

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030420\
 Data File : BF119187.D
 Acq On : 4 Mar 2020 10:53
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 04 13:00:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 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	84	0.00
2	1,4-Dioxane	0.533	0.490	8.1	82	-0.02
3	Pyridine	1.455	1.376	5.4	85	-0.02
4	n-Nitrosodimethylamine	0.441	0.470	-6.6	95	-0.03
5 S	2-Fluorophenol	1.197	1.240	-3.6	91	0.00
6	Aniline	1.796	1.840	-2.4	89	0.00
7 S	Phenol-d6	1.455	1.525	-4.8	92	0.00
8	2-Chlorophenol	1.292	1.378	-6.7	93	0.00
9	Benzaldehyde	0.972	0.875	10.0	79	0.00
10 C	Phenol	1.574	1.643	-4.4	92	0.00
11	bis(2-Chloroethyl)ether	1.204	1.278	-6.1	94	-0.01
12	1,3-Dichlorobenzene	1.548	1.608	-3.9	92	0.00
13 C	1,4-Dichlorobenzene	1.549	1.627	-5.0	92	0.00
14	1,2-Dichlorobenzene	1.413	1.508	-6.7	93	0.00
15	Benzyl Alcohol	1.052	1.190	-13.1	98	0.00
16	2,2'-oxybis(1-Chloropropane	1.043	1.104	-5.8	94	0.00
17	2-Methylphenol	1.047	1.106	-5.6	93	0.00
18	Hexachloroethane	0.554	0.591	-6.7	95	0.00
19 P	n-Nitroso-di-n-propylamine	0.848	0.954	-12.5	101	-0.01
20	3+4-Methylphenols	1.283	1.415	-10.3	100	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
22	Acetophenone	0.463	0.501	-8.2	100	0.00
23 S	Nitrobenzene-d5	0.379	0.402	-6.1	99	0.00
24	Nitrobenzene	0.375	0.394	-5.1	98	0.00
25	Isophorone	0.658	0.694	-5.5	100	0.00
26 C	2-Nitrophenol	0.198	0.209	-5.6	98	0.00
27	2,4-Dimethylphenol	0.269	0.277	-3.0	96	0.00
28	bis(2-Chloroethoxy)methane	0.428	0.441	-3.0	96	-0.01
29 C	2,4-Dichlorophenol	0.317	0.332	-4.7	96	0.00
30	1,2,4-Trichlorobenzene	0.376	0.386	-2.7	95	0.00
31	Naphthalene	1.037	1.073	-3.5	95	0.00
32	Benzoic acid	0.184	0.158	14.1	81	0.00
33	4-Chloroaniline	0.446	0.468	-4.9	96	0.00
34 C	Hexachlorobutadiene	0.234	0.244	-4.3	97	-0.01
35	Caprolactam	0.098	0.103	-5.1	97	-0.01
36 C	4-Chloro-3-methylphenol	0.321	0.347	-8.1	100	0.00
37	2-Methylnaphthalene	0.698	0.729	-4.4	97	0.00
38	1-Methylnaphthalene	0.656	0.691	-5.3	97	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
40	1,2,4,5-Tetrachlorobenzene	0.674	0.684	-1.5	97	0.00
41 P	Hexachlorocyclopentadiene	0.198	0.160	19.2	78	0.00
42 S	2,4,6-Tribromophenol	0.262	0.275	-5.0	100	0.00
43 C	2,4,6-Trichlorophenol	0.426	0.433	-1.6	97	0.00
44	2,4,5-Trichlorophenol	0.447	0.433	3.1	92	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030420\
 Data File : BF119187.D
 Acq On : 4 Mar 2020 10:53
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 04 13:00:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 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 S	2-Fluorobiphenyl	1.358	1.360	-0.1	98	0.00
46	1,1'-Biphenyl	1.616	1.657	-2.5	97	0.00
47	2-Chloronaphthalene	1.303	1.342	-3.0	98	0.00
48	2-Nitroaniline	0.363	0.402	-10.7	105	0.00
49	Acenaphthylene	1.881	1.924	-2.3	97	0.00
50	Dimethylphthalate	1.529	1.597	-4.4	101	-0.01
51	2,6-Dinitrotoluene	0.340	0.350	-2.9	99	0.00
52 C	Acenaphthene	1.134	1.169	-3.1	97	0.00
53	3-Nitroaniline	0.377	0.392	-4.0	98	0.00
54 P	2,4-Dinitrophenol	0.140	0.142	-1.4	95	0.00
55	Dibenzofuran	1.693	1.707	-0.8	94	0.00
56 P	4-Nitrophenol	0.226	0.195	13.7	80	0.00
57	2,4-Dinitrotoluene	0.440	0.445	-1.1	94	0.00
58	Fluorene	1.316	1.411	-7.2	101	0.00
59	2,3,4,6-Tetrachlorophenol	0.377	0.359	4.8	92	0.00
60	Diethylphthalate	1.441	1.547	-7.4	102	0.00
61	4-Chlorophenyl-phenylether	0.696	0.740	-6.3	100	0.00
62	4-Nitroaniline	0.385	0.405	-5.2	99	0.00
63	Azobenzene	1.253	1.354	-8.1	103	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
65	4,6-Dinitro-2-methylphenol	0.121	0.127	-5.0	100	0.00
66 c	n-Nitrosodiphenylamine	0.655	0.688	-5.0	99	0.00
67	4-Bromophenyl-phenylether	0.261	0.274	-5.0	101	0.00
68	Hexachlorobenzene	0.301	0.316	-5.0	100	0.00
69	Atrazine	0.229	0.213	7.0	89	0.00
70 C	Pentachlorophenol	0.121	0.111	8.3	88	0.00
71	Phenanthrene	1.091	1.143	-4.8	99	0.00
72	Anthracene	1.090	1.148	-5.3	98	0.00
73	Carbazole	0.971	1.028	-5.9	100	0.00
74	Di-n-butylphthalate	1.176	1.306	-11.1	105	0.00
75 C	Fluoranthene	1.158	1.224	-5.7	99	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	80	0.00
77	Benzidine	0.813	0.757	6.9	77	0.00
78	Pyrene	1.606	1.762	-9.7	96	0.00
79 S	Terphenyl-d14	1.162	1.332	-14.6	101	0.00
80	Butylbenzylphthalate	0.676	0.802	-18.6	102	0.00
81	Benzo(a)anthracene	1.363	1.496	-9.8	93	0.00
82	3,3'-Dichlorobenzidine	0.529	0.547	-3.4	84	0.00
83	Chrysene	1.358	1.394	-2.7	88	0.00
84	Bis(2-ethylhexyl)phthalate	0.854	1.037	-21.4	104	0.00
85 c	Di-n-octyl phthalate	1.361	1.532	-12.6	91	0.00
86	Indeno(1,2,3-cd)pyrene	1.315	1.112	15.4	56	0.01
87 I	Perylene-d12	1.000	1.000	0.0	45#	0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030420\
Data File : BF119187.D
Acq On : 4 Mar 2020 10:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Mar 04 13:00:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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(b)fluoranthene	1.318	1.499	-13.7	78	0.00
89	Benzo(k)fluoranthene	1.222	1.349	-10.4	69	0.00
90 C	Benzo(a)pyrene	1.190	1.260	-5.9	52	0.00
91	Dibenzo(a,h)anthracene	1.047	1.118	-6.8	55	0.00
92	Benzo(q,h,i)perylene	1.034	1.087	-5.1	55	0.02

(#) = Out of Range

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