

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082825\
 Data File : BG064219.D
 Acq On : 28 Aug 2025 11:32
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC005

Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 08/29/2025

Supervised By :Jagrut Upadhyay 08/29/2025

Quant Time: Aug 28 18:40:09 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	7.859	152	24098	20.000	ng	0.00
21) Naphthalene-d8	10.644	136	109950	20.000	ng	0.00
39) Acenaphthene-d10	14.481	164	78593	20.000	ng	0.00
64) Phenanthrene-d10	17.219	188	199845	20.000	ng	0.00
76) Chrysene-d12	21.449	240	219683	20.000	ng	0.00
86) Perylene-d12	24.458	264	241429	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.439	112	13573	10.105	ng	0.00
7) Phenol-d6	7.019	99	18310	9.922	ng	-0.01
23) Nitrobenzene-d5	9.005	82	18764	9.518	ng	-0.01
42) 2,4,6-Tribromophenol	15.962	330	11040	10.605	ng	-0.01
45) 2-Fluorobiphenyl	13.106	172	53596	10.409	ng	0.00
79) Terphenyl-d14	19.839	244	110867	10.529	ng	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.365	88	3459	6.178	ng	88
3) Pyridine	3.752	79	8730	5.297	ng	# 58
6) Aniline	7.184	93	12228	4.792	ng	89
8) 2-Chlorophenol	7.419	128	7939	4.837	ng	96
10) Phenol	7.043	94	9580	4.709	ng	97
11) bis(2-Chloroethyl)ether	7.278	93	7421	4.854	ng	88
12) 1,3-Dichlorobenzene	7.742	146	8854m	5.071	ng	
13) 1,4-Dichlorobenzene	7.895	146	9030	5.124	ng	93
14) 1,2-Dichlorobenzene	8.206	146	8131	4.781	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.376	45	17609	5.045	ng	95
17) 2-Methylphenol	8.294	107	7155	4.719	ng	# 91
18) Hexachloroethane	8.935	117	3344	4.861	ng	86
19) n-Nitroso-di-n-propyla...	8.653	70	7001	4.957	ng	# 98
22) Acetophenone	8.670	105	14639	5.263	ng	# 99
24) Nitrobenzene	9.046	77	9728	4.723	ng	98
25) Isophorone	9.569	82	20579	5.023	ng	# 95
26) 2-Nitrophenol	9.757	139	3668	4.121	ng	97
27) 2,4-Dimethylphenol	9.816	122	7568	4.843	ng	96
28) bis(2-Chloroethoxy)met...	10.057	93	10885	4.927	ng	# 88
29) 2,4-Dichlorophenol	10.292	162	7945	4.819	ng	91
30) 1,2,4-Trichlorobenzene	10.509	180	9136	5.255	ng	95
31) Naphthalene	10.691	128	30344	5.322	ng	# 94
33) 4-Chloroaniline	10.797	127	10141	4.589	ng	# 91
34) Hexachlorobutadiene	10.985	225	6663	5.185	ng	95
36) 4-Chloro-3-methylphenol	11.925	107	10455	4.887	ng	95
37) 2-Methylnaphthalene	12.301	142	21107	4.980	ng	87
38) 1-Methylnaphthalene	12.525	142	21661m	5.175	ng	
40) 1,2,4,5-Tetrachloroben...	12.671	216	10879	4.797	ng	97
43) 2,4,6-Trichlorophenol	12.912	196	7156	4.721	ng	87
44) 2,4,5-Trichlorophenol	12.977	196	7600	4.622	ng	84
46) 1,1'-Biphenyl	13.312	154	28672	5.009	ng	94
47) 2-Chloronaphthalene	13.359	162	21411	4.975	ng	100
48) 2-Nitroaniline	13.553	65	5960	4.063	ng	89
49) Acenaphthylene	14.199	152	32547	5.116	ng	99
50) Dimethylphthalate	13.935	163	32855	5.523	ng	98
51) 2,6-Dinitrotoluene	14.052	165	4684	4.067	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Acenaphthene	14.546	154	22964	5.115	ng	94
53) 3-Nitroaniline	14.375	138	5404	4.480	ng #	98
55) Dibenzofuran	14.881	168	37579	5.322	ng	95
57) 2,4-Dinitrotoluene	14.834	165	7779	4.467	ng #	94
58) Fluorene	15.527	166	30779	5.318	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.104	232	7695	4.866	ng #	96
60) Diethylphthalate	15.298	149	35947	5.646	ng	97
61) 4-Chlorophenyl-phenyle...	15.521	204	15345	5.318	ng	93
62) 4-Nitroaniline	15.533	138	5254	4.268	ng #	75
63) Azobenzene	15.815	77	29276	5.159	ng	96
66) n-Nitrosodiphenylamine	15.733	169	27609	5.022	ng	98
67) 4-Bromophenyl-phenylether	16.414	248	9745	4.591	ng	98
68) Hexachlorobenzene	16.526	284	12043	5.109	ng	92
69) Atrazine	16.678	200	12001	5.228	ng	97
71) Phenanthrene	17.260	178	55100	5.227	ng	97
72) Anthracene	17.348	178	56156	5.237	ng	99
73) Carbazole	17.619	167	49241	5.213	ng	98
74) Di-n-butylphthalate	18.188	149	59765	5.267	ng	98
75) Fluoranthene	19.270	202	66964	5.403	ng	99
78) Pyrene	19.634	202	70961	5.127	ng	97
80) Butylbenzylphthalate	20.533	149	24638	4.720	ng	98
81) Benzo(a)anthracene	21.432	228	72646	5.169	ng	96
83) Chrysene	21.490	228	70143	5.199	ng	98
84) Bis(2-ethylhexyl)phtha...	21.361	149	39184	5.000	ng	94
87) Indeno(1,2,3-cd)pyrene	27.836	276	79337	4.812	ng	92
88) Benzo(b)fluoranthene	23.506	252	65111	4.778	ng	99
89) Benzo(k)fluoranthene	23.570	252	70084	5.049	ng	98
90) Benzo(a)pyrene	24.311	252	61491	4.851	ng #	95
91) Dibenzo(a,h)anthracene	27.895	278	63017	4.760	ng #	96
92) Benzo(g,h,i)perylene	28.888	276	64396	4.999	ng #	96

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

