

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070325\
 Data File : BP025073.D
 Acq On : 03 Jul 2025 11:02
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Quant Time: Jul 03 15:51:01 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 03 15:38:19 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.449	152	342992	20.000	ng	0.00
21) Naphthalene-d8	10.196	136	1376560	20.000	ng	0.01
39) Acenaphthene-d10	14.096	164	924609	20.000	ng	0.00
64) Phenanthrene-d10	16.913	188	1970410	20.000	ng	0.00
76) Chrysene-d12	21.342	240	2172955	20.000	ng	0.00
86) Perylene-d12	24.489	264	2220570	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.102	112	376746	19.462	ng	0.00
7) Phenol-d6	6.631	99	482699	19.108	ng	0.00
23) Nitrobenzene-d5	8.566	82	381978	18.454	ng	0.00
42) 2,4,6-Tribromophenol	15.619	330	198235	18.149	ng	0.00
45) 2-Fluorobiphenyl	12.701	172	1211432	20.297	ng	0.00
79) Terphenyl-d14	19.672	244	1997838	20.207	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.126	88	77575	10.306	ng 98
3) Pyridine	3.490	79	197662	9.598	ng 97
4) n-Nitrosodimethylamine	3.408	42	85827	9.836	ng # 92
6) Aniline	6.790	93	227023	9.619	ng 98
8) 2-Chlorophenol	7.014	128	206980	9.556	ng 96
9) Benzaldehyde	6.602	77	154921	11.743	ng 98
10) Phenol	6.655	94	250593	9.763	ng 99
11) bis(2-Chloroethyl)ether	6.890	93	195263	9.885	ng 98
12) 1,3-Dichlorobenzene	7.343	146	245857	10.143	ng 98
13) 1,4-Dichlorobenzene	7.478	146	245523	10.004	ng 98
14) 1,2-Dichlorobenzene	7.790	146	237961	10.079	ng 98
15) Benzyl Alcohol	7.678	79	152168	9.025	ng 99
16) 2,2'-oxybis(1-Chloropr...	7.978	45	260708	10.150	ng 97
17) 2-Methylphenol	7.878	107	155778	9.422	ng 98
18) Hexachloroethane	8.502	117	75576	9.453	ng 96
19) n-Nitroso-di-n-propyla...	8.237	70	146424	9.774	ng 97
20) 3+4-Methylphenols	8.196	107	213641	9.225	ng 97
22) Acetophenone	8.243	105	330725	10.159	ng 98
24) Nitrobenzene	8.608	77	201551	9.560	ng 99
25) Isophorone	9.137	82	367740	9.507	ng 100
26) 2-Nitrophenol	9.313	139	63769	9.106	ng 92
27) 2,4-Dimethylphenol	9.390	122	134296	9.307	ng 97
28) bis(2-Chloroethoxy)met...	9.625	93	255336	9.722	ng 99
29) 2,4-Dichlorophenol	9.843	162	181263	9.059	ng 99
30) 1,2,4-Trichlorobenzene	10.066	180	219650	9.913	ng 99
31) Naphthalene	10.243	128	709141	10.005	ng 99
32) Benzoic acid	9.455	122	79225	10.455	ng 95
33) 4-Chloroaniline	10.343	127	254489	9.621	ng 98
34) Hexachlorobutadiene	10.549	225	124190	9.767	ng 99
35) Caprolactam	11.084	113	59547	8.700	ng 94
36) 4-Chloro-3-methylphenol	11.478	107	192224	9.193	ng 99
37) 2-Methylnaphthalene	11.872	142	489608	9.790	ng 99
38) 1-Methylnaphthalene	12.084	142	484403	9.931	ng 100
40) 1,2,4,5-Tetrachloroben...	12.249	216	253871	9.554	ng 99
41) Hexachlorocyclopentadiene	12.243	237	55283	7.965	ng 96
43) 2,4,6-Trichlorophenol	12.490	196	136269	8.469	ng 99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.554	196	161200	8.840	ng	98
46) 1,1'-Biphenyl	12.913	154	663919	9.869	ng	99
47) 2-Chloronaphthalene	12.949	162	503580	9.959	ng	98
48) 2-Nitroaniline	13.143	65	85669	9.283	ng	97
49) Acenaphthylene	13.807	152	750027	9.692	ng	99
50) Dimethylphthalate	13.554	163	636102	9.828	ng	99
51) 2,6-Dinitrotoluene	13.660	165	110177	8.728	ng	95
52) Acenaphthene	14.160	154	496424m	9.953	ng	
53) 3-Nitroaniline	13.990	138	114723	8.481	ng	# 95
54) 2,4-Dinitrophenol	14.201	184	35399	10.737	ng	96
55) Dibenzofuran	14.507	168	838464	10.186	ng	100
56) 4-Nitrophenol	14.301	139	98038	7.853	ng	92
57) 2,4-Dinitrotoluene	14.466	165	131885	9.365	ng	98
58) Fluorene	15.172	166	639683	10.047	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.731	232	146800	8.890	ng	97
60) Diethylphthalate	14.960	149	636288	9.788	ng	98
61) 4-Chlorophenyl-phenyle...	15.184	204	317737	10.084	ng	98
62) 4-Nitroaniline	15.178	138	123493	8.706	ng	93
63) Azobenzene	15.478	77	589360	10.181	ng	97
65) 4,6-Dinitro-2-methylph...	15.243	198	52978	10.800	ng	87
66) n-Nitrosodiphenylamine	15.390	169	539820	9.573	ng	99
67) 4-Bromophenyl-phenylether	16.090	248	186736	9.256	ng	93
68) Hexachlorobenzene	16.195	284	236838	9.604	ng	98
69) Atrazine	16.384	200	164479	9.954	ng	100
70) Pentachlorophenol	16.554	266	129459	7.774	ng	98
71) Phenanthrene	16.960	178	1068228	10.146	ng	99
72) Anthracene	17.042	178	980773	9.710	ng	99
73) Carbazole	17.331	167	979377	9.917	ng	99
74) Di-n-butylphthalate	17.966	149	979540	8.907	ng	100
75) Fluoranthene	19.060	202	1203752	9.781	ng	99
77) Benzidine	19.266	184	192008	5.003	ng	99
78) Pyrene	19.442	202	1286495	9.559	ng	99
80) Butylbenzylphthalate	20.430	149	300028	9.024	ng	95
81) Benzo(a)anthracene	21.325	228	1276106	9.691	ng	99
82) 3,3'-Dichlorobenzidine	21.272	252	270931	6.834	ng	99
83) Chrysene	21.389	228	1246003	9.832	ng	99
84) Bis(2-ethylhexyl)phtha...	21.325	149	564998	7.999	ng	97
85) Di-n-octyl phthalate	22.554	149	662011	10.561	ng	97
87) Indeno(1,2,3-cd)pyrene	28.042	276	1222863	8.642	ng	98
88) Benzo(b)fluoranthene	23.495	252	1180548	9.134	ng	100
89) Benzo(k)fluoranthene	23.560	252	1252989	9.743	ng	99
90) Benzo(a)pyrene	24.342	252	963455	8.758	ng	99
91) Dibenzo(a,h)anthracene	28.118	278	1022797	8.747	ng	100
92) Benzo(g,h,i)perylene	29.095	276	1073371	8.886	ng	99

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

