

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060625\
 Data File : BP024866.D
 Acq On : 06 Jun 2025 14:37
 Operator : RC/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC060

Quant Time: Jun 06 16:04:20 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 15:53:04 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.608	152	233459	20.000	ng	0.00	
21) Naphthalene-d8	10.378	136	1006719	20.000	ng	0.00	
39) Acenaphthene-d10	14.248	164	639758	20.000	ng	0.00	
64) Phenanthrene-d10	17.042	188	1223735	20.000	ng	-0.02	
76) Chrysene-d12	21.477	240	1245531	20.000	ng	0.00	
86) Perylene-d12	24.736	264	1494809	20.000	ng	0.02	
System Monitoring Compounds							
5) 2-Fluorophenol	5.237	112	1754536	125.447	ng	0.00	
7) Phenol-d6	6.814	99	2365958	127.853	ng	0.00	
23) Nitrobenzene-d5	8.760	82	2555584	123.355	ng	0.00	
42) 2,4,6-Tribromophenol	15.772	330	1097449	124.076	ng	-0.01	
45) 2-Fluorobiphenyl	12.860	172	5620380	118.347	ng	0.00	
79) Terphenyl-d14	19.778	244	8373144	120.485	ng	-0.02	
Target Compounds							
2) 1,4-Dioxane	3.167	88	360857	58.637	ng	100	Qvalue
3) Pyridine	3.561	79	957340	64.685	ng	99	
4) n-Nitrosodimethylamine	3.478	42	387978	64.122	ng	99	
6) Aniline	6.955	93	1528129	64.708	ng	100	
8) 2-Chlorophenol	7.190	128	995913	62.813	ng	99	
9) Benzaldehyde	6.766	77	608861	56.921	ng	99	
10) Phenol	6.843	94	1231710	64.565	ng	98	
11) bis(2-Chloroethyl)ether	7.043	93	926134	61.724	ng	98	
12) 1,3-Dichlorobenzene	7.496	146	1063861	60.147	ng	99	
13) 1,4-Dichlorobenzene	7.643	146	1074203	60.190	ng	100	
14) 1,2-Dichlorobenzene	7.955	146	1045089	59.634	ng	99	
15) Benzyl Alcohol	7.861	79	916921	64.561	ng	99	
16) 2,2'-oxybis(1-Chloropr...	8.125	45	1167705	59.460	ng	100	
17) 2-Methylphenol	8.066	107	848195	63.670	ng	100	
18) Hexachloroethane	8.666	117	401855	59.807	ng	96	
19) n-Nitroso-di-n-propyla...	8.413	70	779593	61.970	ng	100	
20) 3+4-Methylphenols	8.396	107	1154457	63.519	ng	96	
22) Acetophenone	8.431	105	1540416	60.563	ng	99	
24) Nitrobenzene	8.808	77	1136226	61.735	ng	99	
25) Isophorone	9.325	82	2191187	61.027	ng	100	
26) 2-Nitrophenol	9.507	139	587760	64.723	ng	99	
27) 2,4-Dimethylphenol	9.578	122	965935	62.151	ng	100	
28) bis(2-Chloroethoxy)met...	9.802	93	1278529	59.694	ng	98	
29) 2,4-Dichlorophenol	10.049	162	945641	63.413	ng	99	
30) 1,2,4-Trichlorobenzene	10.243	180	995211	59.169	ng	99	
31) Naphthalene	10.431	128	3124640	60.566	ng	100	
32) Benzoic acid	9.778	122	710052	68.525	ng	100	
33) 4-Chloroaniline	10.549	127	1397378	64.650	ng	98	
34) Hexachlorobutadiene	10.707	225	598677	59.056	ng	100	
35) Caprolactam	11.360	113	349412	63.654	ng	93	
36) 4-Chloro-3-methylphenol	11.696	107	1096315	63.738	ng	98	
37) 2-Methylnaphthalene	12.043	142	2004422	61.250	ng	99	
38) 1-Methylnaphthalene	12.266	142	2093346	59.832	ng	99	
40) 1,2,4,5-Tetrachloroben...	12.419	216	1100767	60.637	ng	100	
41) Hexachlorocyclopentadiene	12.390	237	707348	63.884	ng	98	
43) 2,4,6-Trichlorophenol	12.672	196	776210	63.669	ng	99	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.760	196	836590	64.077	ng	99
46) 1,1'-Biphenyl	13.066	154	2797746	60.176	ng	100
47) 2-Chloronaphthalene	13.107	162	2180558	61.023	ng	99
48) 2-Nitroaniline	13.331	65	718204	65.376	ng	97
49) Acenaphthylene	13.966	152	3653383	61.318	ng	100
50) Dimethylphthalate	13.707	163	2826983	59.902	ng	99
51) 2,6-Dinitrotoluene	13.831	165	639722	62.878	ng	96
52) Acenaphthene	14.313	154	2051309	60.082	ng	99
53) 3-Nitroaniline	14.172	138	698945	66.199	ng	98
54) 2,4-Dinitrophenol	14.390	184	398250	61.497	ng	97
55) Dibenzofuran	14.654	168	3267254	59.618	ng	99
56) 4-Nitrophenol	14.525	139	539162	61.650	ng	97
57) 2,4-Dinitrotoluene	14.637	165	901515	63.345	ng	98
58) Fluorene	15.319	166	2628887	59.381	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.895	232	730647	62.524	ng	99
60) Diethylphthalate	15.090	149	2772340	58.943	ng	98
61) 4-Chlorophenyl-phenyle...	15.307	204	1276716	58.980	ng	100
62) 4-Nitroaniline	15.360	138	639038	66.714	ng	96
63) Azobenzene	15.613	77	2561443	59.380	ng	98
65) 4,6-Dinitro-2-methylph...	15.419	198	524098	65.178	ng	100
66) n-Nitrosodiphenylamine	15.531	169	2282932	60.188	ng	99
67) 4-Bromophenyl-phenylether	16.225	248	839674	60.812	ng	98
68) Hexachlorobenzene	16.342	284	1008320	60.238	ng	97
69) Atrazine	16.519	200	836217	60.814	ng	99
70) Pentachlorophenol	16.701	266	561603	64.826	ng	99
71) Phenanthrene	17.089	178	4046968	59.844	ng	99
72) Anthracene	17.189	178	4160775	60.751	ng	99
73) Carbazole	17.472	167	3861413	60.827	ng	99
74) Di-n-butylphthalate	18.060	149	4672145	59.401	ng	100
75) Fluoranthene	19.178	202	4656349	59.407	ng	99
77) Benzidine	19.383	184	2477600	64.227	ng	100
78) Pyrene	19.560	202	4751590	61.064	ng	100
80) Butylbenzylphthalate	20.513	149	2225855	62.449	ng	96
81) Benzo(a)anthracene	21.460	228	4896757	61.469	ng	100
82) 3,3'-Dichlorobenzidine	21.389	252	1982344	62.681	ng	100
83) Chrysene	21.530	228	4625245	61.276	ng	99
84) Bis(2-ethylhexyl)phtha...	21.395	149	3103331	60.769	ng	99
85) Di-n-octyl phthalate	22.636	149	5491065	61.004	ng	100
87) Indeno(1,2,3-cd)pyrene	28.436	276	6771619	62.088	ng	100
88) Benzo(b)fluoranthene	23.713	252	5290333	61.842	ng	99
89) Benzo(k)fluoranthene	23.777	252	5191049	59.612	ng	99
90) Benzo(a)pyrene	24.595	252	5093340	60.965	ng	99
91) Dibenzo(a,h)anthracene	28.495	278	5489005	61.830	ng	100
92) Benzo(g,h,i)perylene	29.571	276	5442983	61.794	ng	100

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

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