

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020325\
 Data File : BG063976.D
 Acq On : 3 Feb 2025 11:56
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC005

Quant Time: Feb 04 02:44:51 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 04 02:44:47 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.873	152	44732	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	206910	20.000	ng	0.00
39) Acenaphthene-d10	14.494	164	150510	20.000	ng	0.00
64) Phenanthrene-d10	17.238	188	363622	20.000	ng	0.00
76) Chrysene-d12	21.468	240	424138	20.000	ng	0.00
86) Perylene-d12	24.488	264	422465	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	23509	8.426	ng	0.00
7) Phenol-d6	7.032	99	34391	8.982	ng	0.00
23) Nitrobenzene-d5	0.000	82	0d	0.000	ng	
42) 2,4,6-Tribromophenol	0.000	330	0d	0.000	ng	
45) 2-Fluorobiphenyl	13.119	172	98850	10.063	ng	0.00
79) Terphenyl-d14	19.859	244	213179	10.497	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.384	88	6951	5.246	ng	98
3) Pyridine	3.772	79	17441	4.873	ng	98
4) n-Nitrosodimethylamine	3.683	42	9612	4.672	ng	# 93
6) Aniline	7.197	93	20508	4.998	ng	98
8) 2-Chlorophenol	7.438	128	11388	4.006	ng	94
9) Benzaldehyde	7.009	77	10197	4.879	ng	90
10) Phenol	7.056	94	18408	4.567	ng	95
11) bis(2-Chloroethyl)ether	7.297	93	16734	5.156	ng	93
12) 1,3-Dichlorobenzene	7.761	146	17308	5.116	ng	97
13) 1,4-Dichlorobenzene	7.908	146	17637	5.162	ng	93
14) 1,2-Dichlorobenzene	8.219	146	16538	5.050	ng	94
15) Benzyl Alcohol	8.102	79	12607	4.405	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.401	45	33346	5.103	ng	97
17) 2-Methylphenol	8.307	107	11680	4.481	ng	89
18) Hexachloroethane	8.954	117	5167	4.595	ng	93
19) n-Nitroso-di-n-propyla...	8.672	70	13015	4.909	ng	# 97
20) 3+4-Methylphenols	8.631	107	14960	4.308	ng	96
22) Acetophenone	8.683	105	28008	4.927	ng	98
24) Nitrobenzene	9.065	77	10249	6.222	ng	94
25) Isophorone	9.588	82	35059	4.750	ng	# 98
26) 2-Nitrophenol	9.770	139	2167	3.777	ng	# 83
27) 2,4-Dimethylphenol	9.835	122	8453	3.982	ng	88
28) bis(2-Chloroethoxy)met...	10.076	93	21986	4.682	ng	99
29) 2,4-Dichlorophenol	10.311	162	7983	6.350	ng	99
30) 1,2,4-Trichlorobenzene	10.522	180	17000	4.874	ng	95
31) Naphthalene	10.710	128	55706	5.022	ng	97
33) 4-Chloroaniline	10.816	127	20467	4.815	ng	97
34) Hexachlorobutadiene	11.010	225	10503	4.729	ng	96
36) 4-Chloro-3-methylphenol	11.933	107	13212	3.888	ng	# 94
37) 2-Methylnaphthalene	12.320	142	39386	5.095	ng	98
38) 1-Methylnaphthalene	12.538	142	40640	5.313	ng	98
40) 1,2,4,5-Tetrachloroben...	12.685	216	21324	4.804	ng	98
41) Hexachlorocyclopentadiene	12.679	237	3394	2.834	ng	83
43) 2,4,6-Trichlorophenol	12.925	196	5463	6.777	ng	87
44) 2,4,5-Trichlorophenol	12.996	196	6632	6.014	ng	88
46) 1,1'-Biphenyl	13.331	154	57237	4.995	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.366	162	38700	4.633	ng	99
48) 2-Nitroaniline	13.560	65	4447	6.486	ng	93
49) Acenaphthylene	14.212	152	61714	4.757	ng	99
50) Dimethylphthalate	13.954	163	38723	3.880	ng	97
51) 2,6-Dinitrotoluene	14.059	165	3474	6.308	ng	91
52) Acenaphthene	14.559	154	41919	4.817	ng	95
53) 3-Nitroaniline	14.388	138	4253	7.043	ng	# 87
55) Dibenzofuran	14.894	168	70135	4.963	ng	99
57) 2,4-Dinitrotoluene	14.847	165	4695	6.839	ng	# 87
58) Fluorene	15.546	166	56205	5.146	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.111	232	5485	6.936	ng	# 77
60) Diethylphthalate	15.323	149	39972	3.778	ng	98
61) 4-Chlorophenyl-phenyle...	15.540	204	29007	5.293	ng	98
62) 4-Nitroaniline	15.546	138	6135	6.320	ng	88
63) Azobenzene	15.834	77	63559	5.086	ng	95
66) n-Nitrosodiphenylamine	15.752	169	49093	4.869	ng	96
67) 4-Bromophenyl-phenylether	16.433	248	16670	4.606	ng	97
68) Hexachlorobenzene	16.545	284	20705	4.866	ng	97
69) Atrazine	16.703	200	12851	3.830	ng	86
71) Phenanthrene	17.273	178	95874	5.020	ng	97
72) Anthracene	17.367	178	95726	5.009	ng	98
73) Carbazole	17.632	167	89215	4.987	ng	97
74) Di-n-butylphthalate	18.213	149	60998	3.271	ng	98
75) Fluoranthene	19.289	202	119253	5.128	ng	100
77) Benzidine	19.471	184	29139	3.226	ng	# 96
78) Pyrene	19.647	202	129379	4.941	ng	96
80) Butylbenzylphthalate	20.558	149	15902	7.094	ng	# 86
81) Benzo(a)anthracene	21.445	228	132362	4.915	ng	98
82) 3,3'-Dichlorobenzidine	21.369	252	28005	3.339	ng	97
83) Chrysene	21.510	228	134964	5.036	ng	99
84) Bis(2-ethylhexyl)phtha...	21.398	149	28808	6.250	ng	95
87) Indeno(1,2,3-cd)pyrene	27.890	276	121101	4.443	ng	# 73
88) Benzo(b)fluoranthene	23.537	252	115485	4.642	ng	96
89) Benzo(k)fluoranthene	23.595	252	120118	4.702	ng	96
90) Benzo(a)pyrene	24.342	252	101336	4.564	ng	99
91) Dibenzo(a,h)anthracene	27.949	278	103956	4.517	ng	98
92) Benzo(g,h,i)perylene	28.936	276	103585	4.459	ng	98

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

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