

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012425\
 Data File : BG063954.D
 Acq On : 24 Jan 2025 13:09
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC005

Quant Time: Jan 27 14:57:01 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 14:52:53 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.950	152	71685	20.000	ng	0.00
21) Naphthalene-d8	10.747	136	303409	20.000	ng	0.00
39) Acenaphthene-d10	14.571	164	201031	20.000	ng	0.00
64) Phenanthrene-d10	17.315	188	460358	20.000	ng	0.00
76) Chrysene-d12	21.557	240	499405	20.000	ng	0.00
86) Perylene-d12	24.660	264	513239	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.535	112	54641	10.517	ng	0.00
7) Phenol-d6	7.104	99	73674	10.499	ng	-0.01
23) Nitrobenzene-d5	0.000	82	0d	0.000	ng	
42) 2,4,6-Tribromophenol	16.058	330	16370	7.972	ng	0.00
45) 2-Fluorobiphenyl	13.197	172	145493	10.228	ng	0.00
79) Terphenyl-d14	19.930	244	276404	11.061	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.467	88	11006	4.778	ng	91
3) Pyridine	3.855	79	28595	4.530	ng	89
4) n-Nitrosodimethylamine	3.772	42	16426	4.574	ng	# 86
6) Aniline	7.274	93	29477	5.069	ng	99
8) 2-Chlorophenol	7.509	128	29934	5.488	ng	93
9) Benzaldehyde	7.092	77	24109	5.828	ng	95
10) Phenol	7.133	94	39209	5.290	ng	96
11) bis(2-Chloroethyl)ether	7.374	93	32810	5.657	ng	93
12) 1,3-Dichlorobenzene	7.844	146	30135	5.118	ng	96
13) 1,4-Dichlorobenzene	7.985	146	30834	5.166	ng	93
14) 1,2-Dichlorobenzene	8.302	146	29906	5.157	ng	94
15) Benzyl Alcohol	8.179	79	25265	4.852	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.479	45	52596	4.906	ng	98
17) 2-Methylphenol	8.385	107	22956	4.757	ng	95
18) Hexachloroethane	9.031	117	9779	4.698	ng	# 86
19) n-Nitroso-di-n-propyla...	8.749	70	23374	4.942	ng	# 92
20) 3+4-Methylphenols	8.702	107	30509	4.637	ng	94
22) Acetophenone	8.767	105	45487	5.204	ng	# 97
24) Nitrobenzene	9.143	77	22240	3.868	ng	94
25) Isophorone	9.665	82	62474	5.062	ng	99
26) 2-Nitrophenol	9.859	139	5287	6.286	ng	# 75
27) 2,4-Dimethylphenol	9.906	122	17178	4.803	ng	86
28) bis(2-Chloroethoxy)met...	10.153	93	37317	4.943	ng	# 99
29) 2,4-Dichlorophenol	10.382	162	20606	4.570	ng	91
30) 1,2,4-Trichlorobenzene	10.606	180	29115	5.246	ng	98
31) Naphthalene	10.794	128	89547	5.140	ng	98
33) 4-Chloroaniline	10.893	127	30910	5.009	ng	# 91
34) Hexachlorobutadiene	11.087	225	17846	4.938	ng	99
35) Caprolactam	11.628	113	8263	4.726	ng	91
36) 4-Chloro-3-methylphenol	12.010	107	28161	4.862	ng	91
37) 2-Methylnaphthalene	12.403	142	58471	4.991	ng	97
38) 1-Methylnaphthalene	12.621	142	59574	5.080	ng	92
40) 1,2,4,5-Tetrachloroben...	12.768	216	30557	4.885	ng	99
41) Hexachlorocyclopentadiene	12.750	237	4665	3.635	ng	89
43) 2,4,6-Trichlorophenol	12.997	196	16220	4.259	ng	95
44) 2,4,5-Trichlorophenol	13.067	196	19581	4.460	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.408	154	82563	5.069	ng	96
47) 2-Chloronaphthalene	13.449	162	62508	4.994	ng	97
48) 2-Nitroaniline	13.637	65	8893	7.101	ng	# 86
49) Acenaphthylene	14.295	152	89363	4.952	ng	99
50) Dimethylphthalate	14.025	163	77912	5.121	ng	99
51) 2,6-Dinitrotoluene	14.143	165	5794	6.570	ng	# 87
52) Acenaphthene	14.636	154	61916	4.982	ng	98
53) 3-Nitroaniline	14.466	138	7413	5.495	ng	# 81
55) Dibenzofuran	14.971	168	97212	5.117	ng	99
57) 2,4-Dinitrotoluene	14.918	165	7571	7.436	ng	# 92
58) Fluorene	15.617	166	80501	5.218	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.194	232	14340	3.888	ng	94
60) Diethylphthalate	15.394	149	80950	4.977	ng	97
61) 4-Chlorophenyl-phenyle...	15.617	204	40334	5.326	ng	99
62) 4-Nitroaniline	15.623	138	8975	6.012	ng	95
63) Azobenzene	15.905	77	88069	5.035	ng	97
66) n-Nitrosodiphenylamine	15.829	169	66935	4.905	ng	97
67) 4-Bromophenyl-phenylether	16.510	248	23391	4.896	ng	95
68) Hexachlorobenzene	16.622	284	28616	5.060	ng	99
69) Atrazine	16.775	200	10844	5.499	ng	93
71) Phenanthrene	17.356	178	130792	5.114	ng	100
72) Anthracene	17.445	178	124848	4.976	ng	98
73) Carbazole	17.709	167	123527	4.961	ng	98
74) Di-n-butylphthalate	18.291	149	141253	4.821	ng	99
75) Fluoranthene	19.366	202	162167	5.067	ng	98
77) Benzidine	19.542	184	42716	4.897	ng	96
78) Pyrene	19.724	202	175059	5.239	ng	98
80) Butylbenzylphthalate	20.629	149	51855	4.206	ng	98
81) Benzo(a)anthracene	21.540	228	170553	4.980	ng	98
82) 3,3'-Dichlorobenzidine	21.463	252	51330	4.644	ng	# 96
83) Chrysene	21.604	228	165408	5.000	ng	96
84) Bis(2-ethylhexyl)phtha...	21.487	149	89506	4.530	ng	97
85) Di-n-octyl phthalate	22.668	149	141520	4.359	ng	95
87) Indeno(1,2,3-cd)pyrene	28.150	276	171022	4.766	ng	# 89
88) Benzo(b)fluoranthene	23.678	252	158022	4.776	ng	96
89) Benzo(k)fluoranthene	23.743	252	159060	4.837	ng	97
90) Benzo(a)pyrene	24.513	252	137611	4.739	ng	99
91) Dibenzo(a,h)anthracene	28.214	278	137942	4.613	ng	98
92) Benzo(g,h,i)perylene	29.231	276	141602	4.721	ng	99

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

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