

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082725\
 Data File : BG064208.D
 Acq On : 27 Aug 2025 12:54
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC005

Quant Time: Aug 27 17:46:11 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082725.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 27 17:42:35 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	25638	20.000	ng	0.00
21) Naphthalene-d8	10.643	136	116707	20.000	ng	# 0.00
39) Acenaphthene-d10	14.479	164	87106	20.000	ng	0.00
64) Phenanthrene-d10	17.217	188	208346	20.000	ng	0.00
76) Chrysene-d12	21.448	240	231776	20.000	ng	0.00
86) Perylene-d12	24.456	264	248075	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.443	112	12517	9.353	ng	0.00
7) Phenol-d6	7.018	99	18539	9.368	ng	0.00
23) Nitrobenzene-d5	9.004	82	19996	9.673	ng	0.00
42) 2,4,6-Tribromophenol	15.966	330	10789	9.504	ng	0.00
45) 2-Fluorobiphenyl	13.105	172	56841	10.323	ng	0.00
79) Terphenyl-d14	19.838	244	114453	10.434	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.363	88	3770	6.114	ng	# 81
3) Pyridine	3.757	79	8841	5.090	ng	88
6) Aniline	7.182	93	12983	4.834	ng	95
8) 2-Chlorophenol	7.423	128	8432	4.906	ng	90
10) Phenol	7.041	94	10086	4.624	ng	97
11) bis(2-Chloroethyl)ether	7.282	93	7651	4.854	ng	93
12) 1,3-Dichlorobenzene	7.746	146	9012	4.969	ng	99
13) 1,4-Dichlorobenzene	7.887	146	9501	5.086	ng	90
14) 1,2-Dichlorobenzene	8.210	146	9184	5.086	ng	91
16) 2,2'-oxybis(1-Chloropr...	8.375	45	18736	5.159	ng	98
17) 2-Methylphenol	8.299	107	7648	4.864	ng	# 81
18) Hexachloroethane	8.939	117	3369	4.703	ng	# 81
19) n-Nitroso-di-n-propyla...	8.657	70	7039	4.701	ng	# 84
22) Acetophenone	8.669	105	15391	5.270	ng	96
24) Nitrobenzene	9.045	77	10660	4.962	ng	97
25) Isophorone	9.573	82	21977	5.051	ng	95
26) 2-Nitrophenol	9.756	139	4286	4.627	ng	92
27) 2,4-Dimethylphenol	9.814	122	8464	5.051	ng	# 78
28) bis(2-Chloroethoxy)met...	10.055	93	12096	5.163	ng	# 92
29) 2,4-Dichlorophenol	10.296	162	8093	4.686	ng	94
30) 1,2,4-Trichlorobenzene	10.502	180	9565	5.201	ng	90
31) Naphthalene	10.696	128	30719	5.155	ng	98
33) 4-Chloroaniline	10.801	127	10376	4.457	ng	97
34) Hexachlorobutadiene	10.989	225	6907	5.130	ng	98
36) 4-Chloro-3-methylphenol	11.918	107	10833	4.817	ng	# 85
37) 2-Methylnaphthalene	12.306	142	21649	4.912	ng	93
38) 1-Methylnaphthalene	12.523	142	21784	4.982	ng	97
40) 1,2,4,5-Tetrachloroben...	12.676	216	12044	5.006	ng	# 97
43) 2,4,6-Trichlorophenol	12.911	196	7980	4.831	ng	94
44) 2,4,5-Trichlorophenol	12.981	196	9241	5.050	ng	96
46) 1,1'-Biphenyl	13.316	154	31479	5.097	ng	93
47) 2-Chloronaphthalene	13.351	162	22338	4.846	ng	96
48) 2-Nitroaniline	13.551	65	6741	4.219	ng	# 81
49) Acenaphthylene	14.203	152	32994	4.822	ng	96
50) Dimethylphthalate	13.933	163	34731	5.314	ng	99
51) 2,6-Dinitrotoluene	14.051	165	4980	4.036	ng	# 76

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Acenaphthene	14.544	154	23748	4.943	ng	97
53) 3-Nitroaniline	14.380	138	4798	3.770	ng	98
55) Dibenzofuran	14.879	168	39738	5.213	ng	97
57) 2,4-Dinitrotoluene	14.838	165	7963	4.206	ng #	98
58) Fluorene	15.525	166	31891	5.120	ng #	94
59) 2,3,4,6-Tetrachlorophenol	15.108	232	8362	4.830	ng #	98
60) Diethylphthalate	15.302	149	36349	5.181	ng	96
61) 4-Chlorophenyl-phenyle...	15.525	204	17100	5.320	ng	96
62) 4-Nitroaniline	15.537	138	5280	3.837	ng	85
63) Azobenzene	15.813	77	31827	5.163	ng	94
66) n-Nitrosodiphenylamine	15.731	169	29502	5.233	ng	92
67) 4-Bromophenyl-phenylether	16.418	248	10815	4.939	ng	93
68) Hexachlorobenzene	16.530	284	11738	4.988	ng	98
69) Atrazine	16.683	200	12473	5.229	ng	96
71) Phenanthrene	17.264	178	56100	5.213	ng	96
72) Anthracene	17.353	178	56227	5.159	ng	99
73) Carbazole	17.617	167	50106	5.148	ng	95
74) Di-n-butylphthalate	18.193	149	60247	5.089	ng	98
75) Fluoranthene	19.274	202	67529	5.250	ng	100
78) Pyrene	19.632	202	73745	5.105	ng	97
80) Butylbenzylphthalate	20.531	149	25364	4.681	ng	95
81) Benzo(a)anthracene	21.430	228	74261	5.014	ng	96
83) Chrysene	21.495	228	71506	5.020	ng	98
84) Bis(2-ethylhexyl)phtha...	21.366	149	39181	4.812	ng #	89
87) Indeno(1,2,3-cd)pyrene	27.840	276	82625	4.919	ng #	90
88) Benzo(b)fluoranthene	23.510	252	70389	5.051	ng	95
89) Benzo(k)fluoranthene	23.569	252	69359	4.845	ng	96
90) Benzo(a)pyrene	24.315	252	64554	4.971	ng #	97
91) Dibenzo(a,h)anthracene	27.893	278	65274	4.817	ng #	96
92) Benzo(g,h,i)perylene	28.892	276	65519	5.002	ng #	92

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

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