

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072125\
 Data File : BP025202.D
 Acq On : 21 Jul 2025 18:32
 Operator : CG/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC060

Quant Time: Jul 22 02:33:25 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 02:26:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.437	152	691246	20.000	ng	0.00
21) Naphthalene-d8	10.178	136	2780561	20.000	ng	0.00
39) Acenaphthene-d10	14.078	164	1802837	20.000	ng	0.00
64) Phenanthrene-d10	16.895	188	3596132	20.000	ng	0.00
76) Chrysene-d12	21.330	240	3712729	20.000	ng	0.00
86) Perylene-d12	24.448	264	4449563	20.000	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.102	112	4905756	118.561	ng	0.00
7) Phenol-d6	6.643	99	6110675	119.620	ng	0.00
23) Nitrobenzene-d5	8.572	82	6054858	119.391	ng	0.00
42) 2,4,6-Tribromophenol	15.607	330	3197029	119.256	ng	0.00
45) 2-Fluorobiphenyl	12.678	172	13539099	99.423	ng	0.00
79) Terphenyl-d14	19.648	244	16134390	81.157	ng	-0.01

Target Compounds					Qvalue
2) 1,4-Dioxane	3.119	88	909482	58.484	ng 100
3) Pyridine	3.496	79	2486312	60.365	ng 99
4) n-Nitrosodimethylamine	3.408	42	997579	61.370	ng 99
6) Aniline	6.778	93	4023853	60.573	ng 99
8) 2-Chlorophenol	7.019	128	2831422	60.816	ng 98
9) Benzaldehyde	6.596	77	2060526	56.443	ng 99
10) Phenol	6.666	94	3174369	60.060	ng 100
11) bis(2-Chloroethyl)ether	6.884	93	2426943	59.805	ng 99
12) 1,3-Dichlorobenzene	7.325	146	3115614	58.985	ng 100
13) 1,4-Dichlorobenzene	7.472	146	3138912	58.888	ng 100
14) 1,2-Dichlorobenzene	7.772	146	3018509	58.935	ng 99
15) Benzyl Alcohol	7.672	79	2232586	60.973	ng 99
16) 2,2'-oxybis(1-Chloropr...	7.966	45	3075798	59.089	ng 99
17) 2-Methylphenol	7.878	107	2209131	60.430	ng 100
18) Hexachloroethane	8.484	117	1098658	60.363	ng 100
19) n-Nitroso-di-n-propyla...	8.243	70	1826970	59.342	ng 99
20) 3+4-Methylphenols	8.207	107	3031281	60.708	ng 98
22) Acetophenone	8.243	105	3858483	58.588	ng # 98
24) Nitrobenzene	8.613	77	2713532	60.375	ng 99
25) Isophorone	9.131	82	5301098	59.683	ng 99
26) 2-Nitrophenol	9.307	139	1614578	63.823	ng 99
27) 2,4-Dimethylphenol	9.384	122	2618557	60.802	ng 100
28) bis(2-Chloroethoxy)met...	9.613	93	3327518	59.673	ng 100
29) 2,4-Dichlorophenol	9.837	162	2686031	60.808	ng 99
30) 1,2,4-Trichlorobenzene	10.054	180	2894308	59.549	ng 100
31) Naphthalene	10.225	128	8471283	58.501	ng 100
32) Benzoic acid	9.560	122	2149321	68.334	ng 98
33) 4-Chloroaniline	10.342	127	3853838	61.295	ng 100
34) Hexachlorobutadiene	10.519	225	1757479	60.210	ng 99
35) Caprolactam	11.148	113	911643	62.713	ng 99
36) 4-Chloro-3-methylphenol	11.484	107	2701662	61.242	ng 100
37) 2-Methylnaphthalene	11.860	142	5538936	59.140	ng 100
38) 1-Methylnaphthalene	12.066	142	5796319	58.678	ng 99
40) 1,2,4,5-Tetrachloroben...	12.236	216	3229034	59.145	ng 100
41) Hexachlorocyclopentadiene	12.219	237	2121694	68.531	ng 100
43) 2,4,6-Trichlorophenol	12.484	196	2251566	61.035	ng 100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.560	196	2484748	61.107	ng	100
46) 1,1'-Biphenyl	12.895	154	7671161	57.545	ng	100
47) 2-Chloronaphthalene	12.931	162	6074581	58.817	ng	100
48) 2-Nitroaniline	13.136	65	1603764	61.905	ng	98
49) Acenaphthylene	13.789	152	9744261	57.536	ng	99
50) Dimethylphthalate	13.542	163	7577608	57.607	ng	100
51) 2,6-Dinitrotoluene	13.660	165	1738822	62.447	ng	98
52) Acenaphthene	14.142	154	5579143	57.185	ng	99
53) 3-Nitroaniline	13.983	138	1947927	62.002	ng	96
54) 2,4-Dinitrophenol	14.189	184	1119889	71.772	ng	99
55) Dibenzofuran	14.489	168	8968837	56.778	ng	98
56) 4-Nitrophenol	14.325	139	1695550	63.331	ng	99
57) 2,4-Dinitrotoluene	14.460	165	2469068	63.650	ng	98
58) Fluorene	15.154	166	6997947	55.543	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.730	232	2196650	61.026	ng	100
60) Diethylphthalate	14.948	149	7449753	57.319	ng	99
61) 4-Chlorophenyl-phenyle...	15.160	204	3554251	55.969	ng	99
62) 4-Nitroaniline	15.183	138	1978662	61.804	ng	98
63) Azobenzene	15.460	77	6251887	59.884	ng	100
65) 4,6-Dinitro-2-methylph...	15.242	198	1488428	66.114	ng	99
66) n-Nitrosodiphenylamine	15.378	169	6429114	57.405	ng	99
67) 4-Bromophenyl-phenylether	16.072	248	2534895	59.785	ng	99
68) Hexachlorobenzene	16.189	284	2976346	58.786	ng	96
69) Atrazine	16.366	200	2433839	59.621	ng	99
70) Pentachlorophenol	16.542	266	2118802	62.871	ng	99
71) Phenanthrene	16.942	178	11295817	56.373	ng	99
72) Anthracene	17.030	178	11494384	56.508	ng	98
73) Carbazole	17.319	167	10836403	55.990	ng	98
74) Di-n-butylphthalate	17.936	149	12352965	54.431	ng	# 94
75) Fluoranthene	19.048	202	12783070	53.956	ng	93
77) Benzidine	19.242	184	6968232	49.133	ng	100
78) Pyrene	19.424	202	13220717	56.032	ng	93
80) Butylbenzylphthalate	20.413	149	6306883	60.572	ng	99
81) Benzo(a)anthracene	21.307	228	13600779	56.319	ng	98
82) 3,3'-Dichlorobenzidine	21.254	252	5842950	59.005	ng	100
83) Chrysene	21.371	228	12774794	56.724	ng	98
84) Bis(2-ethylhexyl)phtha...	21.289	149	8682877	57.201	ng	99
85) Di-n-octyl phthalate	22.507	149	15085392	57.015	ng	99
87) Indeno(1,2,3-cd)pyrene	27.994	276	19493176	59.597	ng	100
88) Benzo(b)fluoranthene	23.477	252	14989282	57.763	ng	100
89) Benzo(k)fluoranthene	23.542	252	14928017	57.942	ng	99
90) Benzo(a)pyrene	24.306	252	14791063	58.720	ng	99
91) Dibenzo(a,h)anthracene	28.059	278	15890395	59.240	ng	99
92) Benzo(g,h,i)perylene	29.071	276	15521510	59.072	ng	98

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

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