

Data Path : Z:\HPCHEM1\BNA F\Data\BF031518\
 Data File : BF103674.D
 Acq On : 15 Mar 2018 17:56
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF031518

Quant Time: Mar 15 18:32:15 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 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	97	0.00
2	1,4-Dioxane	0.647	0.636	1.7	94	-0.02
3	Pyridine	1.781	1.776	0.3	97	-0.02
4	n-Nitrosodimethylamine	0.657	0.637	3.0	94	-0.02
5 S	2-Fluorophenol	1.315	1.301	1.1	95	0.00
6	Aniline	2.297	2.200	4.2	92	0.00
7 S	Phenol-d6	1.609	1.573	2.2	95	0.00
8	2-Chlorophenol	1.377	1.374	0.2	96	0.00
9	Benzaldehyde	1.057	1.034	2.2	95	0.00
10 C	Phenol	1.709	1.683	1.5	96	0.00
11	bis(2-Chloroethyl)ether	1.411	1.358	3.8	94	0.00
12	1,3-Dichlorobenzene	1.569	1.510	3.8	94	0.00
13 C	1,4-Dichlorobenzene	1.576	1.531	2.9	94	0.00
14	1,2-Dichlorobenzene	1.463	1.452	0.8	97	0.00
15	Benzyl Alcohol	1.246	1.232	1.1	95	0.00
16	2,2'-oxybis(1-Chloropropane	2.007	1.933	3.7	94	0.00
17	2-Methylphenol	1.227	1.190	3.0	94	0.00
18	Hexachloroethane	0.555	0.560	-0.9	96	0.00
19 P	n-Nitroso-di-n-propylamine	1.031	0.998	3.2	95	0.00
20	3+4-Methylphenols	1.557	1.558	-0.1	96	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22	Acetophenone	0.514	0.493	4.1	95	0.00
23 S	Nitrobenzene-d5	0.287	0.320	-11.5	104	0.00
24	Nitrobenzene	0.337	0.355	-5.3	98	0.00
25	Isophorone	0.664	0.634	4.5	95	0.00
26 C	2-Nitrophenol	0.110	0.136	-23.6#	118	0.00
27	2,4-Dimethylphenol	0.284	0.283	0.4	97	0.00
28	bis(2-Chloroethoxy)methane	0.402	0.388	3.5	95	0.00
29 C	2,4-Dichlorophenol	0.259	0.260	-0.4	95	0.00
30	1,2,4-Trichlorobenzene	0.300	0.290	3.3	95	0.00
31	Naphthalene	1.010	0.976	3.4	96	0.00
32	Benzoic acid	0.171	0.211	-23.4	119	0.00
33	4-Chloroaniline	0.424	0.403	5.0	92	0.00
34 C	Hexachlorobutadiene	0.165	0.159	3.6	95	0.00
35	Caprolactam	0.089	0.089	0.0	97	0.00
36 C	4-Chloro-3-methylphenol	0.285	0.284	0.4	95	0.00
37	2-Methylnaphthalene	0.641	0.618	3.6	95	0.00
38 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
39	1,2,4,5-Tetrachlorobenzene	0.571	0.557	2.5	97	0.00
40 P	Hexachlorocyclopentadiene	0.294	0.317	-7.8	101	0.00
41 S	2,4,6-Tribromophenol	0.172	0.186	-8.1	99	0.00
42 C	2,4,6-Trichlorophenol	0.365	0.379	-3.8	97	0.00
43	2,4,5-Trichlorophenol	0.412	0.432	-4.9	99	0.00
44 S	2-Fluorobiphenyl	1.254	1.205	3.9	98	0.00

Data Path : Z:\HPCHEM1\BNA F\Data\BF031518\
 Data File : BF103674.D
 Acq On : 15 Mar 2018 17:56
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICSVBF031518

Quant Time: Mar 15 18:32:15 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 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	1.668	1.619	2.9	97	0.00
46	2-Chloronaphthalene	1.325	1.282	3.2	96	0.00
47	2-Nitroaniline	0.309	0.371	-20.1	111	0.00
48	Acenaphthylene	2.054	1.982	3.5	97	0.00
49	Dimethylphthalate	1.526	1.484	2.8	96	0.00
50	2,6-Dinitrotoluene	0.274	0.309	-12.8	105	0.00
51 C	Acenaphthene	1.220	1.199	1.7	97	0.00
52	3-Nitroaniline	0.331	0.363	-9.7	102	0.00
53 P	2,4-Dinitrophenol	0.057	0.073	-28.1#	141	0.00
54	Dibenzofuran	1.742	1.675	3.8	96	0.00
55 P	4-Nitrophenol	0.245	0.275	-12.2	104	0.00
56	2,4-Dinitrotoluene	0.336	0.391	-16.4	107	0.00
57	Fluorene	1.352	1.293	4.4	96	0.00
58	2,3,4,6-Tetrachlorophenol	0.292	0.313	-7.2	99	0.00
59	Diethylphthalate	1.514	1.473	2.7	96	0.00
60	4-Chlorophenyl-phenylether	0.618	0.585	5.3	95	0.00
61	4-Nitroaniline	0.319	0.354	-11.0	103	0.00
62	Azobenzene	1.540	1.480	3.9	96	0.00
63 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64	4,6-Dinitro-2-methylphenol	0.065	0.082	-26.2#	124	0.00
65 c	n-Nitrosodiphenylamine	0.681	0.663	2.6	96	0.00
66	4-Bromophenyl-phenylether	0.205	0.201	2.0	95	0.00
67	Hexachlorobenzene	0.234	0.227	3.0	95	0.00
68	Atrazine	0.211	0.203	3.8	94	0.00
69 C	Pentachlorophenol	0.135	0.145	-7.4	100	0.00
70	Phenanthrene	1.094	1.057	3.4	97	0.00
71	Anthracene	1.111	1.070	3.7	96	0.00
72	Carbazole	1.080	1.050	2.8	96	0.00
73	Di-n-butylphthalate	1.296	1.295	0.1	97	0.00
74 C	Fluoranthene	1.127	1.087	3.5	95	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76	Benzidine	0.853	0.815	4.5	93	0.00
77	Pyrene	1.469	1.381	6.0	96	0.00
78 S	Terphenyl-d14	0.862	0.791	8.2	95	0.00
79	Butylbenzylphthalate	0.651	0.662	-1.7	98	0.00
80	Benzo(a)anthracene	1.266	1.215	4.0	96	0.00
81	3,3'-Dichlorobenzidine	0.423	0.434	-2.6	96	0.00
82	Chrysene	1.207	1.179	2.3	99	0.00
83	Bis(2-ethylhexyl)phthalate	0.930	0.964	-3.7	99	0.00
84 c	Di-n-octyl phthalate	1.352	1.488	-10.1	100	0.01
85	Indeno(1,2,3-cd)pyrene	1.054	1.125	-6.7	105	0.02
86 I	Perylene-d12	1.000	1.000	0.0	103	0.01
87	Benzo(b)fluoranthene	1.264	1.231	2.6	102	0.01

Data Path : Z:\HPCHEM1\BNA F\Data\BF031518\
Data File : BF103674.D
Acq On : 15 Mar 2018 17:56
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF031518

Quant Time: Mar 15 18:32:15 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 15 18:31:53 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	1.215	1.138	6.3	96	0.02
89 C	Benzo(a)pyrene	1.146	1.118	2.4	100	0.02
90	Dibenzo(a,h)anthracene	0.992	1.006	-1.4	105	0.02
91	Benzo(a,h,i)perylene	0.965	0.968	-0.3	105	0.02

(#) = Out of Range

SPCC's out = 0 CCC's out = 1