

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061925\
 Data File : BP025006.D
 Acq On : 19 Jun 2025 18:26
 Operator : RC/JU
 Sample : Q2347-01MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-10MS

Quant Time: Jun 19 19:06:11 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	300654	20.000	ng	-0.01
21) Naphthalene-d8	10.372	136	1153563	20.000	ng	0.00
39) Acenaphthene-d10	14.248	164	701293	20.000	ng	0.00
64) Phenanthrene-d10	17.060	188	1369804	20.000	ng	0.00
76) Chrysene-d12	21.495	240	1442401	20.000	ng	0.01
86) Perylene-d12	24.777	264	1737125	20.000	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	1880226	104.388	ng	0.00
7) Phenol-d6	6.814	99	2383086	99.997	ng	0.00
23) Nitrobenzene-d5	8.755	82	1472434	62.025	ng	0.00
42) 2,4,6-Tribromophenol	15.778	330	1014848	104.669	ng	0.00
45) 2-Fluorobiphenyl	12.854	172	3122979	59.990	ng	0.00
79) Terphenyl-d14	19.789	244	4814936	59.828	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.161	88	296522	37.414	ng	98
3) Pyridine	3.561	79	777316	40.783	ng	100
4) n-Nitrosodimethylamine	3.473	42	336607	43.198	ng	96
6) Aniline	6.949	93	848790	27.909	ng	100
8) 2-Chlorophenol	7.184	128	946291	46.344	ng	100
9) Benzaldehyde	6.761	77	617006	44.918	ng	99
10) Phenol	6.837	94	1140925	46.440	ng	97
11) bis(2-Chloroethyl)ether	7.037	93	861256	44.572	ng	98
12) 1,3-Dichlorobenzene	7.490	146	1002153	43.995	ng	98
13) 1,4-Dichlorobenzene	7.637	146	1012546	44.055	ng	99
14) 1,2-Dichlorobenzene	7.949	146	970357	42.995	ng	100
15) Benzyl Alcohol	7.855	79	806234	44.080	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.125	45	1101320	43.546	ng	99
17) 2-Methylphenol	8.066	107	761518	44.388	ng	98
18) Hexachloroethane	8.655	117	380405	43.962	ng	96
19) n-Nitroso-di-n-propyla...	8.408	70	670951	41.414	ng	99
20) 3+4-Methylphenols	8.396	107	1016806	43.442	ng	96
22) Acetophenone	8.425	105	1316378	45.166	ng	# 99
24) Nitrobenzene	8.802	77	970948	46.039	ng	99
25) Isophorone	9.319	82	1839066	44.700	ng	99
26) 2-Nitrophenol	9.502	139	512126	49.216	ng	97
27) 2,4-Dimethylphenol	9.578	122	827479	46.465	ng	97
28) bis(2-Chloroethoxy)met...	9.796	93	1116189	45.480	ng	# 97
29) 2,4-Dichlorophenol	10.060	162	841675	49.257	ng	99
30) 1,2,4-Trichlorobenzene	10.237	180	875913	45.447	ng	99
31) Naphthalene	10.425	128	2707445	45.799	ng	99
32) Benzoic acid	9.790	122	655593	54.476	ng	97
33) 4-Chloroaniline	10.555	127	574847	23.210	ng	99
34) Hexachlorobutadiene	10.702	225	541179	46.589	ng	99
35) Caprolactam	11.360	113	301748	47.973	ng	95
36) 4-Chloro-3-methylphenol	11.707	107	929186	47.144	ng	98
37) 2-Methylnaphthalene	12.043	142	1715743	45.755	ng	100
38) 1-Methylnaphthalene	12.266	142	1791651	44.691	ng	99
40) 1,2,4,5-Tetrachloroben...	12.425	216	944045	47.441	ng	99
41) Hexachlorocyclopentadiene	12.390	237	979716	80.718	ng	98
43) 2,4,6-Trichlorophenol	12.684	196	670333	50.160	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061925\
 Data File : BP025006.D
 Acq On : 19 Jun 2025 18:26
 Operator : RC/JU
 Sample : Q2347-01MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-10MS

Quant Time: Jun 19 19:06:11 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)
44) 2,4,5-Trichlorophenol	12.778	196	737605	51.538	ng	98
46) 1,1'-Biphenyl	13.072	154	2403102	47.152	ng	100
47) 2-Chloronaphthalene	13.113	162	1851737	47.274	ng	100
48) 2-Nitroaniline	13.337	65	581837	48.316	ng	96
49) Acenaphthylene	13.972	152	3018623	46.219	ng	100
50) Dimethylphthalate	13.713	163	2389367	46.186	ng	100
51) 2,6-Dinitrotoluene	13.843	165	530728	47.588	ng	96
52) Acenaphthene	14.319	154	1753441	46.852	ng	100
53) 3-Nitroaniline	14.184	138	335156	28.958	ng	98
54) 2,4-Dinitrophenol	14.413	184	712859	96.898	ng	94
55) Dibenzofuran	14.660	168	2748969	45.759	ng	99
56) 4-Nitrophenol	14.554	139	821107	83.545	ng	97
57) 2,4-Dinitrotoluene	14.654	165	758309	48.607	ng	# 92
58) Fluorene	15.325	166	2256542	46.498	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.907	232	629507	49.229	ng	98
60) Diethylphthalate	15.095	149	2453677	47.591	ng	98
61) 4-Chlorophenyl-phenyle...	15.313	204	1110242	46.789	ng	98
62) 4-Nitroaniline	15.378	138	516167	49.187	ng	91
63) Azobenzene	15.613	77	2509594	53.073	ng	98
65) 4,6-Dinitro-2-methylph...	15.437	198	444463	49.380	ng	95
66) n-Nitrosodiphenylamine	15.537	169	1975658	46.533	ng	99
67) 4-Bromophenyl-phenylether	16.231	248	722139	46.723	ng	97
68) Hexachlorobenzene	16.354	284	859782	45.887	ng	95
69) Atrazine	16.525	200	733050	47.626	ng	99
70) Pentachlorophenol	16.719	266	1102604	113.702	ng	99
71) Phenanthrene	17.101	178	3483796	46.023	ng	99
72) Anthracene	17.195	178	3603129	46.999	ng	100
73) Carbazole	17.484	167	3354983	47.214	ng	100
74) Di-n-butylphthalate	18.066	149	4351849	49.429	ng	100
75) Fluoranthene	19.195	202	4114293	46.894	ng	99
77) Benzidine	19.401	184	1981239	44.350	ng	100
78) Pyrene	19.572	202	4265160	47.331	ng	99
80) Butylbenzylphthalate	20.525	149	2096919	50.802	ng	96
81) Benzo(a)anthracene	21.477	228	4317005	46.795	ng	100
82) 3,3'-Dichlorobenzidine	21.407	252	1194213	32.607	ng	99
83) Chrysene	21.548	228	4112171	47.043	ng	100
84) Bis(2-ethylhexyl)phtha...	21.407	149	3041809	51.435	ng	99
85) Di-n-octyl phthalate	22.648	149	5272581	50.582	ng	100
87) Indeno(1,2,3-cd)pyrene	28.501	276	6063935	47.844	ng	# 90
88) Benzo(b)fluoranthene	23.736	252	4665279m	46.927	ng	
89) Benzo(k)fluoranthene	23.801	252	4698731	46.432	ng	99
90) Benzo(a)pyrene	24.618	252	4613412	47.518	ng	99
91) Dibenzo(a,h)anthracene	28.559	278	4947335	47.955	ng	100
92) Benzo(g,h,i)perylene	29.636	276	4857239	47.452	ng	98

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

