

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061522\
 Data File : BG053916.D
 Acq On : 15 Jun 2022 10:03
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 15 17:15:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 15:49:31 2022
 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	93	0.00
2	1,4-Dioxane	0.457	0.439	3.9	88	0.00
3	Pyridine	1.220	1.214	0.5	89	0.00
4	n-Nitrosodimethylamine	0.537	0.577	-7.4	94	0.00
5 S	2-Fluorophenol	1.046	1.033	1.2	90	0.00
6	Aniline	1.682	1.584	5.8	85	0.00
7 S	Phenol-d6	1.419	1.366	3.7	88	0.00
8	2-Chlorophenol	1.117	1.085	2.9	89	0.00
9	Benzaldehyde	0.838	0.741	11.6	88	0.00
10 C	Phenol	1.438	1.375	4.4	85	0.00
11	bis(2-Chloroethyl)ether	1.104	1.045	5.3	88	0.00
12	1,3-Dichlorobenzene	1.400	1.357	3.1	91	0.00
13 C	1,4-Dichlorobenzene	1.406	1.383	1.6	93	0.00
14	1,2-Dichlorobenzene	1.350	1.312	2.8	90	0.00
15	Benzyl Alcohol	1.061	1.024	3.5	86	0.00
16	2,2'-oxybis(1-Chloropropane	1.650	1.487	9.9	83	-0.01
17	2-Methylphenol	0.999	0.928	7.1	85	0.00
18	Hexachloroethane	0.466	0.459	1.5	91	0.00
19 P	n-Nitroso-di-n-propylamine	0.953	0.863	9.4	82	0.00
20	3+4-Methylphenols	1.401	1.280	8.6	83	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
22	Acetophenone	0.486	0.487	-0.2	84	0.00
23 S	Nitrobenzene-d5	0.356	0.377	-5.9	88	0.00
24	Nitrobenzene	0.350	0.373	-6.6	89	0.00
25	Isophorone	0.665	0.645	3.0	82	0.00
26 C	2-Nitrophenol	0.155	0.174	-12.3	91	0.00
27	2,4-Dimethylphenol	0.251	0.253	-0.8	85	0.00
28	bis(2-Chloroethoxy)methane	0.358	0.347	3.1	82	0.00
29 C	2,4-Dichlorophenol	0.336	0.359	-6.8	89	0.00
30	1,2,4-Trichlorobenzene	0.416	0.446	-7.2	90	0.00
31	Naphthalene	1.010	1.004	0.6	84	0.00
32	Benzoic acid	0.087	0.086	1.1	80	0.02
33	4-Chloroaniline	0.419	0.414	1.2	82	0.00
34 C	Hexachlorobutadiene	0.320	0.368	-15.0	96	0.00
35	Caprolactam	0.092	0.090	2.2	78	0.00
36 C	4-Chloro-3-methylphenol	0.327	0.317	3.1	79	0.00
37	2-Methylnaphthalene	0.747	0.739	1.1	83	0.00
38	1-Methylnaphthalene	0.727	0.722	0.7	84	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	84	0.00
40	1,2,4,5-Tetrachlorobenzene	0.771	0.832	-7.9	90	0.00
41 P	Hexachlorocyclopentadiene	0.359	0.426	-18.7	96	0.00
42 S	2,4,6-Tribromophenol	0.321	0.358	-11.5	91	0.00
43 C	2,4,6-Trichlorophenol	0.433	0.467	-7.9	87	0.00
44	2,4,5-Trichlorophenol	0.483	0.509	-5.4	86	0.00
45 S	2-Fluorobiphenyl	1.375	1.401	-1.9	85	0.00
46	1,1'-Biphenyl	1.393	1.373	1.4	82	0.00
47	2-Chloronaphthalene	1.127	1.149	-2.0	85	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061522\
 Data File : BG053916.D
 Acq On : 15 Jun 2022 10:03
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 15 17:15:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 15:49:31 2022
 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	0.300	0.311	-3.7	83	0.00
49	Acenaphthylene	1.753	1.722	1.8	81	0.00
50	Dimethylphthalate	1.523	1.486	2.4	82	0.00
51	2,6-Dinitrotoluene	0.308	0.328	-6.5	86	0.00
52 C	Acenaphthene	1.118	1.134	-1.4	84	0.00
53	3-Nitroaniline	0.298	0.298	0.0	80	0.00
54 P	2,4-Dinitrophenol	0.140	0.161	-15.0	94	0.00
55	Dibenzofuran	1.801	1.799	0.1	83	0.00
56 P	4-Nitrophenol	0.212	0.220	-3.8	83	0.00
57	2,4-Dinitrotoluene	0.435	0.452	-3.9	84	0.00
58	Fluorene	1.478	1.467	0.7	83	0.00
59	2,3,4,6-Tetrachlorophenol	0.473	0.509	-7.6	87	0.00
60	Diethylphthalate	1.475	1.438	2.5	82	0.00
61	4-Chlorophenyl-phenylether	0.885	0.912	-3.1	86	0.00
62	4-Nitroaniline	0.311	0.312	-0.3	82	0.00
63	Azobenzene	1.187	1.147	3.4	82	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
65	4,6-Dinitro-2-methylphenol	0.100	0.112	-12.0	91	0.00
66 c	n-Nitrosodiphenylamine	0.521	0.515	1.2	81	0.00
67	4-Bromophenyl-phenylether	0.250	0.266	-6.4	88	0.00
68	Hexachlorobenzene	0.286	0.305	-6.6	89	0.00
69	Atrazine	0.216	0.230	-6.5	86	0.00
70 C	Pentachlorophenol	0.133	0.151	-13.5	91	0.00
71	Phenanthrene	1.028	1.027	0.1	83	0.00
72	Anthracene	1.003	1.004	-0.1	83	0.00
73	Carbazole	0.861	0.859	0.2	83	0.00
74	Di-n-butylphthalate	0.996	1.003	-0.7	82	0.00
75 C	Fluoranthene	1.349	1.408	-4.4	87	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
77	Benzidine	0.414	0.433	-4.6	97	0.00
78	Pyrene	1.208	1.115	7.7	87	0.00
79 S	Terphenyl-d14	0.993	0.957	3.6	91	0.00
80	Butylbenzylphthalate	0.387	0.381	1.6	91	0.00
81	Benzo(a)anthracene	1.295	1.303	-0.6	96	0.00
82	3,3'-Dichlorobenzidine	0.387	0.399	-3.1	96	0.00
83	Chrysene	1.234	1.230	0.3	96	0.00
84	Bis(2-ethylhexyl)phthalate	0.554	0.543	2.0	91	0.00
85 c	Di-n-octyl phthalate	0.952	0.941	1.2	92	0.00
86 I	Perylene-d12	1.000	1.000	0.0	97	0.00
87	Indeno(1,2,3-cd)pyrene	1.509	1.546	-2.5	97	0.00
88	Benzo(b)fluoranthene	1.216	1.240	-2.0	96	0.00
89	Benzo(k)fluoranthene	1.204	1.195	0.7	97	0.00
90 C	Benzo(a)pyrene	1.026	1.043	-1.7	97	0.00
91	Dibenzo(a,h)anthracene	1.248	1.275	-2.2	98	0.00
92	Benzo(g,h,i)perylene	1.232	1.262	-2.4	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061522\
Data File : BG053916.D
Acq On : 15 Jun 2022 10:03
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 15 17:15:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 15:49:31 2022
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