

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011525\
 Data File : BF141163.D
 Acq On : 15 Jan 2025 09:34
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 15 10:12:02 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 22:54:19 2025
 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	90	0.00
2	1,4-Dioxane	0.577	0.535	7.3	85	0.00
3	Pyridine	1.414	1.322	6.5	87	0.00
4	n-Nitrosodimethylamine	0.691	0.663	4.1	88	0.00
5 S	2-Fluorophenol	1.295	1.263	2.5	91	0.00
6	Aniline	1.456	1.367	6.1	86	0.00
7 S	Phenol-d6	1.640	1.570	4.3	90	0.00
8	2-Chlorophenol	1.389	1.347	3.0	90	0.00
9	Benzaldehyde	0.966	0.866	10.4	92	0.00
10 C	Phenol	1.756	1.667	5.1	88	0.00
11	bis(2-Chloroethyl)ether	1.308	1.229	6.0	88	0.00
12	1,3-Dichlorobenzene	1.549	1.497	3.4	91	0.00
13 C	1,4-Dichlorobenzene	1.561	1.517	2.8	91	0.00
14	1,2-Dichlorobenzene	1.456	1.415	2.8	92	0.00
15	Benzyl Alcohol	1.184	1.160	2.0	90	0.00
16	2,2'-oxybis(1-Chloropropane	1.826	1.648	9.7	84	0.00
17	2-Methylphenol	1.121	1.090	2.8	90	0.00
18	Hexachloroethane	0.565	0.551	2.5	90	0.00
19 P	n-Nitroso-di-n-propylamine	0.966	0.881	8.8	87	0.00
20	3+4-Methylphenols	1.415	1.349	4.7	88	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
22	Acetophenone	0.475	0.457	3.8	92	0.00
23 S	Nitrobenzene-d5	0.377	0.370	1.9	90	0.00
24	Nitrobenzene	0.393	0.385	2.0	91	0.00
25	Isophorone	0.645	0.612	5.1	89	0.00
26 C	2-Nitrophenol	0.179	0.184	-2.8	92	0.00
27	2,4-Dimethylphenol	0.241	0.232	3.7	88	0.00
28	bis(2-Chloroethoxy)methane	0.405	0.380	6.2	88	0.00
29 C	2,4-Dichlorophenol	0.287	0.284	1.0	91	0.00
30	1,2,4-Trichlorobenzene	0.328	0.317	3.4	90	0.00
31	Naphthalene	1.055	1.026	2.7	92	0.00
32	Benzoic acid	0.231	0.241	-4.3	91	0.01
33	4-Chloroaniline	0.365	0.344	5.8	89	0.00
34 C	Hexachlorobutadiene	0.197	0.196	0.5	93	0.00
35	Caprolactam	0.094	0.092	2.1	90	0.00
36 C	4-Chloro-3-methylphenol	0.313	0.306	2.2	91	0.00
37	2-Methylnaphthalene	0.672	0.639	4.9	90	0.00
38	1-Methylnaphthalene	0.662	0.629	5.0	90	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
40	1,2,4,5-Tetrachlorobenzene	0.574	0.557	3.0	91	0.00
41 P	Hexachlorocyclopentadiene	0.188	0.199	-5.9	91	0.00
42 S	2,4,6-Tribromophenol	0.228	0.234	-2.6	97	0.00
43 C	2,4,6-Trichlorophenol	0.387	0.370	4.4	89	0.00
44	2,4,5-Trichlorophenol	0.409	0.404	1.2	92	0.00
45 S	2-Fluorobiphenyl	1.310	1.231	6.0	91	0.00
46	1,1'-Biphenyl	1.553	1.467	5.5	90	0.00
47	2-Chloronaphthalene	1.210	1.166	3.6	91	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011525\
 Data File : BF141163.D
 Acq On : 15 Jan 2025 09:34
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 15 10:12:02 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 22:54:19 2025
 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	0.359	0.353	1.7	92	0.00
49	Acenaphthylene	1.728	1.651	4.5	90	0.00
50	Dimethylphthalate	1.343	1.288	4.1	90	0.00
51	2,6-Dinitrotoluene	0.312	0.308	1.3	90	0.00
52 C	Acenaphthene	1.208	1.176	2.6	92	0.00
53	3-Nitroaniline	0.317	0.324	-2.2	93	0.00
54 P	2,4-Dinitrophenol	0.148	0.159	-7.4	93	0.00
55	Dibenzofuran	1.642	1.566	4.6	90	0.00
56 P	4-Nitrophenol	0.248	0.265	-6.9	97	0.00
57	2,4-Dinitrotoluene	0.402	0.413	-2.7	95	0.00
58	Fluorene	1.276	1.223	4.2	93	0.00
59	2,3,4,6-Tetrachlorophenol	0.340	0.329	3.2	90	0.00
60	Diethylphthalate	1.311	1.242	5.3	90	0.00
61	4-Chlorophenyl-phenylether	0.622	0.595	4.3	91	0.00
62	4-Nitroaniline	0.311	0.328	-5.5	99	0.00
63	Azobenzene	1.293	1.220	5.6	89	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
65	4,6-Dinitro-2-methylphenol	0.116	0.131	-12.9	96	0.00
66 c	n-Nitrosodiphenylamine	0.645	0.629	2.5	92	0.00
67	4-Bromophenyl-phenylether	0.230	0.229	0.4	94	0.00
68	Hexachlorobenzene	0.256	0.257	-0.4	95	0.00
69	Atrazine	0.174	0.143	17.8	96	0.00
70 C	Pentachlorophenol	0.164	0.178	-8.5	97	0.00
71	Phenanthrene	1.074	1.070	0.4	95	0.00
72	Anthracene	1.044	1.044	0.0	95	0.00
73	Carbazole	0.957	0.989	-3.3	99	0.00
74	Di-n-butylphthalate	1.096	1.092	0.4	95	0.00
75 C	Fluoranthene	1.071	1.131	-5.6	103	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	107	0.00
77	Benzidine	0.371	0.551	-48.5#	144	0.00
78	Pyrene	1.910	1.798	5.9	104	0.00
79 S	Terphenyl-d14	1.346	1.266	5.9	107	0.00
80	Butylbenzylphthalate	0.637	0.622	2.4	109	0.00
81	Benzo(a)anthracene	1.361	1.300	4.5	112	0.00
82	3,3'-Dichlorobenzidine	0.433	0.415	4.2	106	0.00
83	Chrysene	1.274	1.151	9.7	102	0.00
84	Bis(2-ethylhexyl)phthalate	0.764	0.720	5.8	106	0.00
85 c	Di-n-octyl phthalate	1.154	0.997	13.6	93	0.00
86 I	Perylene-d12	1.000	1.000	0.0	89	0.00
87	Indeno(1,2,3-cd)pyrene	1.465	1.419	3.1	85	0.00
88	Benzo(b)fluoranthene	1.290	1.198	7.1	88	0.00
89	Benzo(k)fluoranthene	1.090	1.161	-6.5	99	0.00
90 C	Benzo(a)pyrene	1.064	1.042	2.1	89	0.00
91	Dibenzo(a,h)anthracene	1.185	1.151	2.9	85	0.00
92	Benzo(g,h,i)perylene	1.233	1.220	1.1	86	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011525\
Data File : BF141163.D
Acq On : 15 Jan 2025 09:34
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jan 15 10:12:02 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 10 22:54:19 2025
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