

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052025\
 Data File : BF142474.D
 Acq On : 20 May 2025 15:31
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Quant Time: May 20 16:55:40 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 20 16:26:47 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	92899	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	358380	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	188198	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	287337	20.000	ng	0.00
76) Chrysene-d12	14.074	240	154555	20.000	ng	0.00
86) Perylene-d12	15.568	264	196342	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	815861	147.959	ng	0.00
7) Phenol-d6	6.534	99	986649	148.644	ng	0.00
23) Nitrobenzene-d5	7.475	82	993768	151.205	ng	0.00
42) 2,4,6-Tribromophenol	10.739	330	304807	145.927	ng	0.00
45) 2-Fluorobiphenyl	9.269	172	1908541	136.080	ng	0.00
79) Terphenyl-d14	13.015	244	1392704	123.140	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.663	88	169363	76.756	ng	100
3) Pyridine	3.440	79	442247	78.706	ng	98
4) n-Nitrosodimethylamine	3.404	42	230814	79.114	ng	97
6) Aniline	6.569	93	671713	75.263	ng	99
8) 2-Chlorophenol	6.686	128	454577	75.687	ng	98
9) Benzaldehyde	6.451	77	219574	54.999	ng	97
10) Phenol	6.551	94	552574	73.985	ng	87
11) bis(2-Chloroethyl)ether	6.639	93	410501	76.354	ng	100
12) 1,3-Dichlorobenzene	6.845	146	489261	73.004	ng	100
13) 1,4-Dichlorobenzene	6.922	146	495948	73.465	ng	99
14) 1,2-Dichlorobenzene	7.075	146	477210	73.875	ng	100
15) Benzyl Alcohol	7.045	79	381381	77.750	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.181	45	663576	73.497	ng	99
17) 2-Methylphenol	7.151	107	358569	76.467	ng	99
18) Hexachloroethane	7.416	117	177426	75.267	ng	98
19) n-Nitroso-di-n-propyla...	7.328	70	304362	73.987	ng	97
20) 3+4-Methylphenols	7.310	107	426933	71.097	ng	94
22) Acetophenone	7.316	105	571777	72.072	ng	# 95
24) Nitrobenzene	7.492	77	452330	76.309	ng	98
25) Isophorone	7.733	82	838652	76.321	ng	99
26) 2-Nitrophenol	7.804	139	248427	78.434	ng	99
27) 2,4-Dimethylphenol	7.839	122	422403	75.623	ng	99
28) bis(2-Chloroethoxy)met...	7.939	93	513063	74.744	ng	98
29) 2,4-Dichlorophenol	8.045	162	381590	75.121	ng	99
30) 1,2,4-Trichlorobenzene	8.128	180	406857	73.801	ng	98
31) Naphthalene	8.210	128	1277048	72.369	ng	99
32) Benzoic acid	7.975	122	288731	84.918	ng	99
33) 4-Chloroaniline	8.257	127	492168	68.830	ng	100
34) Hexachlorobutadiene	8.328	225	252445	73.338	ng	98
35) Caprolactam	8.645	113	109232	76.565	ng	100
36) 4-Chloro-3-methylphenol	8.739	107	388740	74.674	ng	99
37) 2-Methylnaphthalene	8.898	142	796463	71.742	ng	100
38) 1-Methylnaphthalene	8.998	142	811747	70.689	ng	99
40) 1,2,4,5-Tetrachloroben...	9.069	216	409177	75.740	ng	100
41) Hexachlorocyclopentadiene	9.051	237	309537	84.936	ng	100
43) 2,4,6-Trichlorophenol	9.175	196	285072	78.254	ng	97

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44) 2,4,5-Trichlorophenol	9.216	196	287041	74.712	ng	99
46) 1,1'-Biphenyl	9.369	154	1059787	72.584	ng	100
47) 2-Chloronaphthalene	9.392	162	798539	74.024	ng	99
48) 2-Nitroaniline	9.486	65	241162	77.343	ng	99
49) Acenaphthylene	9.810	152	1313668	72.118	ng	99
50) Dimethylphthalate	9.669	163	911534	73.405	ng	100
51) 2,6-Dinitrotoluene	9.727	165	201545	75.198	ng	96
52) Acenaphthene	9.980	154	810811	72.896	ng	100
53) 3-Nitroaniline	9.898	138	216422	73.930	ng	100
54) 2,4-Dinitrophenol	10.004	184	119082	85.999	ng	94
55) Dibenzofuran	10.151	168	1136276	70.990	ng	99
56) 4-Nitrophenol	10.051	139	165537	76.638	ng	99
57) 2,4-Dinitrotoluene	10.133	165	259035	74.005	ng	99
58) Fluorene	10.498	166	857882	69.377	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.269	232	238382	74.324	ng	99
60) Diethylphthalate	10.369	149	858326	70.773	ng	100
61) 4-Chlorophenyl-phenyle...	10.486	204	426386	70.193	ng	99
62) 4-Nitroaniline	10.516	138	188782	71.105	ng	98
63) Azobenzene	10.645	77	794029	72.893	ng	98
65) 4,6-Dinitro-2-methylph...	10.539	198	141863	88.468	ng	99
66) n-Nitrosodiphenylamine	10.604	169	758366	77.145	ng	99
67) 4-Bromophenyl-phenylether	10.974	248	264739	77.920	ng	98
68) Hexachlorobenzene	11.045	284	287676	76.555	ng	98
69) Atrazine	11.127	200	199929	76.581	ng	99
70) Pentachlorophenol	11.233	266	177504	83.666	ng	99
71) Phenanthrene	11.457	178	1107683	72.134	ng	100
72) Anthracene	11.510	178	1144182	73.009	ng	99
73) Carbazole	11.663	167	956069	71.360	ng	99
74) Di-n-butylphthalate	11.992	149	1043468	71.153	ng	100
75) Fluoranthene	12.645	202	972105	67.750	ng	99
77) Benzidine	12.763	184	343909	59.795	ng	100
78) Pyrene	12.874	202	931844	64.632	ng	100
80) Butylbenzylphthalate	13.492	149	319716	80.252	ng	98
81) Benzo(a)anthracene	14.062	228	765441	74.167	ng	99
82) 3,3'-Dichlorobenzidine	14.027	252	257542	82.274	ng	99
83) Chrysene	14.104	228	715970	77.206	ng	100
84) Bis(2-ethylhexyl)phtha...	14.051	149	449654	88.168	ng	99
85) Di-n-octyl phthalate	14.668	149	888230	89.073	ng	99
87) Indeno(1,2,3-cd)pyrene	17.092	276	1126401	76.417	ng	100
88) Benzo(b)fluoranthene	15.127	252	874640	75.267	ng	98
89) Benzo(k)fluoranthene	15.162	252	803775	73.850	ng	99
90) Benzo(a)pyrene	15.509	252	833783	76.121	ng	99
91) Dibenzo(a,h)anthracene	17.115	278	902757	75.513	ng	99
92) Benzo(g,h,i)perylene	17.556	276	909649	76.073	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

