

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030521\
 Data File : BM028992.D
 Acq On : 05 Mar 2021 12:35
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 05 13:17:27 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 05 13:17:12 2021
 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	67	0.00
2	1,4-Dioxane	40.000	44.053	-10.1	73	0.00
3	Pyridine	40.000	42.771	-6.9	71	0.00
4	n-Nitrosodimethylamine	40.000	55.785	-39.5#	91	0.00
5 S	2-Fluorophenol	80.000	82.367	-3.0	68	0.00
6	Aniline	40.000	40.160	-0.4	67	0.00
7 S	Phenol-d6	80.000	79.462	0.7	67	0.00
8	2-Chlorophenol	40.000	39.317	1.7	66	0.00
9	Benzaldehyde	40.000	34.237	14.4	65	0.00
10 C	Phenol	40.000	39.355	1.6	66	0.00
11	bis(2-Chloroethyl)ether	40.000	38.436	3.9	65	0.00
12	1,3-Dichlorobenzene	40.000	40.256	-0.6	68	0.00
13 C	1,4-Dichlorobenzene	40.000	39.697	0.8	66	0.00
14	1,2-Dichlorobenzene	40.000	39.955	0.1	68	0.00
15	Benzyl Alcohol	40.000	40.629	-1.6	67	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	43.978	-9.9	75	0.00
17	2-Methylphenol	40.000	39.304	1.7	65	0.00
18	Hexachloroethane	40.000	41.094	-2.7	69	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	40.295	-0.7	67	0.00
20	3+4-Methylphenols	40.000	40.100	-0.3	67	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	65	0.00
22	Acetophenone	40.000	41.474	-3.7	67	0.00
23 S	Nitrobenzene-d5	80.000	86.104	-7.6	70	0.00
24	Nitrobenzene	40.000	42.699	-6.7	69	0.00
25	Isophorone	40.000	40.705	-1.8	66	0.00
26 C	2-Nitrophenol	40.000	41.159	-2.9	65	0.00
27	2,4-Dimethylphenol	40.000	40.213	-0.5	66	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.939	2.7	63	0.00
29 C	2,4-Dichlorophenol	40.000	41.058	-2.6	66	0.00
30	1,2,4-Trichlorobenzene	40.000	41.097	-2.7	66	0.00
31	Naphthalene	40.000	39.815	0.5	66	0.00
32	Benzoic acid	40.000	42.908	-7.3	65	0.00
33	4-Chloroaniline	40.000	40.227	-0.6	65	0.00
34 C	Hexachlorobutadiene	40.000	41.661	-4.2	68	0.00
35	Caprolactam	40.000	39.512	1.2	63	0.00
36 C	4-Chloro-3-methylphenol	40.000	41.809	-4.5	68	0.00
37	2-Methylnaphthalene	40.000	39.595	1.0	65	0.00
38	1-Methylnaphthalene	40.000	39.112	2.2	64	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	65	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.530	1.2	65	0.00
41 P	Hexachlorocyclopentadiene	40.000	40.301	-0.8	64	0.00
42 S	2,4,6-Tribromophenol	80.000	84.747	-5.9	67	0.00
43 C	2,4,6-Trichlorophenol	40.000	41.457	-3.6	66	0.00
44	2,4,5-Trichlorophenol	40.000	41.745	-4.4	68	0.00
45 S	2-Fluorobiphenyl	80.000	80.181	-0.2	67	0.00
46	1,1'-Biphenyl	40.000	39.445	1.4	65	0.00
47	2-Chloronaphthalene	40.000	40.049	-0.1	66	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030521\
 Data File : BM028992.D
 Acq On : 05 Mar 2021 12:35
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 05 13:17:27 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 05 13:17:12 2021
 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	46.204	-15.5	73	0.00
49	Acenaphthylene	40.000	40.490	-1.2	66	0.00
50	Dimethylphthalate	40.000	41.413	-3.5	67	0.00
51	2,6-Dinitrotoluene	40.000	41.824	-4.6	66	0.00
52 C	Acenaphthene	40.000	39.888	0.3	66	0.00
53	3-Nitroaniline	40.000	43.239	-8.1	68	0.00
54 P	2,4-Dinitrophenol	40.000	43.162	-7.9	67	0.00
55	Dibenzofuran	40.000	41.296	-3.2	68	0.00
56 P	4-Nitrophenol	40.000	42.229	-5.6	68	0.00
57	2,4-Dinitrotoluene	40.000	44.571	-11.4	70	0.00
58	Fluorene	40.000	41.545	-3.9	68	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	41.900	-4.7	67	0.00
60	Diethylphthalate	40.000	42.190	-5.5	68	0.00
61	4-Chlorophenyl-phenylether	40.000	41.613	-4.0	69	0.00
62	4-Nitroaniline	40.000	44.404	-11.0	68	0.00
63	Azobenzene	40.000	44.143	-10.4	71	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	69	0.00
65	4,6-Dinitro-2-methylphenol	40.000	40.258	-0.6	68	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.165	4.6	66	0.00
67	4-Bromophenyl-phenylether	40.000	38.621	3.4	67	0.00
68	Hexachlorobenzene	40.000	38.624	3.4	68	0.00
69	Atrazine	40.000	40.456	-1.1	69	0.00
70 C	Pentachlorophenol	40.000	44.015	-10.0	71	0.00
71	Phenanthrene	40.000	39.461	1.3	69	0.00
72	Anthracene	40.000	39.961	0.1	69	0.00
73	Carbazole	40.000	40.294	-0.7	69	0.00
74	Di-n-butylphthalate	40.000	41.005	-2.5	68	0.00
75 C	Fluoranthene	40.000	41.792	-4.5	71	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	73	0.00
77	Benzydine	40.000	36.416	9.0	58	0.00
78	Pyrene	40.000	38.390	4.0	72	0.00
79 S	Terphenyl-d14	80.000	77.915	2.6	73	0.00
80	Butylbenzylphthalate	40.000	39.567	1.1	71	0.00
81	Benzo(a)anthracene	40.000	39.979	0.1	74	0.00
82	3,3'-Dichlorobenzidine	40.000	40.012	-0.0	72	0.00
83	Chrysene	40.000	40.093	-0.2	74	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.046	-0.1	71	0.00
85 c	Di-n-octyl phthalate	40.000	40.392	-1.0	71	0.00
86 I	Perylene-d12	20.000	20.000	0.0	74	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	38.447	3.9	70	0.00
88	Benzo(b)fluoranthene	40.000	40.142	-0.4	73	0.00
89	Benzo(k)fluoranthene	40.000	38.761	3.1	73	0.00
90 C	Benzo(a)pyrene	40.000	39.225	1.9	73	0.00
91	Dibenzo(a,h)anthracene	40.000	38.370	4.1	70	0.00
92	Benzo(g,h,i)perylene	40.000	38.588	3.5	71	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030521\
Data File : BM028992.D
Acq On : 05 Mar 2021 12:35
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 05 13:17:27 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 05 13:17:12 2021
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