

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111025\
 Data File : BG064732.D
 Acq On : 10 Nov 2025 14:07
 Operator : RC/JU
 Sample : Q3562-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EO-03-110625MSD

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 11/11/2025
 Supervised By :Jagrut Upadhyay 11/11/2025

Quant Time: Nov 10 14:45:01 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 03 18:16:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.206	152	42939	20.000	ng	-0.02
21) Naphthalene-d8	11.032	136	168015	20.000	ng	#-0.02
39) Acenaphthene-d10	14.828	164	131298	20.000	ng	-0.02
64) Phenanthrene-d10	17.566	188	287909	20.000	ng	-0.02
76) Chrysene-d12	21.843	240	317067	20.000	ng	#-0.02
86) Perylene-d12	25.180	264	351711	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.738	112	157381	61.521	ng	-0.01
7) Phenol-d6	7.348	99	215146	61.289	ng	0.00
23) Nitrobenzene-d5	9.369	82	145395	51.976	ng	-0.02
42) 2,4,6-Tribromophenol	16.308	330	161107	85.016	ng	-0.01
45) 2-Fluorobiphenyl	13.459	172	400011	46.179	ng	-0.02
79) Terphenyl-d14	20.151	244	760736	45.279	ng	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.564	88	27939	24.669	ng	100
3) Pyridine	3.976	79	81511	23.966	ng	98
4) n-Nitrosodimethylamine	3.887	42	55211	30.271	ng	95
6) Aniline	7.519	93	72736	15.227	ng	99
8) 2-Chlorophenol	7.765	128	116852	43.640	ng	97
9) Benzaldehyde	7.330	77	65303	34.075	ng	96
10) Phenol	7.377	94	157387	40.541	ng	96
11) bis(2-Chloroethyl)ether	7.607	93	101631	35.607	ng	88
12) 1,3-Dichlorobenzene	8.094	146	130056	42.200	ng	99
13) 1,4-Dichlorobenzene	8.241	146	134715	42.990	ng	94
14) 1,2-Dichlorobenzene	8.564	146	132940	43.983	ng	96
15) Benzyl Alcohol	8.435	79	115801	40.004	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.729	45	248666	33.213	ng	98
17) 2-Methylphenol	8.641	107	110102	42.716	ng	96
18) Hexachloroethane	9.305	117	51421	46.760	ng	89
19) n-Nitroso-di-n-propyla...	9.011	70	91621	37.766	ng	92
20) 3+4-Methylphenols	8.970	107	147897	40.801	ng	96
22) Acetophenone	9.028	105	191927	41.788	ng	94
24) Nitrobenzene	9.416	77	140741	46.258	ng	94
25) Isophorone	9.939	82	254019	37.844	ng	99
26) 2-Nitrophenol	10.133	139	69681	58.617	ng	# 93
27) 2,4-Dimethylphenol	10.186	122	120027	48.072	ng	98
28) bis(2-Chloroethoxy)met...	10.415	93	144956	37.929	ng	99
29) 2,4-Dichlorophenol	10.668	162	137504	51.173	ng	97
30) 1,2,4-Trichlorobenzene	10.891	180	165611	54.245	ng	95
31) Naphthalene	11.085	128	407526	43.748	ng	100
32) Benzoic acid	10.327	122	85556	45.338	ng	93
33) 4-Chloroaniline	11.179	127	57865	15.982	ng	97
34) Hexachlorobutadiene	11.373	225	133859	56.064	ng	94
35) Caprolactam	11.954	113	38291m	37.154	ng	
36) 4-Chloro-3-methylphenol	12.289	107	146511	45.361	ng	96
37) 2-Methylnaphthalene	12.677	142	275575	40.293	ng	99
38) 1-Methylnaphthalene	12.895	142	286379	42.298	ng	96
40) 1,2,4,5-Tetrachloroben...	13.036	216	231530	52.543	ng	98
41) Hexachlorocyclopentadiene	13.018	237	122898	61.964	ng	92
43) 2,4,6-Trichlorophenol	13.271	196	139804	55.466	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111025\
 Data File : BG064732.D
 Acq On : 10 Nov 2025 14:07
 Operator : RC/JU
 Sample : Q3562-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EO-03-110625MSD

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 11/11/2025
 Supervised By :Jagrut Upadhyay 11/11/2025

Quant Time: Nov 10 14:45:01 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 03 18:16:26 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.341	196	153087	52.631	ng	94
46) 1,1'-Biphenyl	13.670	154	421935	45.202	ng	96
47) 2-Chloronaphthalene	13.711	162	319400	45.331	ng	99
48) 2-Nitroaniline	13.899	65	108101	45.856	ng	96
49) Acenaphthylene	14.557	152	556743	50.095	ng	99
50) Dimethylphthalate	14.275	163	417333	41.538	ng	100
51) 2,6-Dinitrotoluene	14.387	165	87860	52.667	ng	94
52) Acenaphthene	14.892	154	348658	45.918	ng	98
53) 3-Nitroaniline	14.722	138	61566	35.657	ng	# 93
54) 2,4-Dinitrophenol	14.922	184	77444	76.401	ng	# 36
55) Dibenzofuran	15.227	168	532222	43.356	ng	99
56) 4-Nitrophenol	15.016	139	139316	87.028	ng	91
57) 2,4-Dinitrotoluene	15.174	165	125370	48.855	ng	# 93
58) Fluorene	15.873	166	410434	42.812	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.445	232	154053	55.386	ng	96
60) Diethylphthalate	15.633	149	402116	39.199	ng	98
61) 4-Chlorophenyl-phenyle...	15.862	204	244320	45.412	ng	96
62) 4-Nitroaniline	15.873	138	75550	39.331	ng	90
63) Azobenzene	16.150	77	364286	38.703	ng	97
65) 4,6-Dinitro-2-methylph...	15.938	198	61533	48.506	ng	96
66) n-Nitrosodiphenylamine	16.067	169	362884	46.660	ng	98
67) 4-Bromophenyl-phenylether	16.749	248	176506	51.734	ng	97
68) Hexachlorobenzene	16.878	284	208669	53.610	ng	97
69) Atrazine	17.007	200	166417	49.724	ng	98
70) Pentachlorophenol	17.213	266	250808	103.332	ng	99
71) Phenanthrene	17.613	178	766837	48.705	ng	100
72) Anthracene	17.701	178	731474	45.672	ng	100
73) Carbazole	17.959	167	586704	41.542	ng	100
74) Di-n-butylphthalate	18.512	149	661745	40.519	ng	99
75) Fluoranthene	19.604	202	1056870	53.583	ng	97
77) Benzidine	19.769	184	131027	18.445	ng	96
78) Pyrene	19.969	202	1071671	51.733	ng	99
80) Butylbenzylphthalate	20.832	149	285276	45.564	ng	95
81) Benzo(a)anthracene	21.819	228	1130975	53.796	ng	99
82) 3,3'-Dichlorobenzidine	21.725	252	163431	23.280	ng	98
83) Chrysene	21.890	228	1043300	52.154	ng	97
84) Bis(2-ethylhexyl)phtha...	21.714	149	408760	43.692	ng	# 96
85) Di-n-octyl phthalate	22.971	149	715969	45.392	ng	96
87) Indeno(1,2,3-cd)pyrene	29.023	276	946868	36.597	ng	# 81
88) Benzo(b)fluoranthene	24.123	252	1078174	49.992	ng	# 96
89) Benzo(k)fluoranthene	24.193	252	1157843	53.649	ng	# 95
90) Benzo(a)pyrene	25.027	252	1060123	53.475	ng	# 97
91) Dibenzo(a,h)anthracene	29.087	278	744532	35.420	ng	99
92) Benzo(g,h,i)perylene	30.198	276	653722	32.556	ng	# 96

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111025\
Data File : BG064732.D
Acq On : 10 Nov 2025 14:07
Operator : RC/JU
Sample : Q3562-01MSD
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-03-110625MSD

Quant Time: Nov 10 14:45:01 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 03 18:16:26 2025
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Rahul Chavli 11/11/2025
Supervised By :Jagrut Upadhyay 11/11/2025

