

Data Path : Z:\HPCHEM1\BNA F\Data\BF050218\
 Data File : BF105047.D
 Acq On : 2 May 2018 23:24
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: May 03 04:50:35 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 02 17:25:18 2018
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	76	0.00
2	1,4-Dioxane	0.609	0.477	21.7	59	-0.06
3	Pyridine	1.714	1.328	22.5	59	-0.05
4	n-Nitrosodimethylamine	0.599	0.454	24.2	57	-0.06
5 S	2-Fluorophenol	1.308	1.168	10.7	69	0.00
6	Aniline	2.233	1.868	16.3	63	0.00
7 S	Phenol-d6	1.607	1.422	11.5	68	0.00
8	2-Chlorophenol	1.381	1.288	6.7	71	0.00
9	Benzaldehyde	0.802	0.702	12.5	64	0.00
10 C	Phenol	1.705	1.575	7.6	72	0.00
11	bis(2-Chloroethyl)ether	1.361	1.209	11.2	67	0.00
12	1,3-Dichlorobenzene	1.564	1.450	7.3	71	0.00
13 C	1,4-Dichlorobenzene	1.559	1.482	4.9	72	0.00
14	1,2-Dichlorobenzene	1.445	1.384	4.2	72	0.00
15	Benzyl Alcohol	1.221	1.008	17.4	63	0.00
16	2,2'-oxybis(1-Chloropropane	1.827	1.354	25.9#	56	0.00
17	2-Methylphenol	1.200	1.092	9.0	69	0.00
18	Hexachloroethane	0.556	0.448	19.4	62	0.00
19 P	n-Nitroso-di-n-propylamine	1.050	0.826	21.3	63	0.00
20	3+4-Methylphenols	1.488	1.344	9.7	69	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	77	0.00
22	Acetophenone	0.492	0.434	11.8	68	0.00
23 S	Nitrobenzene-d5	0.306	0.304	0.7	74	0.00
24	Nitrobenzene	0.338	0.324	4.1	70	0.00
25	Isophorone	0.648	0.539	16.8	64	0.00
26 C	2-Nitrophenol	0.138	0.165	-19.6	87	0.00
27	2,4-Dimethylphenol	0.286	0.239	16.4	63	0.00
28	bis(2-Chloroethoxy)methane	0.396	0.344	13.1	67	0.00
29 C	2,4-Dichlorophenol	0.273	0.255	6.6	71	0.00
30	1,2,4-Trichlorobenzene	0.299	0.282	5.7	73	0.00
31	Naphthalene	0.996	0.903	9.3	70	0.00
32	Benzoic acid	0.228	0.168	26.3#	53	-0.02
33	4-Chloroaniline	0.427	0.390	8.7	70	0.00
34 C	Hexachlorobutadiene	0.164	0.159	3.0	75	0.00
35	Caprolactam	0.095	0.081	14.7	64	-0.01
36 C	4-Chloro-3-methylphenol	0.292	0.266	8.9	69	0.00
37	2-Methylnaphthalene	0.644	0.586	9.0	70	0.00
38 I	Acenaphthene-d10	1.000	1.000	0.0	71	0.00
39	1,2,4,5-Tetrachlorobenzene	0.573	0.597	-4.2	76	0.00
40 P	Hexachlorocyclopentadiene	0.321	0.013#	96.0#	3#	0.00
41 S	2,4,6-Tribromophenol	0.207	0.182	12.1	62	0.00
42 C	2,4,6-Trichlorophenol	0.397	0.384	3.3	67	0.00
43	2,4,5-Trichlorophenol	0.445	0.435	2.2	70	0.00
44 S	2-Fluorobiphenyl	1.266	1.257	0.7	72	0.00

Data Path : Z:\HPCHEM1\BNA F\Data\BF050218\
 Data File : BF105047.D
 Acq On : 2 May 2018 23:24
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: May 03 04:50:35 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 02 17:25:18 2018
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	1,1'-Biphenyl	1.680	1.637	2.6	70	0.00
46	2-Chloronaphthalene	1.327	1.269	4.4	69	0.00
47	2-Nitroaniline	0.328	0.365	-11.3	74	0.00
48	Acenaphthylene	2.082	1.858	10.8	64	0.00
49	Dimethylphthalate	1.577	1.524	3.4	70	0.00
50	2,6-Dinitrotoluene	0.292	0.317	-8.6	71	0.00
51 C	Acenaphthene	1.270	1.105	13.0	63	0.00
52	3-Nitroaniline	0.342	0.376	-9.9	73	0.00
53 P	2,4-Dinitrophenol	0.089	0.022#	75.3#	18#	0.00
54	Dibenzofuran	1.770	1.546	12.7	62	0.00
55 P	4-Nitrophenol	0.268	0.190	29.1#	46#	0.00
56	2,4-Dinitrotoluene	0.383	0.371	3.1	66	0.00
57	Fluorene	1.289	1.257	2.5	69	0.00
58	2,3,4,6-Tetrachlorophenol	0.332	0.280	15.7	60	0.00
59	Diethylphthalate	1.571	1.435	8.7	66	0.00
60	4-Chlorophenyl-phenylether	0.574	0.595	-3.7	73	0.00
61	4-Nitroaniline	0.334	0.356	-6.6	69	0.00
62	Azobenzene	1.501	1.206	19.7	58	0.00
63 I	Phenanthrene-d10	1.000	1.000	0.0	64	0.00
64	4,6-Dinitro-2-methylphenol	0.087	0.038	56.3#	26#	0.00
65 c	n-Nitrosodiphenylamine	0.693	0.688	0.7	64	0.00
66	4-Bromophenyl-phenylether	0.213	0.215	-0.9	66	0.00
67	Hexachlorobenzene	0.239	0.232	2.9	63	0.00
68	Atrazine	0.209	0.188	10.0	58	0.00
69 C	Pentachlorophenol	0.148	0.095	35.8#	39#	0.00
70	Phenanthrene	1.114	0.993	10.9	58	0.00
71	Anthracene	1.132	1.014	10.4	59	0.00
72	Carbazole	1.106	0.959	13.3	57	0.00
73	Di-n-butylphthalate	1.351	1.268	6.1	61	0.00
74 C	Fluoranthene	1.164	0.971	16.6	54	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	63	-0.02
76	Benzidine	0.692	0.675	2.5	62	0.00
77	Pyrene	1.482	1.269	14.4	55	0.00
78 S	Terphenyl-d14	0.877	0.784	10.6	59	0.00
79	Butylbenzylphthalate	0.698	0.627	10.2	57	-0.01
80	Benzo(a)anthracene	1.195	1.160	2.9	63	-0.02
81	3,3'-Dichlorobenzidine	0.445	0.429	3.6	60	-0.02
82	Chrysene	1.227	1.111	9.5	58	-0.02
83	Bis(2-ethylhexyl)phthalate	0.898	0.884	1.6	63	-0.02
84 c	Di-n-octyl phthalate	1.501	1.406	6.3	58	-0.03
85	Indeno(1,2,3-cd)pyrene	1.043	0.826	20.8	52	-0.05
86 I	Perylene-d12	1.000	1.000	0.0	56	-0.04
87	Benzo(b)fluoranthene	1.263	1.160	8.2	51	-0.04

Data Path : Z:\HPCHEM1\BNA F\Data\BF050218\
Data File : BF105047.D
Acq On : 2 May 2018 23:24
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 03 04:50:35 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 02 17:25:18 2018
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)
88	Benzo(k)fluoranthene	1.193	1.230	-3.1	59	-0.04
89 C	Benzo(a)pyrene	1.130	1.048	7.3	52	-0.04
90	Dibenzo(a,h)anthracene	0.928	0.840	9.5	53	-0.05
91	Benzo(a,h,i)perylene	0.899	0.788	12.3	53	-0.05

(#) = Out of Range

SPCC's out = 2 CCC's out = 1