

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072425\
 Data File : BP025235.D
 Acq On : 24 Jul 2025 09:54
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 24 10:38:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 22 02:55:21 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	91	0.00
2	1,4-Dioxane	0.444	0.444	0.0	94	0.00
3	Pyridine	1.188	1.160	2.4	90	0.00
4	n-Nitrosodimethylamine	0.471	0.462	1.9	89	0.00
5 S	2-Fluorophenol	1.184	1.179	0.4	92	0.00
6	Aniline	1.916	1.830	4.5	86	0.00
7 S	Phenol-d6	1.466	1.413	3.6	87	0.00
8	2-Chlorophenol	1.343	1.316	2.0	89	0.00
9	Benzaldehyde	0.994	0.976	1.8	80	0.00
10 C	Phenol	1.520	1.460	3.9	87	0.00
11	bis(2-Chloroethyl)ether	1.166	1.120	3.9	87	0.00
12	1,3-Dichlorobenzene	1.510	1.481	1.9	90	0.00
13 C	1,4-Dichlorobenzene	1.525	1.506	1.2	91	0.00
14	1,2-Dichlorobenzene	1.467	1.432	2.4	90	0.00
15	Benzyl Alcohol	1.057	0.993	6.1	85	0.00
16	2,2'-oxybis(1-Chloropropane	1.491	1.409	5.5	86	0.00
17	2-Methylphenol	1.054	1.005	4.6	86	0.00
18	Hexachloroethane	0.524	0.523	0.2	91	0.00
19 P	n-Nitroso-di-n-propylamine	0.882	0.807	8.5	82	0.00
20	3+4-Methylphenols	1.438	1.347	6.3	85	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	82	0.00
22	Acetophenone	0.465	0.470	-1.1	86	0.00
23 S	Nitrobenzene-d5	0.360	0.366	-1.7	85	0.00
24	Nitrobenzene	0.321	0.327	-1.9	85	0.00
25	Isophorone	0.633	0.613	3.2	81	-0.01
26 C	2-Nitrophenol	0.184	0.194	-5.4	87	0.00
27	2,4-Dimethylphenol	0.308	0.306	0.6	83	0.00
28	bis(2-Chloroethoxy)methane	0.396	0.389	1.8	83	0.00
29 C	2,4-Dichlorophenol	0.317	0.316	0.3	83	0.00
30	1,2,4-Trichlorobenzene	0.346	0.349	-0.9	86	0.00
31	Naphthalene	1.023	1.013	1.0	84	0.00
32	Benzoic acid	0.235	0.233	0.9	81	-0.01
33	4-Chloroaniline	0.450	0.439	2.4	82	0.00
34 C	Hexachlorobutadiene	0.208	0.212	-1.9	86	0.00
35	Caprolactam	0.105	0.097	7.6	79	0.00
36 C	4-Chloro-3-methylphenol	0.317	0.305	3.8	82	0.00
37	2-Methylnaphthalene	0.665	0.643	3.3	82	0.00
38	1-Methylnaphthalene	0.698	0.683	2.1	82	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	80	0.00
40	1,2,4,5-Tetrachlorobenzene	0.599	0.609	-1.7	82	0.00
41 P	Hexachlorocyclopentadiene	0.358	0.367	-2.5	83	0.00
42 S	2,4,6-Tribromophenol	0.295	0.289	2.0	80	0.00
43 C	2,4,6-Trichlorophenol	0.410	0.416	-1.5	81	0.00
44	2,4,5-Trichlorophenol	0.451	0.451	0.0	80	0.00
45 S	2-Fluorobiphenyl	1.419	1.457	-2.7	83	0.00
46	1,1'-Biphenyl	1.450	1.441	0.6	80	0.00
47	2-Chloronaphthalene	1.132	1.131	0.1	82	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072425\
 Data File : BP025235.D
 Acq On : 24 Jul 2025 09:54
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 24 10:38:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 22 02:55:21 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	0.289	0.291	-0.7	81	0.00
49	Acenaphthylene	1.831	1.853	-1.2	82	0.00
50	Dimethylphthalate	1.430	1.358	5.0	78	0.00
51	2,6-Dinitrotoluene	0.311	0.307	1.3	79	0.00
52 C	Acenaphthene	1.059	1.038	2.0	80	0.00
53	3-Nitroaniline	0.350	0.345	1.4	80	0.00
54 P	2,4-Dinitrophenol	0.184	0.184	0.0	81	0.00
55	Dibenzofuran	1.710	1.655	3.2	79	0.00
56 P	4-Nitrophenol	0.300	0.285	5.0	78	0.00
57	2,4-Dinitrotoluene	0.435	0.437	-0.5	80	0.00
58	Fluorene	1.358	1.327	2.3	80	0.00
59	2,3,4,6-Tetrachlorophenol	0.399	0.387	3.0	79	0.00
60	Diethylphthalate	1.410	1.362	3.4	79	0.00
61	4-Chlorophenyl-phenylether	0.686	0.669	2.5	79	0.00
62	4-Nitroaniline	0.356	0.349	2.0	80	0.00
63	Azobenzene	1.149	1.121	2.4	80	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	78	0.00
65	4,6-Dinitro-2-methylphenol	0.129	0.135	-4.7	80	0.00
66 c	n-Nitrosodiphenylamine	0.607	0.612	-0.8	80	0.00
67	4-Bromophenyl-phenylether	0.233	0.234	-0.4	79	0.00
68	Hexachlorobenzene	0.277	0.277	0.0	77	0.00
69	Atrazine	0.225	0.225	0.0	80	0.00
70 C	Pentachlorophenol	0.188	0.191	-1.6	79	0.00
71	Phenanthrene	1.079	1.083	-0.4	79	-0.01
72	Anthracene	1.087	1.110	-2.1	80	0.00
73	Carbazole	1.037	1.015	2.1	77	0.00
74	Di-n-butylphthalate	1.197	1.236	-3.3	80	0.00
75 C	Fluoranthene	1.270	1.259	0.9	79	0.01
76 I	Chrysene-d12	1.000	1.000	0.0	79	0.01
77	Benzidine	0.702	0.767	-9.3	72	0.01
78	Pyrene	1.233	1.255	-1.8	79	0.00
79 S	Terphenyl-d14	1.013	1.033	-2.0	92	0.01
80	Butylbenzylphthalate	0.558	0.558	0.0	78	0.00
81	Benzo(a)anthracene	1.263	1.256	0.6	80	0.01
82	3,3'-Dichlorobenzidine	0.524	0.513	2.1	80	0.01
83	Chrysene	1.179	1.185	-0.5	81	0.01
84	Bis(2-ethylhexyl)phthalate	0.798	0.791	0.9	78	0.01
85 c	Di-n-octyl phthalate	1.378	1.388	-0.7	80	0.01
86 I	Perylene-d12	1.000	1.000	0.0	79	0.02
87	Indeno(1,2,3-cd)pyrene	1.462	1.456	0.4	81	0.05
88	Benzo(b)fluoranthene	1.145	1.151	-0.5	80	0.03
89	Benzo(k)fluoranthene	1.130	1.127	0.3	78	0.02
90 C	Benzo(a)pyrene	1.117	1.110	0.6	79	0.03
91	Dibenzo(a,h)anthracene	1.192	1.186	0.5	80	0.05
92	Benzo(g,h,i)perylene	1.172	1.160	1.0	80	0.05

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072425\
Data File : BP025235.D
Acq On : 24 Jul 2025 09:54
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jul 24 10:38:58 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 22 02:55:21 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