

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100820\
 Data File : BG046818.D
 Acq On : 8 Oct 2020 16:48
 Operator : CG/JU
 Sample : L4293-05MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B-6-MW-2MSD

Quant Time: Oct 08 17:40:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 07 11:38:17 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.109	152	65045	20.00	ng	# 0.00
21) Naphthalene-d8	10.917	136	252721	20.00	ng	0.00
39) Acenaphthene-d10	14.731	164	165739	20.00	ng	0.00
64) Phenanthrene-d10	17.474	188	404016	20.00	ng	0.00
76) Chrysene-d12	21.752	240	469216	20.00	ng	# 0.00
86) Perylene-d12	25.019	264	499001	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.671	112	222588	54.82	ng	0.00
7) Phenol-d6	7.257	99	181161	31.63	ng	0.00
23) Nitrobenzene-d5	9.272	82	503070	105.90	ng	0.00
42) 2,4,6-Tribromophenol	16.217	330	350546	160.40	ng	0.00
45) 2-Fluorobiphenyl	13.356	172	1085559	92.73	ng	0.00
79) Terphenyl-d14	20.077	244	1928092	82.36	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.567	88	29201	15.41	ng	94
3) Pyridine	3.961	79	82981	15.48	ng	96
4) n-Nitrosodimethylamine	3.873	42	46612	16.62	ng	# 66
6) Aniline	7.427	93	159128	22.99	ng	93
8) 2-Chlorophenol	7.668	128	158347	39.14	ng	93
9) Benzaldehyde	7.239	77	56309	13.73	ng	85
10) Phenol	7.281	94	79541	13.95	ng	79
11) bis(2-Chloroethyl)ether	7.521	93	190572	43.25	ng	90
12) 1,3-Dichlorobenzene	7.997	146	206633	39.55	ng	# 93
13) 1,4-Dichlorobenzene	8.144	146	210260	40.61	ng	98
14) 1,2-Dichlorobenzene	8.461	146	206039	42.67	ng	95
15) Benzyl Alcohol	8.338	79	131491	27.93	ng	# 88
16) 2,2'-oxybis(1-Chloropr...	8.632	45	277804	35.93	ng	98
17) 2-Methylphenol	8.544	107	110850	30.28	ng	# 89
18) Hexachloroethane	9.202	117	74642	38.22	ng	92
19) n-Nitroso-di-n-propyla...	8.908	70	170614	43.44	ng	95
20) 3+4-Methylphenols	8.867	107	134720	27.00	ng	91
22) Acetophenone	8.926	105	315893	43.03	ng	# 93
24) Nitrobenzene	9.313	77	253158	48.82	ng	# 86
25) Isophorone	9.836	82	449985	43.40	ng	94
26) 2-Nitrophenol	10.024	139	111325	59.39	ng	# 79
27) 2,4-Dimethylphenol	10.077	122	165007	44.04	ng	90
28) bis(2-Chloroethoxy)met...	10.318	93	262978	41.11	ng	98
29) 2,4-Dichlorophenol	10.559	162	199471	44.25	ng	95
30) 1,2,4-Trichlorobenzene	10.782	180	218810	40.67	ng	96
31) Naphthalene	10.970	128	580159	42.32	ng	100
32) Benzoic acid	10.160	122	30890	16.60	ng	94
33) 4-Chloroaniline	11.070	127	139256	22.93	ng	# 94
34) Hexachlorobutadiene	11.252	225	154267	39.11	ng	97
35) Caprolactam	11.840	113	11032	7.23	ng	98
36) 4-Chloro-3-methylphenol	12.175	107	195178	40.28	ng	89
37) 2-Methylnaphthalene	12.568	142	445583	43.85	ng	# 95
38) 1-Methylnaphthalene	12.786	142	417413	44.25	ng	100
40) 1,2,4,5-Tetrachloroben...	12.933	216	281943	47.32	ng	# 96
41) Hexachlorocyclopentadiene	12.915	237	339335	97.92	ng	99
43) 2,4,6-Trichlorophenol	13.162	196	187951	50.16	ng	92

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44) 2,4,5-Trichlorophenol	13.232	196	200601	53.20	ng	#	96
46) 1,1'-Biphenyl	13.567	154	602747	47.30	ng		96
47) 2-Chloronaphthalene	13.614	162	466597	45.06	ng		97
48) 2-Nitroaniline	13.802	65	165998	60.39	ng	#	79
49) Acenaphthylene	14.454	152	710315	46.62	ng		97
50) Dimethylphthalate	14.184	163	659116	49.79	ng		99
51) 2,6-Dinitrotoluene	14.296	165	140043	65.09	ng	#	75
52) Acenaphthene	14.795	154	573335	55.40	ng		96
53) 3-Nitroaniline	14.625	138	80433	33.08	ng	#	70
54) 2,4-Dinitrophenol	14.831	184	176283	150.03	ng	#	76
55) Dibenzofuran	15.130	168	754033	48.26	ng		95
56) 4-Nitrophenol	14.925	139	79049	39.48	ng	#	66
57) 2,4-Dinitrotoluene	15.083	165	201166	70.61	ng	#	84
58) Fluorene	15.776	166	616462	49.27	ng		98
59) 2,3,4,6-Tetrachlorophenol	15.348	232	203717	54.05	ng	#	95
60) Diethylphthalate	15.541	149	611716	45.96	ng		98
61) 4-Chlorophenyl-phenyle...	15.765	204	345236	48.33	ng		97
62) 4-Nitroaniline	15.788	138	137453	50.92	ng	#	66
63) Azobenzene	16.058	77	582587	44.53	ng		90
65) 4,6-Dinitro-2-methylph...	15.847	198	131830	78.61	ng		88
66) n-Nitrosodiphenylamine	15.982	169	561813	46.29	ng		98
67) 4-Bromophenyl-phenylether	16.664	248	233448	45.73	ng		97
68) Hexachlorobenzene	16.781	284	248458	45.24	ng		95
69) Atrazine	16.928	200	185351	38.01	ng		98
70) Pentachlorophenol	17.122	266	354785	111.90	ng		91
71) Phenanthrene	17.516	178	1025588	46.75	ng		98
72) Anthracene	17.610	178	1054831	48.74	ng		98
73) Carbazole	17.874	167	977966	49.65	ng		98
74) Di-n-butylphthalate	18.432	149	1070450	44.22	ng	#	97
75) Fluoranthene	19.519	202	1387233	49.66	ng		97
77) Benzidine	19.695	184	532601	36.82	ng		98
78) Pyrene	19.883	202	1389145	46.06	ng		99
80) Butylbenzylphthalate	20.765	149	538429	46.20	ng		90
81) Benzo(a)anthracene	21.734	228	1447343	46.20	ng		99
82) 3,3'-Dichlorobenzidine	21.640	252	322643	26.54	ng		96
83) Chrysene	21.799	228	1390539	46.46	ng		99
84) Bis(2-ethylhexyl)phtha...	21.634	149	759894	44.38	ng		97
85) Di-n-octyl phthalate	22.868	149	1325145	45.01	ng	#	94
87) Indeno(1,2,3-cd)pyrene	28.767	276	1722267	46.65	ng	#	88
88) Benzo(b)fluoranthene	23.984	252	1527090	46.70	ng	#	98
89) Benzo(k)fluoranthene	24.055	252	1465582	46.62	ng	#	97
90) Benzo(a)pyrene	24.872	252	1367960	44.46	ng	#	98
91) Dibenzo(a,h)anthracene	28.832	278	1428200	47.29	ng	#	96
92) Benzo(g,h,i)perylene	29.936	276	1440573	48.78	ng	#	94

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

