

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061521\
 Data File : BP005922.D
 Acq On : 15 Jun 2021 10:12
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 15 11:31:56 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 14 13:15:03 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	76	0.00
2	1,4-Dioxane	0.563	0.586	-4.1	83	0.00
3	Pyridine	1.323	1.559	-17.8	89	0.00
4	n-Nitrosodimethylamine	0.615	0.614	0.2	80	0.00
5 S	2-Fluorophenol	1.246	1.406	-12.8	89	0.00
6	Aniline	1.804	2.041	-13.1	90	0.00
7 S	Phenol-d6	1.616	1.822	-12.7	89	0.00
8	2-Chlorophenol	1.429	1.610	-12.7	91	0.00
9	Benzaldehyde	0.689	0.797	-15.7	84	0.00
10 C	Phenol	1.614	1.826	-13.1	91	0.00
11	bis(2-Chloroethyl)ether	1.223	1.392	-13.8	92	0.00
12	1,3-Dichlorobenzene	1.598	1.790	-12.0	90	0.00
13 C	1,4-Dichlorobenzene	1.630	1.796	-10.2	89	0.00
14	1,2-Dichlorobenzene	1.558	1.736	-11.4	90	0.00
15	Benzyl Alcohol	1.217	1.348	-10.8	87	0.00
16	2,2'-oxybis(1-Chloropropane	1.130	1.169	-3.5	82	0.00
17	2-Methylphenol	1.100	1.265	-15.0	91	0.00
18	Hexachloroethane	0.589	0.638	-8.3	86	0.00
19 P	n-Nitroso-di-n-propylamine	0.994	1.099	-10.6	88	0.00
20	3+4-Methylphenols	1.485	1.725	-16.2	92	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	77	0.00
22	Acetophenone	0.526	0.568	-8.0	87	0.00
23 S	Nitrobenzene-d5	0.408	0.422	-3.4	84	0.00
24	Nitrobenzene	0.424	0.437	-3.1	86	0.00
25	Isophorone	0.753	0.793	-5.3	87	0.00
26 C	2-Nitrophenol	0.202	0.223	-10.4	91	0.00
27	2,4-Dimethylphenol	0.301	0.318	-5.6	88	0.00
28	bis(2-Chloroethoxy)methane	0.393	0.428	-8.9	92	0.00
29 C	2,4-Dichlorophenol	0.362	0.388	-7.2	90	0.00
30	1,2,4-Trichlorobenzene	0.407	0.423	-3.9	88	0.00
31	Naphthalene	1.157	1.228	-6.1	90	0.00
32	Benzoic acid	0.193	0.233	-20.7	85	0.00
33	4-Chloroaniline	0.467	0.507	-8.6	91	0.00
34 C	Hexachlorobutadiene	0.275	0.287	-4.4	89	0.00
35	Caprolactam	0.104	0.114	-9.6	85	0.00
36 C	4-Chloro-3-methylphenol	0.367	0.398	-8.4	90	0.00
37	2-Methylnaphthalene	0.824	0.867	-5.2	89	0.00
38	1-Methylnaphthalene	0.776	0.827	-6.6	89	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	77	0.00
40	1,2,4,5-Tetrachlorobenzene	0.704	0.720	-2.3	84	0.00
41 P	Hexachlorocyclopentadiene	0.297	0.302	-1.7	79	0.00
42 S	2,4,6-Tribromophenol	0.343	0.372	-8.5	88	0.00
43 C	2,4,6-Trichlorophenol	0.483	0.488	-1.0	83	0.00
44	2,4,5-Trichlorophenol	0.520	0.535	-2.9	84	0.00
45 S	2-Fluorobiphenyl	1.523	1.573	-3.3	86	0.00
46	1,1'-Biphenyl	1.613	1.659	-2.9	84	0.00
47	2-Chloronaphthalene	1.317	1.361	-3.3	88	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061521\
 Data File : BP005922.D
 Acq On : 15 Jun 2021 10:12
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 15 11:31:56 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 14 13:15:03 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48	2-Nitroaniline	0.346	0.368	-6.4	87	0.00
49	Acenaphthylene	2.020	2.109	-4.4	87	0.00
50	Dimethylphthalate	1.644	1.717	-4.4	89	0.00
51	2,6-Dinitrotoluene	0.374	0.391	-4.5	87	0.00
52 C	Acenaphthene	1.227	1.276	-4.0	88	0.00
53	3-Nitroaniline	0.362	0.390	-7.7	89	0.00
54 P	2,4-Dinitrophenol	0.191	0.242	-26.7#	96	0.00
55	Dibenzofuran	2.018	2.041	-1.1	86	0.00
56 P	4-Nitrophenol	0.294	0.308	-4.8	84	0.00
57	2,4-Dinitrotoluene	0.500	0.538	-7.6	88	0.00
58	Fluorene	1.572	1.621	-3.1	87	0.00
59	2,3,4,6-Tetrachlorophenol	0.458	0.459	-0.2	83	0.00
60	Diethylphthalate	1.650	1.726	-4.6	88	0.00
61	4-Chlorophenyl-phenylether	0.815	0.833	-2.2	87	0.00
62	4-Nitroaniline	0.373	0.409	-9.7	91	0.00
63	Azobenzene	1.493	1.501	-0.5	84	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	75	0.00
65	4,6-Dinitro-2-methylphenol	0.147	0.140	4.8	76	0.00
66 c	n-Nitrosodiphenylamine	0.663	0.691	-4.2	86	0.00
67	4-Bromophenyl-phenylether	0.257	0.270	-5.1	87	0.00
68	Hexachlorobenzene	0.308	0.318	-3.2	86	0.00
69	Atrazine	0.206	0.186	9.7	63	0.00
70 C	Pentachlorophenol	0.171	0.179	-4.7	79	0.00
71	Phenanthrene	1.201	1.243	-3.5	85	0.00
72	Anthracene	1.182	1.245	-5.3	86	0.00
73	Carbazole	1.135	1.208	-6.4	87	0.00
74	Di-n-butylphthalate	1.315	1.450	-10.3	90	0.00
75 C	Fluoranthene	1.459	1.560	-6.9	89	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	84	0.00
77	Benzydine	0.416	0.385	7.5	68	0.00
78	Pyrene	1.396	1.386	0.7	92	0.00
79 S	Terphenyl-d14	1.068	1.101	-3.1	94	0.00
80	Butylbenzylphthalate	0.574	0.615	-7.1	97	0.00
81	Benzo(a)anthracene	1.444	1.477	-2.3	95	0.00
82	3,3'-Dichlorobenzidine	0.499	0.527	-5.6	99	0.00
83	Chrysene	1.398	1.452	-3.9	97	0.00
84	Bis(2-ethylhexyl)phthalate	0.842	0.918	-9.0	100	0.00
85 c	Di-n-octyl phthalate	1.507	1.630	-8.2	99	0.00
86 I	Perylene-d12	1.000	1.000	0.0	81	0.00
87	Indeno(1,2,3-cd)pyrene	1.710	1.844	-7.8	96	0.00
88	Benzo(b)fluoranthene	1.370	1.515	-10.6	98	0.00
89	Benzo(k)fluoranthene	1.328	1.455	-9.6	97	0.00
90 C	Benzo(a)pyrene	1.288	1.414	-9.8	97	0.00
91	Dibenzo(a,h)anthracene	1.472	1.582	-7.5	96	0.00
92	Benzo(g,h,i)perylene	1.454	1.544	-6.2	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061521\
Data File : BP005922.D
Acq On : 15 Jun 2021 10:12
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jun 15 11:31:56 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 14 13:15:03 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
----------	-------	------	------	-------	----------

(#) = Out of Range

SPCC's out = 0 CCC's out = 0