

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090425\
 Data File : BG064283.D
 Acq On : 4 Sep 2025 15:01
 Operator : RC/JU
 Sample : Q3003-01MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

VNJ-231MS

Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 09/05/2025

Supervised By :Jagrut Upadhyay 09/05/2025

Quant Time: Sep 04 15:39:06 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Aug 28 17:41:58 2025

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	30578	20.000	ng	0.00
21) Naphthalene-d8	10.640	136	135975	20.000	ng	#-0.01
39) Acenaphthene-d10	14.482	164	94592	20.000	ng	0.00
64) Phenanthrene-d10	17.220	188	216727	20.000	ng	0.00
76) Chrysene-d12	21.457	240	219333	20.000	ng	0.00
86) Perylene-d12	24.465	264	242390	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	221834	130.156	ng	0.00
7) Phenol-d6	7.026	99	302571	129.220	ng	0.00
23) Nitrobenzene-d5	9.006	82	212192	87.031	ng	-0.01
42) 2,4,6-Tribromophenol	15.969	330	175356	139.959	ng	0.00
45) 2-Fluorobiphenyl	13.107	172	536494	86.570	ng	0.00
79) Terphenyl-d14	19.841	244	947126	90.093	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.354	88	25204	35.479	ng	94
3) Pyridine	3.742	79	70496	33.712	ng	# 59
4) n-Nitrosodimethylamine	3.660	42	60147	43.485	ng	95
6) Aniline	7.185	93	73910	22.826	ng	98
8) 2-Chlorophenol	7.420	128	101088	48.537	ng	97
9) Benzaldehyde	6.991	77	39856	20.683	ng	97
10) Phenol	7.056	94	123506	47.842	ng	94
11) bis(2-Chloroethyl)ether	7.279	93	79944	41.210	ng	98
12) 1,3-Dichlorobenzene	7.743	146	103998	46.937	ng	97
13) 1,4-Dichlorobenzene	7.890	146	103777	46.409	ng	97
14) 1,2-Dichlorobenzene	8.207	146	103108	47.776	ng	96
15) Benzyl Alcohol	8.090	79	99037	45.833	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.378	45	161800	36.532	ng	97
17) 2-Methylphenol	8.295	107	83891	43.603	ng	96
18) Hexachloroethane	8.936	117	39072	44.762	ng	92
19) n-Nitroso-di-n-propyla...	8.660	70	71540	39.921	ng	93
20) 3+4-Methylphenols	8.624	107	115211	43.095	ng	94
22) Acetophenone	8.672	105	152098	44.214	ng	# 98
24) Nitrobenzene	9.048	77	114588	44.983	ng	99
25) Isophorone	9.570	82	207759	41.001	ng	98
26) 2-Nitrophenol	9.758	139	53856	48.928	ng	88
27) 2,4-Dimethylphenol	9.823	122	99925	51.708	ng	98
28) bis(2-Chloroethoxy)met...	10.052	93	112855	41.304	ng	97
29) 2,4-Dichlorophenol	10.293	162	100213	49.147	ng	97
30) 1,2,4-Trichlorobenzene	10.505	180	107351	49.932	ng	97
31) Naphthalene	10.693	128	311069	44.117	ng	97
32) Benzoic acid	9.988	122	80535m	52.500	ng	
33) 4-Chloroaniline	10.798	127	37620	13.766	ng	93
34) Hexachlorobutadiene	10.986	225	76849	48.356	ng	# 88
35) Caprolactam	11.580	113	32609m	41.302	ng	
36) 4-Chloro-3-methylphenol	11.932	107	121603	45.965	ng	96
37) 2-Methylnaphthalene	12.303	142	213233	40.685	ng	97
38) 1-Methylnaphthalene	12.520	142	224333	43.335	ng	99
40) 1,2,4,5-Tetrachloroben...	12.673	216	136168	49.892	ng	99
41) Hexachlorocyclopentadiene	12.655	237	181167	139.667	ng	96
43) 2,4,6-Trichlorophenol	12.914	196	92345	50.622	ng	91

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.984	196	99224	50.142	ng	97
46) 1,1'-Biphenyl	13.313	154	320552	46.527	ng	98
47) 2-Chloronaphthalene	13.354	162	237488	45.851	ng	96
48) 2-Nitroaniline	13.554	65	85862	48.637	ng	95
49) Acenaphthylene	14.200	152	394070	51.468	ng	97
50) Dimethylphthalate	13.942	163	306481	42.806	ng	100
51) 2,6-Dinitrotoluene	14.053	165	68426	49.366	ng	98
52) Acenaphthene	14.547	154	243642	45.091	ng	98
53) 3-Nitroaniline	14.382	138	41699	28.720	ng	97
54) 2,4-Dinitrophenol	14.594	184	96406	104.586	ng	# 83
55) Dibenzofuran	14.882	168	391719	46.090	ng	97
56) 4-Nitrophenol	14.700	139	118839	99.780	ng	90
57) 2,4-Dinitrotoluene	14.841	165	103822	49.538	ng	# 94
58) Fluorene	15.528	166	326199	46.828	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.105	232	100972	53.052	ng	97
60) Diethylphthalate	15.305	149	331930	43.318	ng	99
61) 4-Chlorophenyl-phenyle...	15.522	204	160268	46.145	ng	98
62) 4-Nitroaniline	15.546	138	70046	47.278	ng	98
63) Azobenzene	15.816	77	336575	49.282	ng	96
65) 4,6-Dinitro-2-methylph...	15.605	198	68651	55.173	ng	97
66) n-Nitrosodiphenylamine	15.740	169	282417	47.370	ng	98
67) 4-Bromophenyl-phenylether	16.415	248	111645	48.501	ng	97
68) Hexachlorobenzene	16.533	284	122542	47.938	ng	94
69) Atrazine	16.692	200	119597	48.043	ng	96
70) Pentachlorophenol	16.874	266	170377	109.924	ng	92
71) Phenanthrene	17.267	178	637909	55.806	ng	99
72) Anthracene	17.355	178	549505	47.258	ng	99
73) Carbazole	17.626	167	485803	47.427	ng	99
74) Di-n-butylphthalate	18.190	149	573286	46.584	ng	99
75) Fluoranthene	19.277	202	832099	61.909	ng	98
77) Benzidine	19.453	184	74321	16.108	ng	97
78) Pyrene	19.635	202	807627	58.446	ng	99
80) Butylbenzylphthalate	20.534	149	254673	48.864	ng	92
81) Benzo(a)anthracene	21.433	228	766323	54.613	ng	100
82) 3,3'-Dichlorobenzidine	21.351	252	73303	16.040	ng	99
83) Chrysene	21.498	228	724115	53.753	ng	99
84) Bis(2-ethylhexyl)phtha...	21.363	149	376359	48.098	ng	98
85) Di-n-octyl phthalate	22.502	149	657256	48.206	ng	97
87) Indeno(1,2,3-cd)pyrene	27.873	276	872399	52.703	ng	# 97
88) Benzo(b)fluoranthene	23.525	252	788825	57.662	ng	99
89) Benzo(k)fluoranthene	23.583	252	737256	52.899	ng	# 97
90) Benzo(a)pyrene	24.330	252	717378	56.373	ng	98
91) Dibenzo(a,h)anthracene	27.925	278	662768	49.867	ng	98
92) Benzo(g,h,i)perylene	28.930	276	703986	54.435	ng	96

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

