

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061925\
 Data File : BP025007.D
 Acq On : 19 Jun 2025 19:08
 Operator : RC/JU
 Sample : Q2347-01MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-10MSD

Quant Time: Jun 20 01:13:54 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 16:20:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.602	152	256103	20.000	ng	-0.01
21) Naphthalene-d8	10.372	136	1009701	20.000	ng	0.00
39) Acenaphthene-d10	14.248	164	615822	20.000	ng	0.00
64) Phenanthrene-d10	17.054	188	1249087	20.000	ng	0.00
76) Chrysene-d12	21.495	240	1308170	20.000	ng	0.01
86) Perylene-d12	24.765	264	1527853	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	1617814	105.444	ng	0.00
7) Phenol-d6	6.813	99	2087706	102.842	ng	0.00
23) Nitrobenzene-d5	8.754	82	1269394	61.091	ng	0.00
42) 2,4,6-Tribromophenol	15.783	330	928676	109.075	ng	0.00
45) 2-Fluorobiphenyl	12.854	172	2748203	60.117	ng	0.00
79) Terphenyl-d14	19.783	244	4387058	60.105	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.161	88	261359	38.714	ng	99
3) Pyridine	3.561	79	674552	41.548	ng	99
4) n-Nitrosodimethylamine	3.472	42	288015	43.392	ng	93
6) Aniline	6.943	93	751711	29.016	ng	100
8) 2-Chlorophenol	7.184	128	820397	47.168	ng	100
9) Benzaldehyde	6.760	77	535651	45.779	ng	100
10) Phenol	6.837	94	987188	47.172	ng	97
11) bis(2-Chloroethyl)ether	7.037	93	727221	44.182	ng	100
12) 1,3-Dichlorobenzene	7.490	146	851172	43.867	ng	99
13) 1,4-Dichlorobenzene	7.637	146	857924	43.821	ng	99
14) 1,2-Dichlorobenzene	7.949	146	829329	43.138	ng	99
15) Benzyl Alcohol	7.854	79	704875	45.243	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.119	45	938524	43.564	ng	99
17) 2-Methylphenol	8.066	107	671577	45.955	ng	98
18) Hexachloroethane	8.654	117	325767	44.197	ng	96
19) n-Nitroso-di-n-propyla...	8.401	70	575544	41.705	ng	100
20) 3+4-Methylphenols	8.396	107	896635	44.971	ng	97
22) Acetophenone	8.425	105	1142688	44.793	ng	# 99
24) Nitrobenzene	8.801	77	852776	46.197	ng	98
25) Isophorone	9.319	82	1608274	44.660	ng	100
26) 2-Nitrophenol	9.501	139	441229	48.444	ng	99
27) 2,4-Dimethylphenol	9.578	122	737667	47.323	ng	98
28) bis(2-Chloroethoxy)met...	9.796	93	976691	45.467	ng	98
29) 2,4-Dichlorophenol	10.060	162	739099	49.416	ng	99
30) 1,2,4-Trichlorobenzene	10.237	180	763077	45.234	ng	100
31) Naphthalene	10.419	128	2353821	45.490	ng	99
32) Benzoic acid	9.784	122	575259	54.612	ng	99
33) 4-Chloroaniline	10.554	127	483590	22.307	ng	99
34) Hexachlorobutadiene	10.695	225	465748	45.808	ng	99
35) Caprolactam	11.360	113	269335	48.921	ng	96
36) 4-Chloro-3-methylphenol	11.707	107	828727	48.038	ng	99
37) 2-Methylnaphthalene	12.042	142	1505667	45.874	ng	99
38) 1-Methylnaphthalene	12.266	142	1576082	44.915	ng	98
40) 1,2,4,5-Tetrachloroben...	12.419	216	844665	48.338	ng	99
41) Hexachlorocyclopentadiene	12.384	237	844810	79.264	ng	99
43) 2,4,6-Trichlorophenol	12.684	196	600238	51.149	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.778	196	658524	52.398	ng	98
46) 1,1'-Biphenyl	13.072	154	2135117	47.709	ng	99
47) 2-Chloronaphthalene	13.113	162	1628681	47.350	ng	99
48) 2-Nitroaniline	13.342	65	530296	50.148	ng	95
49) Acenaphthylene	13.966	152	2727641	47.560	ng	100
50) Dimethylphthalate	13.707	163	2131155	46.913	ng	99
51) 2,6-Dinitrotoluene	13.842	165	481655	49.182	ng	96
52) Acenaphthene	14.319	154	1574892	47.921	ng	99
53) 3-Nitroaniline	14.189	138	292308	28.762	ng	99
54) 2,4-Dinitrophenol	14.413	184	647775	100.079	ng	96
55) Dibenzofuran	14.660	168	2485168	47.109	ng	100
56) 4-Nitrophenol	14.560	139	735984	85.165	ng	96
57) 2,4-Dinitrotoluene	14.654	165	687020	50.150	ng	95
58) Fluorene	15.319	166	2041782	47.912	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.907	232	569986	50.761	ng	99
60) Diethylphthalate	15.089	149	2175202	48.045	ng	99
61) 4-Chlorophenyl-phenyle...	15.307	204	992505	47.632	ng	97
62) 4-Nitroaniline	15.372	138	482604	52.371	ng	93
63) Azobenzene	15.613	77	2253120	54.262	ng	98
65) 4,6-Dinitro-2-methylph...	15.436	198	404357	49.266	ng	98
66) n-Nitrosodiphenylamine	15.542	169	1818455	46.969	ng	100
67) 4-Bromophenyl-phenylether	16.230	248	654553	46.443	ng	99
68) Hexachlorobenzene	16.354	284	779037	45.596	ng	96
69) Atrazine	16.525	200	670358	47.763	ng	99
70) Pentachlorophenol	16.713	266	997318	112.784	ng	99
71) Phenanthrene	17.101	178	3181572	46.092	ng	99
72) Anthracene	17.195	178	3335317	47.710	ng	99
73) Carbazole	17.483	167	3132092	48.337	ng	100
74) Di-n-butylphthalate	18.060	149	3936416	49.032	ng	99
75) Fluoranthene	19.189	202	3829480	47.866	ng	99
77) Benzidine	19.395	184	1908036	47.093	ng	100
78) Pyrene	19.566	202	3981423	48.716	ng	100
80) Butylbenzylphthalate	20.519	149	1878758	50.187	ng	96
81) Benzo(a)anthracene	21.477	228	3927926	46.947	ng	99
82) 3,3'-Dichlorobenzidine	21.401	252	1083650	32.624	ng	100
83) Chrysene	21.542	228	3732223	47.078	ng	99
84) Bis(2-ethylhexyl)phtha...	21.401	149	2704452	50.423	ng	99
85) Di-n-octyl phthalate	22.636	149	4660187	49.294	ng	99
87) Indeno(1,2,3-cd)pyrene	28.495	276	5453869	48.924	ng	# 91
88) Benzo(b)fluoranthene	23.730	252	4168657	47.675	ng	98
89) Benzo(k)fluoranthene	23.801	252	4054783	45.557	ng	98
90) Benzo(a)pyrene	24.612	252	4026608	47.154	ng	99
91) Dibenzo(a,h)anthracene	28.559	278	4443108	48.966	ng	99
92) Benzo(g,h,i)perylene	29.647	276	4438842	49.304	ng	98

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

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