

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111225\
 Data File : BG064748.D
 Acq On : 12 Nov 2025 13:20
 Operator : RC/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC020

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 11/14/2025

Supervised By :Sohil Jodhani 11/14/2025

Quant Time: Nov 12 16:56:17 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	8.167	152	30983	20.000	ng	0.00
21) Naphthalene-d8	10.982	136	122144	20.000	ng	0.00
39) Acenaphthene-d10	14.777	164	99587	20.000	ng	0.00
64) Phenanthrene-d10	17.504	188	258312	20.000	ng	0.00
76) Chrysene-d12	21.757	240	361266	20.000	ng	0.00
86) Perylene-d12	25.018	264	380612	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.700	112	72167	40.674	ng	0.00
7) Phenol-d6	7.304	99	95345	40.402	ng	0.00
23) Nitrobenzene-d5	9.325	82	100042	40.496	ng	0.00
42) 2,4,6-Tribromophenol	16.246	330	81708	42.069	ng	0.00
45) 2-Fluorobiphenyl	13.408	172	300532	42.108	ng	0.00
79) Terphenyl-d14	20.089	244	766056	41.712	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.532	88	13788	20.073	ng	95
3) Pyridine	3.943	79	40729	19.901	ng	97
4) n-Nitrosodimethylamine	3.849	42	20577	19.754	ng	100
6) Aniline	7.480	93	65205	20.729	ng	96
8) 2-Chlorophenol	7.727	128	39290	20.167	ng	97
9) Benzaldehyde	7.292	77	30372	19.661	ng	96
10) Phenol	7.333	94	51616	20.103	ng	97
11) bis(2-Chloroethyl)ether	7.574	93	37916	20.752	ng	98
12) 1,3-Dichlorobenzene	8.056	146	46058	20.545	ng	96
13) 1,4-Dichlorobenzene	8.203	146	47387	20.759	ng	96
14) 1,2-Dichlorobenzene	8.526	146	44745	20.216	ng	96
15) Benzyl Alcohol	8.397	79	38065	19.651	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.696	45	91675	21.198	ng	98
17) 2-Methylphenol	8.596	107	37068	20.403	ng	93
18) Hexachloroethane	9.266	117	17722	20.498	ng	97
19) n-Nitroso-di-n-propyla...	8.967	70	34572	21.667	ng	94
20) 3+4-Methylphenols	8.925	107	50448	20.269	ng	88
22) Acetophenone	8.984	105	70914	21.310	ng	99
24) Nitrobenzene	9.366	77	52662	21.097	ng	93
25) Isophorone	9.895	82	97191	20.659	ng	95
26) 2-Nitrophenol	10.089	139	23040	20.521	ng	96
27) 2,4-Dimethylphenol	10.136	122	41015	20.563	ng	97
28) bis(2-Chloroethoxy)met...	10.371	93	54247	21.004	ng	100
29) 2,4-Dichlorophenol	10.623	162	44768	20.905	ng	95
30) 1,2,4-Trichlorobenzene	10.847	180	54086	20.916	ng	99
31) Naphthalene	11.035	128	142862	20.633	ng	99
32) Benzoic acid	10.218	122	20309m	18.429	ng	
33) 4-Chloroaniline	11.129	127	58239	21.026	ng	98
34) Hexachlorobutadiene	11.323	225	45189	20.367	ng	98
35) Caprolactam	11.875	113	12678	19.811	ng	95
36) 4-Chloro-3-methylphenol	12.233	107	48730	20.444	ng	98
37) 2-Methylnaphthalene	12.627	142	110208	21.269	ng	99
38) 1-Methylnaphthalene	12.844	142	107491	20.884	ng	96
40) 1,2,4,5-Tetrachloroben...	12.991	216	80735	20.831	ng	99
41) Hexachlorocyclopentadiene	12.974	237	54509	19.622	ng	98
43) 2,4,6-Trichlorophenol	13.214	196	46989	20.658	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.291	196	54340	21.300	ng	95
46) 1,1'-Biphenyl	13.614	154	149187	20.857	ng	97
47) 2-Chloronaphthalene	13.661	162	116198	20.620	ng	97
48) 2-Nitroaniline	13.849	65	37146	20.495	ng	97
49) Acenaphthylene	14.501	152	198809	20.651	ng	98
50) Dimethylphthalate	14.219	163	164729	20.940	ng	99
51) 2,6-Dinitrotoluene	14.331	165	31143	20.573	ng	96
52) Acenaphthene	14.836	154	116463	20.176	ng	97
53) 3-Nitroaniline	14.666	138	29942	20.041	ng	95
54) 2,4-Dinitrophenol	14.871	184	11387	18.387	ng	94
55) Dibenzofuran	15.171	168	198574	20.809	ng	97
56) 4-Nitrophenol	14.954	139	22343	18.849	ng	# 86
57) 2,4-Dinitrotoluene	15.112	165	47500	20.931	ng	98
58) Fluorene	15.817	166	158837	21.222	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.388	232	51855	20.601	ng	97
60) Diethylphthalate	15.571	149	157815	20.942	ng	99
61) 4-Chlorophenyl-phenyle...	15.806	204	93601	20.839	ng	97
62) 4-Nitroaniline	15.811	138	31639	20.442	ng	90
63) Azobenzene	16.093	77	130344	20.795	ng	97
65) 4,6-Dinitro-2-methylph...	15.876	198	25607	17.855	ng	96
66) n-Nitrosodiphenylamine	16.011	169	140419	20.753	ng	100
67) 4-Bromophenyl-phenylether	16.693	248	69609	20.520	ng	98
68) Hexachlorobenzene	16.816	284	84192	20.485	ng	98
69) Atrazine	16.945	200	66348	20.339	ng	97
70) Pentachlorophenol	17.157	266	35678	19.686	ng	99
71) Phenanthrene	17.545	178	290163	20.531	ng	100
72) Anthracene	17.639	178	297476	20.642	ng	100
73) Carbazole	17.897	167	253683	20.477	ng	99
74) Di-n-butylphthalate	18.450	149	285465	20.654	ng	98
75) Fluoranthene	19.542	202	395447	20.462	ng	98
77) Benzidine	19.707	184	228040	21.077	ng	99
78) Pyrene	19.901	202	413710	20.433	ng	97
80) Butylbenzylphthalate	20.770	149	145343	20.408	ng	97
81) Benzo(a)anthracene	21.734	228	503863	20.512	ng	99
82) 3,3'-Dichlorobenzidine	21.640	252	186980	20.680	ng	99
83) Chrysene	21.804	228	470111	20.299	ng	99
84) Bis(2-ethylhexyl)phtha...	21.640	149	216073	20.763	ng	98
85) Di-n-octyl phthalate	22.862	149	370539	20.598	ng	99
87) Indeno(1,2,3-cd)pyrene	28.737	276	528109	20.196	ng	# 96
88) Benzo(b)fluoranthene	23.978	252	494588	20.419	ng	99
89) Benzo(k)fluoranthene	24.049	252	502781	20.661	ng	98
90) Benzo(a)pyrene	24.860	252	455638	20.279	ng	# 97
91) Dibenzo(a,h)anthracene	28.808	278	429357	20.043	ng	93
92) Benzo(g,h,i)perylene	29.901	276	415387	20.148	ng	99

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

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