

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
 Data File : BF142914.D
 Acq On : 30 Jun 2025 13:17
 Operator : RC/JU
 Sample : Q2430-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-E/FMSD

Quant Time: Jun 30 13:49:55 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 24 05:28:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	110394	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	429125	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	237255	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	439577	20.000	ng	0.00
76) Chrysene-d12	14.063	240	267586	20.000	ng	0.00
86) Perylene-d12	15.551	264	284395	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	589585	88.919	ng	0.03
7) Phenol-d6	6.522	99	718865	91.894	ng	0.01
23) Nitrobenzene-d5	7.445	82	446908	57.124	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	251453	102.991	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	964853	53.424	ng	0.00
79) Terphenyl-d14	12.998	244	1063719	54.926	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.828	88	97687	36.834	ng	98
3) Pyridine	3.569	79	262762	39.521	ng	96
4) n-Nitrosodimethylamine	3.534	42	144546	42.042	ng	99
6) Aniline	6.545	93	302665	29.078	ng	# 69
8) 2-Chlorophenol	6.669	128	317052	44.327	ng	98
9) Benzaldehyde	6.434	77	132287	28.503	ng	100
10) Phenol	6.534	94	365970	42.087	ng	83
11) bis(2-Chloroethyl)ether	6.616	93	272133	43.787	ng	98
12) 1,3-Dichlorobenzene	6.822	146	336218	42.031	ng	100
13) 1,4-Dichlorobenzene	6.898	146	339085	42.015	ng	99
14) 1,2-Dichlorobenzene	7.051	146	324004	42.326	ng	99
15) Benzyl Alcohol	7.028	79	258053	45.630	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.151	45	393052	39.364	ng	71
17) 2-Methylphenol	7.139	107	240647	44.398	ng	95
18) Hexachloroethane	7.392	117	121020	42.871	ng	96
19) n-Nitroso-di-n-propyla...	7.298	70	205016	45.061	ng	97
20) 3+4-Methylphenols	7.292	107	303469	44.904	ng	99
22) Acetophenone	7.292	105	412283	43.568	ng	# 98
24) Nitrobenzene	7.463	77	302074	42.660	ng	99
25) Isophorone	7.704	82	561857	44.769	ng	100
26) 2-Nitrophenol	7.781	139	178819	46.358	ng	99
27) 2,4-Dimethylphenol	7.822	122	292307	43.748	ng	99
28) bis(2-Chloroethoxy)met...	7.910	93	348800	43.716	ng	100
29) 2,4-Dichlorophenol	8.028	162	270738	44.689	ng	99
30) 1,2,4-Trichlorobenzene	8.104	180	285079	42.793	ng	99
31) Naphthalene	8.186	128	898275	42.765	ng	100
32) Benzoic acid	7.957	122	144109	42.317	ng	97
33) 4-Chloroaniline	8.234	127	117242	13.727	ng	99
34) Hexachlorobutadiene	8.298	225	175332	43.028	ng	99
35) Caprolactam	8.628	113	78994m	49.413	ng	
36) 4-Chloro-3-methylphenol	8.722	107	285073	47.347	ng	98
37) 2-Methylnaphthalene	8.875	142	569691	43.747	ng	99
38) 1-Methylnaphthalene	8.981	142	591674	43.891	ng	99
40) 1,2,4,5-Tetrachloroben...	9.045	216	292815	41.827	ng	98
41) Hexachlorocyclopentadiene	9.028	237	307393	77.376	ng	99
43) 2,4,6-Trichlorophenol	9.157	196	198396	43.491	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.204	196	212273	44.205	ng	99
46) 1,1'-Biphenyl	9.345	154	779647	41.599	ng	100
47) 2-Chloronaphthalene	9.369	162	589551	41.918	ng	99
48) 2-Nitroaniline	9.469	65	180886	45.875	ng	95
49) Acenaphthylene	9.786	152	977274	42.209	ng	100
50) Dimethylphthalate	9.651	163	721236	46.963	ng	100
51) 2,6-Dinitrotoluene	9.710	165	158408	46.968	ng	97
52) Acenaphthene	9.957	154	633829m	44.869	ng	
53) 3-Nitroaniline	9.880	138	106442	28.665	ng	99
54) 2,4-Dinitrophenol	9.992	184	110371	65.356	ng	97
55) Dibenzofuran	10.133	168	856627	42.289	ng	100
56) 4-Nitrophenol	10.057	139	268103	105.463	ng	97
57) 2,4-Dinitrotoluene	10.122	165	217368	48.521	ng	# 88
58) Fluorene	10.475	166	690003	43.616	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.251	232	180513	46.635	ng	97
60) Diethylphthalate	10.345	149	734120	48.753	ng	99
61) 4-Chlorophenyl-phenyle...	10.463	204	329879	43.698	ng	98
62) 4-Nitroaniline	10.504	138	176766	52.050	ng	98
63) Azobenzene	10.622	77	654963	48.667	ng	99
65) 4,6-Dinitro-2-methylph...	10.533	198	101246	38.079	ng	98
66) n-Nitrosodiphenylamine	10.586	169	622140	40.837	ng	99
67) 4-Bromophenyl-phenylether	10.957	248	210753	42.127	ng	99
68) Hexachlorobenzene	11.022	284	232514	41.827	ng	99
69) Atrazine	11.116	200	213613	54.298	ng	98
70) Pentachlorophenol	11.222	266	229534	82.864	ng	99
71) Phenanthrene	11.439	178	1002078	42.083	ng	99
72) Anthracene	11.492	178	1029114	42.140	ng	99
73) Carbazole	11.645	167	959921	45.549	ng	99
74) Di-n-butylphthalate	11.969	149	1187601	54.094	ng	100
75) Fluoranthene	12.627	202	1052967	46.594	ng	99
77) Benzidine	12.745	184	317334	31.564	ng	100
78) Pyrene	12.857	202	1050214	41.871	ng	99
80) Butylbenzylphthalate	13.468	149	445864	61.282	ng	99
81) Benzo(a)anthracene	14.051	228	793975	43.984	ng	100
82) 3,3'-Dichlorobenzidine	14.010	252	107304	18.326	ng	99
83) Chrysene	14.086	228	707737	43.938	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	570793	54.528	ng	99
85) Di-n-octyl phthalate	14.645	149	844081	42.763	ng	99
87) Indeno(1,2,3-cd)pyrene	17.068	276	655902	30.279	ng	98
88) Benzo(b)fluoranthene	15.115	252	740550	45.001	ng	100
89) Benzo(k)fluoranthene	15.145	252	727830	47.273	ng	99
90) Benzo(a)pyrene	15.492	252	717112	45.107	ng	100
91) Dibenzo(a,h)anthracene	17.086	278	540782	30.725	ng	98
92) Benzo(g,h,i)perylene	17.527	276	451541	25.727	ng	98

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

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