

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111225\
 Data File : BG064747.D
 Acq On : 12 Nov 2025 12:39
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 11/14/2025

Supervised By :Sohil Jodhani 11/14/2025

Quant Time: Nov 12 16:55:23 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Nov 12 16:45:57 2025

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.164	152	28063	20.000	ng	0.00
21) Naphthalene-d8	10.985	136	110132	20.000	ng	0.00
39) Acenaphthene-d10	14.774	164	90796	20.000	ng	0.00
64) Phenanthrene-d10	17.506	188	239033	20.000	ng	0.00
76) Chrysene-d12	21.754	240	366018	20.000	ng	0.00
86) Perylene-d12	25.021	264	382667	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.703	112	31652	19.695	ng	0.00
7) Phenol-d6	7.307	99	41135	19.245	ng	0.00
23) Nitrobenzene-d5	9.328	82	41577	18.666	ng	0.00
42) 2,4,6-Tribromophenol	16.249	330	34875	19.694	ng	0.00
45) 2-Fluorobiphenyl	13.405	172	131226	20.167	ng	0.00
79) Terphenyl-d14	20.092	244	376175	20.217	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.529	88	6388	10.268	ng	97
3) Pyridine	3.946	79	18201	9.819	ng	92
4) n-Nitrosodimethylamine	3.852	42	9221	9.773	ng	# 98
6) Aniline	7.477	93	28080	9.856	ng	98
8) 2-Chlorophenol	7.730	128	17126	9.705	ng	94
9) Benzaldehyde	7.295	77	14410	10.299	ng	96
10) Phenol	7.330	94	22227	9.557	ng	94
11) bis(2-Chloroethyl)ether	7.571	93	16009	9.674	ng	96
12) 1,3-Dichlorobenzene	8.059	146	19968	9.834	ng	98
13) 1,4-Dichlorobenzene	8.206	146	20506	9.918	ng	98
14) 1,2-Dichlorobenzene	8.523	146	19237	9.596	ng	96
15) Benzyl Alcohol	8.394	79	14805	8.438	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.687	45	39228	10.014	ng	96
17) 2-Methylphenol	8.599	107	16118	9.795	ng	97
18) Hexachloroethane	9.263	117	7708	9.843	ng	98
19) n-Nitroso-di-n-propyla...	8.969	70	14240	9.853	ng	94
20) 3+4-Methylphenols	8.928	107	20773	9.215	ng	97
22) Acetophenone	8.987	105	29535	9.843	ng	93
24) Nitrobenzene	9.369	77	21900	9.730	ng	99
25) Isophorone	9.898	82	41875	9.872	ng	96
26) 2-Nitrophenol	10.086	139	9640	9.523	ng	95
27) 2,4-Dimethylphenol	10.139	122	16597	9.228	ng	93
28) bis(2-Chloroethoxy)met...	10.368	93	23070	9.907	ng	97
29) 2,4-Dichlorophenol	10.626	162	18228	9.440	ng	98
30) 1,2,4-Trichlorobenzene	10.844	180	22952	9.844	ng	99
31) Naphthalene	11.032	128	64016	10.254	ng	98
32) Benzoic acid	10.233	122	7448m	10.914	ng	
33) 4-Chloroaniline	11.132	127	24891	9.967	ng	97
34) Hexachlorobutadiene	11.325	225	19432	9.713	ng	97
35) Caprolactam	11.878	113	4958	8.592	ng	# 92
36) 4-Chloro-3-methylphenol	12.236	107	21366	9.942	ng	95
37) 2-Methylnaphthalene	12.624	142	46349	9.921	ng	97
38) 1-Methylnaphthalene	12.841	142	46327m	9.982	ng	
40) 1,2,4,5-Tetrachloroben...	12.988	216	33859	9.582	ng	96
41) Hexachlorocyclopentadiene	12.971	237	21693	8.565	ng	91
43) 2,4,6-Trichlorophenol	13.217	196	19600	9.451	ng	95

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111225\
 Data File : BG064747.D
 Acq On : 12 Nov 2025 12:39
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 11/14/2025
 Supervised By :Sohil Jodhani 11/14/2025

Quant Time: Nov 12 16:55:23 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Nov 12 16:45:57 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.288	196	20868	8.972	ng	93
46) 1,1'-Biphenyl	13.617	154	64329	9.864	ng	97
47) 2-Chloronaphthalene	13.664	162	50487	9.827	ng	98
48) 2-Nitroaniline	13.852	65	15648	9.469	ng	87
49) Acenaphthylene	14.498	152	87532	9.973	ng	99
50) Dimethylphthalate	14.216	163	71209	9.928	ng	97
51) 2,6-Dinitrotoluene	14.334	165	12760	9.246	ng	92
52) Acenaphthene	14.839	154	50131	9.526	ng	98
53) 3-Nitroaniline	14.663	138	12586	9.240	ng	87
54) 2,4-Dinitrophenol	14.874	184	2717	11.149	ng	# 84
55) Dibenzofuran	15.168	168	90369	10.387	ng	97
56) 4-Nitrophenol	14.951	139	8225	7.611	ng	# 83
57) 2,4-Dinitrotoluene	15.115	165	19317	9.336	ng	95
58) Fluorene	15.814	166	71244	10.441	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.391	232	21790	9.495	ng	# 97
60) Diethylphthalate	15.568	149	69072	10.053	ng	97
61) 4-Chlorophenyl-phenyle...	15.803	204	42098	10.280	ng	98
62) 4-Nitroaniline	15.814	138	13078	9.268	ng	95
63) Azobenzene	16.090	77	58632	10.260	ng	98
65) 4,6-Dinitro-2-methylph...	15.873	198	8710	6.563	ng	92
66) n-Nitrosodiphenylamine	16.014	169	63405	10.127	ng	96
67) 4-Bromophenyl-phenylether	16.696	248	31456	10.021	ng	96
68) Hexachlorobenzene	16.813	284	37251	9.795	ng	93
69) Atrazine	16.948	200	30006	9.940	ng	96
70) Pentachlorophenol	17.154	266	10561	10.475	ng	95
71) Phenanthrene	17.548	178	133643	10.219	ng	99
72) Anthracene	17.636	178	134138	10.059	ng	99
73) Carbazole	17.900	167	117923	10.286	ng	99
74) Di-n-butylphthalate	18.452	149	132248	10.340	ng	99
75) Fluoranthene	19.539	202	186154	10.409	ng	99
77) Benzidine	19.710	184	123305	11.249	ng	96
78) Pyrene	19.898	202	199370	9.719	ng	97
80) Butylbenzylphthalate	20.773	149	71488	9.908	ng	95
81) Benzo(a)anthracene	21.737	228	250849	10.079	ng	99
82) 3,3'-Dichlorobenzidine	21.643	252	91769	10.018	ng	100
83) Chrysene	21.801	228	236517	10.080	ng	99
84) Bis(2-ethylhexyl)phtha...	21.637	149	106405	10.092	ng	97
85) Di-n-octyl phthalate	22.865	149	181658	9.967	ng	99
87) Indeno(1,2,3-cd)pyrene	28.740	276	253326	9.636	ng	# 98
88) Benzo(b)fluoranthene	23.981	252	239260	9.825	ng	99
89) Benzo(k)fluoranthene	24.046	252	248008	10.137	ng	100
90) Benzo(a)pyrene	24.863	252	224991	9.960	ng	98
91) Dibenzo(a,h)anthracene	28.805	278	208966	9.702	ng	# 96
92) Benzo(g,h,i)perylene	29.904	276	199793	9.639	ng	# 96

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

Manual Integrations

Quant Time: Nov 12 16:55:23 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 12 16:45:57 2025
Response via : Initial Calibration

