

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011022\
 Data File : BP008744.D
 Acq On : 10 Jan 2022 10:51
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 10 16:32:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 07 07:45:52 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	108	0.00
2	1,4-Dioxane	40.000	35.539	11.2	108	0.00
3	Pyridine	40.000	35.489	11.3	103	0.00
4	n-Nitrosodimethylamine	40.000	35.544	11.1	106	0.00
5 S	2-Fluorophenol	80.000	71.043	11.2	106	0.00
6	Aniline	40.000	35.180	12.1	105	0.00
7 S	Phenol-d6	80.000	70.237	12.2	104	0.00
8	2-Chlorophenol	40.000	35.430	11.4	106	0.00
9	Benzaldehyde	40.000	34.807	13.0	104	0.00
10 C	Phenol	40.000	34.973	12.6	104	0.00
11	bis(2-Chloroethyl)ether	40.000	35.467	11.3	106	0.00
12	1,3-Dichlorobenzene	40.000	35.600	11.0	107	0.00
13 C	1,4-Dichlorobenzene	40.000	35.517	11.2	107	0.00
14	1,2-Dichlorobenzene	40.000	35.351	11.6	107	0.00
15	Benzyl Alcohol	40.000	34.755	13.1	103	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	35.335	11.7	107	0.00
17	2-Methylphenol	40.000	34.981	12.5	103	0.00
18	Hexachloroethane	40.000	36.279	9.3	109	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	34.519	13.7	103	0.00
20	3+4-Methylphenols	40.000	34.737	13.2	103	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22	Acetophenone	40.000	34.607	13.5	102	0.00
23 S	Nitrobenzene-d5	80.000	71.671	10.4	104	0.00
24	Nitrobenzene	40.000	35.799	10.5	104	0.00
25	Isophorone	40.000	35.597	11.0	103	0.00
26 C	2-Nitrophenol	40.000	36.705	8.2	104	0.00
27	2,4-Dimethylphenol	40.000	35.323	11.7	101	0.00
28	bis(2-Chloroethoxy)methane	40.000	35.904	10.2	105	0.00
29 C	2,4-Dichlorophenol	40.000	35.604	11.0	102	0.00
30	1,2,4-Trichlorobenzene	40.000	35.981	10.0	104	0.00
31	Naphthalene	40.000	35.742	10.6	104	0.00
32	Benzoic acid	40.000	34.504	13.7	97	0.00
33	4-Chloroaniline	40.000	34.933	12.7	101	0.00
34 C	Hexachlorobutadiene	40.000	36.184	9.5	107	0.00
35	Caprolactam	40.000	35.000	12.5	100	0.00
36 C	4-Chloro-3-methylphenol	40.000	35.095	12.3	102	0.00
37	2-Methylnaphthalene	40.000	35.205	12.0	102	0.00
38	1-Methylnaphthalene	40.000	35.085	12.3	102	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	35.208	12.0	101	0.00
41 P	Hexachlorocyclopentadiene	40.000	34.348	14.1	96	-0.01
42 S	2,4,6-Tribromophenol	80.000	66.071	17.4	97	0.00
43 C	2,4,6-Trichlorophenol	40.000	35.528	11.2	101	0.00
44	2,4,5-Trichlorophenol	40.000	35.409	11.5	101	0.00
45 S	2-Fluorobiphenyl	80.000	69.854	12.7	101	0.00
46	1,1'-Biphenyl	40.000	35.149	12.1	102	0.00
47	2-Chloronaphthalene	40.000	35.369	11.6	102	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011022\
 Data File : BP008744.D
 Acq On : 10 Jan 2022 10:51
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 10 16:32:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 07 07:45:52 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	37.175	7.1	103	0.00
49	Acenaphthylene	40.000	35.410	11.5	102	0.00
50	Dimethylphthalate	40.000	34.485	13.8	101	0.00
51	2,6-Dinitrotoluene	40.000	35.531	11.2	102	0.00
52 C	Acenaphthene	40.000	35.526	11.2	101	0.00
53	3-Nitroaniline	40.000	35.756	10.6	100	0.00
54 P	2,4-Dinitrophenol	40.000	33.910	15.2	93	0.00
55	Dibenzofuran	40.000	35.135	12.2	101	0.00
56 P	4-Nitrophenol	40.000	34.814	13.0	96	0.00
57	2,4-Dinitrotoluene	40.000	36.242	9.4	103	0.00
58	Fluorene	40.000	35.408	11.5	112	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	34.938	12.7	100	0.00
60	Diethylphthalate	40.000	35.519	11.2	103	0.00
61	4-Chlorophenyl-phenylether	40.000	35.374	11.6	111	0.00
62	4-Nitroaniline	40.000	35.894	10.3	109	0.00
63	Azobenzene	40.000	37.465	6.3	121	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	107	0.00
65	4,6-Dinitro-2-methylphenol	40.000	36.986	7.5	103	0.00
66 c	n-Nitrosodiphenylamine	40.000	36.287	9.3	111	0.00
67	4-Bromophenyl-phenylether	40.000	35.337	11.7	105	0.00
68	Hexachlorobenzene	40.000	34.661	13.3	102	0.00
69	Atrazine	40.000	36.236	9.4	109	0.00
70 C	Pentachlorophenol	40.000	33.327	16.7	94	0.00
71	Phenanthrene	40.000	35.708	10.7	109	0.00
72	Anthracene	40.000	35.983	10.0	109	0.00
73	Carbazole	40.000	35.564	11.1	107	0.00
74	Di-n-butylphthalate	40.000	37.469	6.3	112	0.00
75 C	Fluoranthene	40.000	36.666	8.3	107	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	104	0.00
77	Benzydine	40.000	32.038	19.9	92	0.00
78	Pyrene	40.000	36.219	9.5	106	0.00
79 S	Terphenyl-d14	80.000	74.961	6.3	103	0.00
80	Butylbenzylphthalate	40.000	38.094	4.8	110	0.00
81	Benzo(a)anthracene	40.000	35.990	10.0	104	0.00
82	3,3'-Dichlorobenzidine	40.000	35.073	12.3	100	0.00
83	Chrysene	40.000	35.937	10.2	105	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.602	3.5	113	0.00
85 c	Di-n-octyl phthalate	40.000	38.402	4.0	111	0.00
86 I	Perylene-d12	20.000	20.000	0.0	95	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	34.830	12.9	92	-0.01
88	Benzo(b)fluoranthene	40.000	36.369	9.1	99	0.00
89	Benzo(k)fluoranthene	40.000	37.880	5.3	100	0.00
90 C	Benzo(a)pyrene	40.000	36.850	7.9	98	0.00
91	Dibenzo(a,h)anthracene	40.000	34.736	13.2	92	-0.01
92	Benzo(g,h,i)perylene	40.000	34.184	14.5	91	-0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011022\
Data File : BP008744.D
Acq On : 10 Jan 2022 10:51
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jan 10 16:32:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 07 07:45:52 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