

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
 Data File : BF083733.D
 Acq On : 13 Dec 2015 17:05
 Operator : UM/IZ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF121315

Quant Time: Dec 14 02:29:15 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 01:30:12 2015
 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	104	0.00
2	1,4-Dioxane	0.599	0.594	0.8	96	-0.02
3	Pyridine	1.508	1.549	-2.7	96	-0.01
4	n-Nitrosodimethylamine	0.571	0.571	0.0	96	-0.01
5 S	2-Fluorophenol	1.240	1.209	2.5	99	-0.01
6	Aniline	2.133	2.143	-0.5	97	0.00
7 S	Phenol-d6	1.570	1.530	2.5	97	0.00
8	2-Chlorophenol	1.457	1.527	-4.8	97	0.00
9	Benzaldehyde	0.843	0.867	-2.8	96	0.00
10 C	Phenol	1.807	1.802	0.3	97	0.00
11	bis(2-Chloroethyl)ether	1.309	1.246	4.8	100	-0.01
12	1,3-Dichlorobenzene	1.546	1.587	-2.7	97	0.00
13 C	1,4-Dichlorobenzene	1.562	1.525	2.4	99	-0.01
14	1,2-Dichlorobenzene	1.404	1.501	-6.9	99	-0.01
15	Benzyl Alcohol	1.153	1.243	-7.8	97	0.00
16	2,2'-oxybis(1-Chloropropane	1.699	1.726	-1.6	100	-0.01
17	2-Methylphenol	1.171	1.211	-3.4	98	0.00
18	Hexachloroethane	0.555	0.564	-1.6	99	-0.01
19 P	n-Nitroso-di-n-propylamine	0.961	0.892	7.2	99	-0.01
20	3+4-Methylphenols	1.383	1.326	4.1	97	-0.01
21 I	Naphthalene-d8	1.000	1.000	0.0	102	-0.01
22	Acetophenone	0.491	0.515	-4.9	100	0.00
23 S	Nitrobenzene-d5	0.357	0.365	-2.2	98	0.00
24	Nitrobenzene	0.372	0.389	-4.6	103	0.00
25	Isophorone	0.701	0.723	-3.1	98	0.00
26 C	2-Nitrophenol	0.179	0.179	0.0	100	-0.01
27	2,4-Dimethylphenol	0.351	0.362	-3.1	100	0.00
28	bis(2-Chloroethoxy)methane	0.400	0.388	3.0	97	-0.01
29 C	2,4-Dichlorophenol	0.288	0.283	1.7	96	-0.01
30	1,2,4-Trichlorobenzene	0.299	0.302	-1.0	100	-0.01
31	Naphthalene	1.043	1.025	1.7	97	-0.01
32	Benzoic acid	0.219	0.222	-1.4	98	0.00
33	4-Chloroaniline	0.426	0.464	-8.9	97	0.00
34 C	Hexachlorobutadiene	0.164	0.175	-6.7	101	-0.01
35	Caprolactam	0.093	0.091	2.2	95	0.00
36 C	4-Chloro-3-methylphenol	0.337	0.364	-8.0	97	0.00
37	2-Methylnaphthalene	0.655	0.661	-0.9	100	-0.01
38 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
39	1,2,4,5-Tetrachlorobenzene	0.590	0.576	2.4	99	-0.01
40 P	Hexachlorocyclopentadiene	0.299	0.331	-10.7	100	0.00
41 S	2,4,6-Tribromophenol	0.204	0.201	1.5	101	-0.01
42 C	2,4,6-Trichlorophenol	0.432	0.435	-0.7	99	0.00
43	2,4,5-Trichlorophenol	0.410	0.450	-9.8	99	0.00
44 S	2-Fluorobiphenyl	1.406	1.345	4.3	97	-0.01

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
 Data File : BF083733.D
 Acq On : 13 Dec 2015 17:05
 Operator : UM/IZ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICSVBF121315

Quant Time: Dec 14 02:29:15 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 01:30:12 2015
 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	1.773	1.736	2.1	98	-0.01
46	2-Chloronaphthalene	1.308	1.301	0.5	100	-0.01
47	2-Nitroaniline	0.371	0.439	-18.3	101	0.00
48	Acenaphthylene	2.201	2.138	2.9	97	-0.01
49	Dimethylphthalate	1.559	1.513	3.0	99	-0.01
50	2,6-Dinitrotoluene	0.331	0.336	-1.5	100	-0.01
51 C	Acenaphthene	1.331	1.332	-0.1	99	0.00
52	3-Nitroaniline	0.382	0.430	-12.6	98	0.00
53 P	2,4-Dinitrophenol	0.109	0.124	-13.8	102	-0.01
54	Dibenzofuran	1.763	1.688	4.3	98	-0.01
55 P	4-Nitrophenol	0.283	0.344	-21.6	99	0.00
56	2,4-Dinitrotoluene	0.432	0.439	-1.6	100	-0.01
57	Fluorene	1.389	1.293	6.9	99	-0.01
58	2,3,4,6-Tetrachlorophenol	0.307	0.308	-0.3	99	-0.01
59	Diethylphthalate	1.579	1.597	-1.1	98	-0.01
60	4-Chlorophenyl-phenylether	0.633	0.677	-7.0	99	0.00
61	4-Nitroaniline	0.336	0.345	-2.7	101	0.00
62	Azobenzene	1.439	1.518	-5.5	99	-0.01
63 I	Phenanthrene-d10	1.000	1.000	0.0	102	-0.01
64	4,6-Dinitro-2-methylphenol	0.114	0.134	-17.5	100	0.00
65 c	n-Nitrosodiphenylamine	0.684	0.730	-6.7	99	0.00
66	4-Bromophenyl-phenylether	0.224	0.222	0.9	98	-0.01
67	Hexachlorobenzene	0.242	0.235	2.9	98	-0.01
68	Atrazine	0.195	0.230	-17.9	101	0.00
69 C	Pentachlorophenol	0.136	0.153	-12.5	100	0.00
70	Phenanthrene	1.062	1.057	0.5	101	-0.01
71	Anthracene	1.103	1.161	-5.3	100	0.00
72	Carbazole	1.028	1.094	-6.4	99	0.00
73	Di-n-butylphthalate	1.249	1.241	0.6	101	-0.01
74 C	Fluoranthene	1.098	1.103	-0.5	93	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	102	-0.01
76	Benzidine	0.609	0.683	-12.2	91	0.00
77	Pyrene	1.581	1.608	-1.7	98	-0.01
78 S	Terphenyl-d14	0.930	0.881	5.3	97	-0.01
79	Butylbenzylphthalate	0.678	0.730	-7.7	93	-0.01
80	Benzo(a)anthracene	1.137	1.135	0.2	95	-0.01
81	3,3'-Dichlorobenzidine	0.410	0.412	-0.5	91	0.00
82	Chrysene	1.109	1.101	0.7	95	-0.01
83	Bis(2-ethylhexyl)phthalate	0.772	0.793	-2.7	93	-0.01
84 c	Di-n-octyl phthalate	1.227	1.278	-4.2	96	-0.01
85	Indeno(1,2,3-cd)pyrene	1.089	1.127	-3.5	103	-0.02
86 I	Perylene-d12	1.000	1.000	0.0	106	-0.01
87	Benzo(b)fluoranthene	1.276	1.227	3.8	97	-0.01

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
Data File : BF083733.D
Acq On : 13 Dec 2015 17:05
Operator : UM/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF121315

Quant Time: Dec 14 02:29:15 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 14 01:30:12 2015
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	1.156	1.282	-10.9	96	-0.01
89 C	Benzo(a)pyrene	1.139	1.184	-4.0	101	-0.01
90	Dibenzo(a,h)anthracene	1.132	1.229	-8.6	103	-0.01
91	Benzo(a,h,i)perylene	1.164	1.264	-8.6	102	-0.02

(#) = Out of Range

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