

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061825\
 Data File : BP024995.D
 Acq On : 18 Jun 2025 15:26
 Operator : RC/JU
 Sample : Q2311-04MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP03-MH2MH3-WCMS

Quant Time: Jun 18 15:48:34 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	242047	20.000	ng	-0.01
21) Naphthalene-d8	10.372	136	965848	20.000	ng	0.00
39) Acenaphthene-d10	14.248	164	614166	20.000	ng	0.00
64) Phenanthrene-d10	17.060	188	1216080	20.000	ng	0.00
76) Chrysene-d12	21.501	240	1258513	20.000	ng	0.02
86) Perylene-d12	24.771	264	1511485	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	1970717	135.904	ng	0.00
7) Phenol-d6	6.814	99	2335784	121.744	ng	0.00
23) Nitrobenzene-d5	8.761	82	1683911	84.720	ng	0.00
42) 2,4,6-Tribromophenol	15.778	330	1231452	145.027	ng	0.00
45) 2-Fluorobiphenyl	12.854	172	3713121	81.444	ng	0.00
79) Terphenyl-d14	19.789	244	6194695	88.219	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.173	88	233736	36.633	ng	# 58
3) Pyridine	3.573	79	607536	39.593	ng	99
4) n-Nitrosodimethylamine	3.478	42	263396	41.987	ng	95
6) Aniline	6.949	93	613124	25.041	ng	99
8) 2-Chlorophenol	7.184	128	773411	47.049	ng	100
9) Benzaldehyde	6.761	77	417644	37.766	ng	100
10) Phenol	6.843	94	822820	41.601	ng	98
11) bis(2-Chloroethyl)ether	7.037	93	714257	45.914	ng	99
12) 1,3-Dichlorobenzene	7.496	146	782730	42.683	ng	99
13) 1,4-Dichlorobenzene	7.637	146	798490	43.154	ng	100
14) 1,2-Dichlorobenzene	7.949	146	766159	42.167	ng	99
15) Benzyl Alcohol	7.855	79	689279	46.811	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.131	45	903934	44.396	ng	99
17) 2-Methylphenol	8.072	107	637440	46.152	ng	98
18) Hexachloroethane	8.661	117	298119	42.794	ng	98
19) n-Nitroso-di-n-propyla...	8.408	70	569886	43.693	ng	99
20) 3+4-Methylphenols	8.402	107	851471	45.186	ng	96
22) Acetophenone	8.425	105	1120345	45.911	ng	99
24) Nitrobenzene	8.802	77	834079	47.236	ng	99
25) Isophorone	9.319	82	1622372	47.097	ng	99
26) 2-Nitrophenol	9.502	139	426264	48.926	ng	98
27) 2,4-Dimethylphenol	9.578	122	735556	49.330	ng	98
28) bis(2-Chloroethoxy)met...	9.796	93	987853	48.074	ng	99
29) 2,4-Dichlorophenol	10.060	162	730516	51.060	ng	99
30) 1,2,4-Trichlorobenzene	10.243	180	722759	44.789	ng	98
31) Naphthalene	10.425	128	2255528	45.570	ng	100
32) Benzoic acid	9.766	122	306657	30.434	ng	94
33) 4-Chloroaniline	10.566	127	325962	15.719	ng	100
34) Hexachlorobutadiene	10.696	225	433953	44.618	ng	98
35) Caprolactam	11.354	113	219676	41.713	ng	97
36) 4-Chloro-3-methylphenol	11.707	107	832967	50.476	ng	100
37) 2-Methylnaphthalene	12.043	142	1450864	46.211	ng	99
38) 1-Methylnaphthalene	12.266	142	1532263	45.649	ng	99
40) 1,2,4,5-Tetrachloroben...	12.425	216	804930	46.188	ng	100
41) Hexachlorocyclopentadiene	12.390	237	835920	78.641	ng	99
43) 2,4,6-Trichlorophenol	12.678	196	589260	50.349	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.778	196	638845	50.970	ng	99
46) 1,1'-Biphenyl	13.066	154	2127848	47.674	ng	99
47) 2-Chloronaphthalene	13.113	162	1599570	46.629	ng	99
48) 2-Nitroaniline	13.343	65	538789	51.088	ng	96
49) Acenaphthylene	13.972	152	2703658	47.269	ng	100
50) Dimethylphthalate	13.707	163	2311406	51.018	ng	100
51) 2,6-Dinitrotoluene	13.837	165	491960	50.370	ng	99
52) Acenaphthene	14.313	154	1561197	47.633	ng	99
53) 3-Nitroaniline	14.190	138	257325	25.388	ng	99
54) 2,4-Dinitrophenol	14.413	184	479264	75.680	ng	95
55) Dibenzofuran	14.660	168	2493777	47.400	ng	98
56) 4-Nitrophenol	14.560	139	629869	73.850	ng	91
57) 2,4-Dinitrotoluene	14.654	165	701358	51.335	ng	96
58) Fluorene	15.319	166	2057394	48.409	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.907	232	580077	51.799	ng	99
60) Diethylphthalate	15.095	149	2311971	51.204	ng	99
61) 4-Chlorophenyl-phenyle...	15.313	204	1001717	48.204	ng	99
62) 4-Nitroaniline	15.372	138	461293	50.193	ng	94
63) Azobenzene	15.613	77	2064593	49.856	ng	99
65) 4,6-Dinitro-2-methylph...	15.437	198	367659	46.011	ng	94
66) n-Nitrosodiphenylamine	15.543	169	1838746	48.782	ng	99
67) 4-Bromophenyl-phenylether	16.231	248	662017	48.247	ng	96
68) Hexachlorobenzene	16.354	284	790416	47.517	ng	96
69) Atrazine	16.525	200	705811	51.653	ng	99
70) Pentachlorophenol	16.725	266	972643	112.979	ng	99
71) Phenanthrene	17.101	178	3205790	47.704	ng	99
72) Anthracene	17.195	178	3290810	48.351	ng	99
73) Carbazole	17.489	167	3156761	50.040	ng	100
74) Di-n-butylphthalate	18.066	149	4078294	52.178	ng	100
75) Fluoranthene	19.195	202	3807278	48.880	ng	100
77) Benzidine	19.401	184	1005765	25.803	ng	100
78) Pyrene	19.572	202	3945102	50.176	ng	100
80) Butylbenzylphthalate	20.519	149	1942979	53.950	ng	98
81) Benzo(a)anthracene	21.477	228	3990097	49.571	ng	99
82) 3,3'-Dichlorobenzidine	21.407	252	959911	30.039	ng	99
83) Chrysene	21.548	228	3719434	48.768	ng	100
84) Bis(2-ethylhexyl)phtha...	21.407	149	2808178	54.422	ng	99
85) Di-n-octyl phthalate	22.654	149	4901102	53.888	ng	100
87) Indeno(1,2,3-cd)pyrene	28.483	276	5450032	49.419	ng	# 86
88) Benzo(b)fluoranthene	23.736	252	4207871	48.644	ng	99
89) Benzo(k)fluoranthene	23.807	252	4337889	49.265	ng	100
90) Benzo(a)pyrene	24.618	252	4110605	48.659	ng	99
91) Dibenzo(a,h)anthracene	28.554	278	4445031	49.518	ng	99
92) Benzo(g,h,i)perylene	29.659	276	4368895	49.053	ng	99

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

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