	0.00
63	Azobenzene	40.000	41.298	-3.2	131	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	141	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.793	-4.5	156	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.473	3.8	137	0.00
67	4-Bromophenyl-phenylether	40.000	38.480	3.8	138	0.00
68	Hexachlorobenzene	40.000	38.895	2.8	141	0.00
69	Atrazine	40.000	43.635	-9.1	149	0.00
70 C	Pentachlorophenol	40.000	41.042	-2.6	140	0.00
71	Phenanthrene	40.000	39.400	1.5	139	0.00
72	Anthracene	40.000	40.741	-1.9	142	0.00
73	Carbazole	40.000	43.811	-9.5	147	0.00
74	Di-n-butylphthalate	40.000	44.894	-12.2	147	0.00
75 C	Fluoranthene	40.000	43.355	-8.4	145	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	142	0.00
77	Benzydine	40.000	49.584	-24.0	175	0.00
78	Pyrene	40.000	39.710	0.7	145	0.00
79 S	Terphenyl-d14	80.000	79.851	0.2	145	0.00
80	Butylbenzylphthalate	40.000	45.621	-14.1	155	0.00
81	Benzo(a)anthracene	40.000	40.231	-0.6	142	0.00
82	3,3'-Dichlorobenzidine	40.000	43.262	-8.2	144	0.00
83	Chrysene	40.000	39.634	0.9	141	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	45.692	-14.2	150	0.00
85 c	Di-n-octyl phthalate	40.000	45.926	-14.8	151	0.00
86 I	Perylene-d12	20.000	20.000	0.0	135	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	35.483	11.3	119	0.01
88	Benzo(b)fluoranthene	40.000	40.672	-1.7	133	0.00
89	Benzo(k)fluoranthene	40.000	40.458	-1.1	138	0.00
90 C	Benzo(a)pyrene	40.000	40.680	-1.7	134	0.00
91	Dibenzo(a,h)anthracene	40.000	36.026	9.9	121	0.00
92	Benzo(g,h,i)perylene	40.000	34.099	14.8	116	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060122\
Data File : BP010644.D
Acq On : 01 Jun 2022 10:57
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jun 02 05:00:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 28 02:18:04 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