

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092525\
 Data File : BG064443.D
 Acq On : 25 Sep 2025 13:18
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 10/03/2025

Supervised By :Jagrut Upadhyay 10/03/2025

Quant Time: Sep 25 15:30:10 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG092525.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Sep 25 15:27:57 2025

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.841	152	15905	20.000	ng	0.00
21) Naphthalene-d8	10.626	136	69680	20.000	ng	0.00
39) Acenaphthene-d10	14.462	164	56405	20.000	ng	0.00
64) Phenanthrene-d10	17.206	188	143699	20.000	ng	# 0.00
76) Chrysene-d12	21.430	240	158110	20.000	ng	# 0.00
86) Perylene-d12	24.427	264	170967	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.426	112	16143	19.033	ng	0.00
7) Phenol-d6	7.012	99	21666	18.632	ng	0.00
23) Nitrobenzene-d5	8.986	82	25372	19.515	ng	0.00
42) 2,4,6-Tribromophenol	15.949	330	19015	19.090	ng	0.00
45) 2-Fluorobiphenyl	13.087	172	73015	19.299	ng	0.00
79) Terphenyl-d14	19.821	244	157036	19.446	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.346	88	3008	8.494	ng	86
3) Pyridine	3.734	79	9315	9.570	ng	93
4) n-Nitrosodimethylamine	3.651	42	6760	8.723	ng	98
6) Aniline	7.165	93	15311	10.165	ng	# 85
8) 2-Chlorophenol	7.412	128	10220	9.652	ng	86
9) Benzaldehyde	6.977	77	7785	7.486	ng	90
10) Phenol	7.036	94	11539	9.422	ng	90
11) bis(2-Chloroethyl)ether	7.265	93	8253	9.948	ng	95
12) 1,3-Dichlorobenzene	7.729	146	11972	10.360	ng	99
13) 1,4-Dichlorobenzene	7.876	146	11746	9.860	ng	96
14) 1,2-Dichlorobenzene	8.187	146	11269	9.888	ng	86
15) Benzyl Alcohol	8.076	79	10061	9.356	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.352	45	14740	9.894	ng	97
17) 2-Methylphenol	8.275	107	8505	9.742	ng	88
18) Hexachloroethane	8.916	117	4654	10.025	ng	91
19) n-Nitroso-di-n-propyla...	8.640	70	8091	10.060	ng	88
20) 3+4-Methylphenols	8.604	107	11454	9.290	ng	82
22) Acetophenone	8.657	105	16490	9.728	ng	# 90
24) Nitrobenzene	9.033	77	12674	9.960	ng	97
25) Isophorone	9.556	82	22504	9.462	ng	# 96
26) 2-Nitrophenol	9.738	139	5991	9.821	ng	94
27) 2,4-Dimethylphenol	9.803	122	9142	9.583	ng	97
28) bis(2-Chloroethoxy)met...	10.038	93	12010	10.017	ng	# 91
29) 2,4-Dichlorophenol	10.279	162	10370	9.303	ng	97
30) 1,2,4-Trichlorobenzene	10.490	180	11743	9.636	ng	# 89
31) Naphthalene	10.672	128	34751	9.625	ng	99
32) Benzoic acid	9.909	122	5839m	7.598	ng	
33) 4-Chloroaniline	10.784	127	13041	9.868	ng	96
34) Hexachlorobutadiene	10.966	225	10433	10.257	ng	96
35) Caprolactam	11.530	113	4162	10.264	ng	# 74
36) 4-Chloro-3-methylphenol	11.912	107	13688	9.794	ng	# 94
37) 2-Methylnaphthalene	12.282	142	26943	9.813	ng	96
38) 1-Methylnaphthalene	12.506	142	27771	10.481	ng	99
40) 1,2,4,5-Tetrachloroben...	12.658	216	16576	9.231	ng	98
41) Hexachlorocyclopentadiene	12.641	237	7251	8.011	ng	98
43) 2,4,6-Trichlorophenol	12.893	196	10895m	9.356	ng	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.976	196	11557	8.898	ng	# 88
46) 1,1'-Biphenyl	13.299	154	38218	9.503	ng	98
47) 2-Chloronaphthalene	13.340	162	29207	9.554	ng	97
48) 2-Nitroaniline	13.540	65	9517	8.929	ng	# 79
49) Acenaphthylene	14.186	152	41938	9.244	ng	# 97
50) Dimethylphthalate	13.916	163	44365	10.132	ng	99
51) 2,6-Dinitrotoluene	14.039	165	8462	9.421	ng	89
52) Acenaphthene	14.527	154	29647	9.497	ng	93
53) 3-Nitroaniline	14.368	138	8216	9.206	ng	# 75
54) 2,4-Dinitrophenol	14.574	184	3414	5.825	ng	85
55) Dibenzofuran	14.862	168	50579	9.854	ng	98
56) 4-Nitrophenol	14.685	139	5352	7.255	ng	# 79
57) 2,4-Dinitrotoluene	14.826	165	12656	8.780	ng	# 90
58) Fluorene	15.508	166	42191	9.634	ng	89
59) 2,3,4,6-Tetrachlorophenol	15.097	232	12813	9.528	ng	# 98
60) Diethylphthalate	15.285	149	48258	9.980	ng	98
61) 4-Chlorophenyl-phenyle...	15.508	204	23011	10.003	ng	95
62) 4-Nitroaniline	15.526	138	8441	8.732	ng	98
63) Azobenzene	15.796	77	35685	9.336	ng	94
65) 4,6-Dinitro-2-methylph...	15.584	198	8001	7.990	ng	81
66) n-Nitrosodiphenylamine	15.720	169	36342	9.678	ng	96
67) 4-Bromophenyl-phenylether	16.401	248	15668	9.655	ng	95
68) Hexachlorobenzene	16.513	284	18397	9.912	ng	99
69) Atrazine	16.665	200	17134	9.538	ng	82
70) Pentachlorophenol	16.865	266	8494	7.394	ng	# 83
71) Phenanthrene	17.247	178	74868	9.927	ng	98
72) Anthracene	17.335	178	72371	9.399	ng	97
73) Carbazole	17.606	167	66699	9.683	ng	97
74) Di-n-butylphthalate	18.170	149	77066	9.147	ng	98
75) Fluoranthene	19.257	202	92336	9.585	ng	97
77) Benzidine	19.439	184	33941	8.331	ng	97
78) Pyrene	19.621	202	95803	9.638	ng	98
80) Butylbenzylphthalate	20.514	149	35243	9.452	ng	96
81) Benzo(a)anthracene	21.413	228	97105	9.585	ng	95
82) 3,3'-Dichlorobenzidine	21.336	252	34415	9.950	ng	98
83) Chrysene	21.477	228	92795	9.684	ng	99
84) Bis(2-ethylhexyl)phtha...	21.336	149	49375	9.158	ng	98
85) Di-n-octyl phthalate	22.470	149	85942	9.459	ng	99
87) Indeno(1,2,3-cd)pyrene	27.794	276	113528	9.435	ng	# 98
88) Benzo(b)fluoranthene	23.481	252	93375m	9.454	ng	
89) Benzo(k)fluoranthene	23.546	252	95222m	9.677	ng	
90) Benzo(a)pyrene	24.280	252	89595	9.818	ng	95
91) Dibenzo(a,h)anthracene	27.858	278	91287	9.400	ng	# 95
92) Benzo(g,h,i)perylene	28.845	276	90607	9.736	ng	# 96

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

