

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090325\
 Data File : BG064250.D
 Acq On : 3 Sep 2025 12:04
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 09/04/2025

Supervised By :Jagrut Upadhyay 09/04/2025

Quant Time: Sep 03 12:58:14 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	7.853	152	38309	20.000	ng	0.00
21) Naphthalene-d8	10.644	136	190272	20.000	ng	0.00
39) Acenaphthene-d10	14.481	164	144796	20.000	ng	0.00
64) Phenanthrene-d10	17.225	188	325433	20.000	ng	0.00
76) Chrysene-d12	21.455	240	228359	20.000	ng	0.00
86) Perylene-d12	24.469	264	252610	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.438	112	175991	82.420	ng	0.00
7) Phenol-d6	7.025	99	269625	91.912	ng	0.00
23) Nitrobenzene-d5	9.011	82	288769	84.641	ng	0.00
42) 2,4,6-Tribromophenol	15.973	330	160034	83.443	ng	0.00
45) 2-Fluorobiphenyl	13.106	172	737502	77.743	ng	0.00
79) Terphenyl-d14	19.845	244	929215	84.896	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.358	88	35305	39.668	ng	89
3) Pyridine	3.746	79	109062	41.629	ng	# 58
4) n-Nitrosodimethylamine	3.664	42	66615	38.442	ng	# 87
6) Aniline	7.183	93	176418	43.488	ng	97
8) 2-Chlorophenol	7.424	128	115727	44.352	ng	99
9) Benzaldehyde	6.995	77	106319m	44.040	ng	
10) Phenol	7.054	94	151604	46.875	ng	98
11) bis(2-Chloroethyl)ether	7.277	93	107157	44.091	ng	97
12) 1,3-Dichlorobenzene	7.747	146	113967	41.056	ng	96
13) 1,4-Dichlorobenzene	7.888	146	116247	41.494	ng	97
14) 1,2-Dichlorobenzene	8.206	146	117287	43.378	ng	97
15) Benzyl Alcohol	8.088	79	119278	44.061	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.376	45	227806	41.055	ng	98
17) 2-Methylphenol	8.294	107	106486	44.177	ng	93
18) Hexachloroethane	8.934	117	45935	42.005	ng	90
19) n-Nitroso-di-n-propyla...	8.658	70	102191	45.517	ng	97
20) 3+4-Methylphenols	8.629	107	147569	44.060	ng	92
22) Acetophenone	8.676	105	199646	41.475	ng	# 95
24) Nitrobenzene	9.052	77	146907	41.213	ng	100
25) Isophorone	9.575	82	304138	42.893	ng	97
26) 2-Nitrophenol	9.757	139	68162	44.254	ng	93
27) 2,4-Dimethylphenol	9.821	122	118013	43.641	ng	98
28) bis(2-Chloroethoxy)met...	10.056	93	164805	43.105	ng	# 94
29) 2,4-Dichlorophenol	10.297	162	121120	42.450	ng	94
30) 1,2,4-Trichlorobenzene	10.509	180	120973	40.211	ng	98
31) Naphthalene	10.697	128	399340	40.474	ng	98
32) Benzoic acid	9.986	122	92463m	43.075	ng	
33) 4-Chloroaniline	10.803	127	167102	43.697	ng	99
34) Hexachlorobutadiene	10.985	225	86498	38.896	ng	90
35) Caprolactam	11.584	113	44550	40.324	ng	# 92
36) 4-Chloro-3-methylphenol	11.931	107	164546	44.448	ng	98
37) 2-Methylnaphthalene	12.307	142	309518	42.204	ng	99
38) 1-Methylnaphthalene	12.524	142	302302	41.732	ng	100
40) 1,2,4,5-Tetrachloroben...	12.677	216	165591	39.636	ng	99
41) Hexachlorocyclopentadiene	12.659	237	81570	41.081	ng	94
43) 2,4,6-Trichlorophenol	12.918	196	114233	40.909	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.988	196	126630	41.804	ng	96
46) 1,1'-Biphenyl	13.317	154	424781	40.278	ng	98
47) 2-Chloronaphthalene	13.358	162	316208	39.882	ng	98
48) 2-Nitroaniline	13.558	65	119620	44.265	ng	90
49) Acenaphthylene	14.205	152	478582	40.834	ng	98
50) Dimethylphthalate	13.940	163	433350	39.540	ng	99
51) 2,6-Dinitrotoluene	14.058	165	90664	42.731	ng	92
52) Acenaphthene	14.545	154	330706	39.984	ng	96
53) 3-Nitroaniline	14.387	138	92706	41.712	ng	83
54) 2,4-Dinitrophenol	14.592	184	49696	38.968	ng	93
55) Dibenzofuran	14.880	168	519494	39.931	ng	99
56) 4-Nitrophenol	14.698	139	67715	37.142	ng	91
57) 2,4-Dinitrotoluene	14.845	165	138285	43.105	ng	# 92
58) Fluorene	15.532	166	426768	40.023	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.109	232	116971	40.149	ng	98
60) Diethylphthalate	15.309	149	461968	39.385	ng	99
61) 4-Chlorophenyl-phenyle...	15.527	204	213406	40.141	ng	100
62) 4-Nitroaniline	15.550	138	89072	39.275	ng	91
63) Azobenzene	15.820	77	425374	40.689	ng	98
65) 4,6-Dinitro-2-methylph...	15.609	198	81520	43.631	ng	99
66) n-Nitrosodiphenylamine	15.744	169	372610	41.621	ng	98
67) 4-Bromophenyl-phenylether	16.420	248	145919	42.216	ng	98
68) Hexachlorobenzene	16.537	284	162012	42.208	ng	94
69) Atrazine	16.690	200	123522	33.045	ng	98
70) Pentachlorophenol	16.878	266	95296	40.946	ng	93
71) Phenanthrene	17.266	178	685447	39.934	ng	98
72) Anthracene	17.360	178	701562	40.181	ng	99
73) Carbazole	17.624	167	551475	35.854	ng	98
74) Di-n-butylphthalate	18.194	149	594738	32.184	ng	100
75) Fluoranthene	19.275	202	645991	32.008	ng	99
77) Benzidine	19.457	184	202472	42.148	ng	100
78) Pyrene	19.639	202	649982	45.179	ng	99
80) Butylbenzylphthalate	20.532	149	247977	45.699	ng	95
81) Benzo(a)anthracene	21.437	228	599916	41.064	ng	98
82) 3,3'-Dichlorobenzidine	21.355	252	202059	42.467	ng	97
83) Chrysene	21.496	228	565854	40.344	ng	98
84) Bis(2-ethylhexyl)phtha...	21.361	149	361781	44.407	ng	99
85) Di-n-octyl phthalate	22.501	149	629646	44.356	ng	99
87) Indeno(1,2,3-cd)pyrene	27.865	276	712898	41.325	ng	# 92
88) Benzo(b)fluoranthene	23.517	252	584279	40.982	ng	99
89) Benzo(k)fluoranthene	23.582	252	602620	41.489	ng	100
90) Benzo(a)pyrene	24.328	252	546166	41.182	ng	99
91) Dibenzo(a,h)anthracene	27.924	278	577561	41.698	ng	99
92) Benzo(g,h,i)perylene	28.923	276	555765	41.235	ng	98

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

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