

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
 Data File : BG048038.D
 Acq On : 27 Jan 2021 21:37
 Operator : CG/JU
 Sample : M1220-03MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP1-BMSD

Manual Integrations
 APPROVED

mohammad
 1/28/2021 11:46:40 AM

Quant Time: Jan 27 23:31:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 25 19:39:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.123	152	57381	20.00	ng	0.00
21) Naphthalene-d8	10.937	136	223722	20.00	ng	0.00
39) Acenaphthene-d10	14.751	164	160789	20.00	ng	0.00
64) Phenanthrene-d10	17.489	188	380907	20.00	ng	# 0.00
76) Chrysene-d12	21.766	240	408335	20.00	ng	# 0.00
86) Perylene-d12	25.050	264	398958	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.685	112	361577	96.84	ng	0.00
7) Phenol-d6	7.277	99	495482	87.28	ng	0.00
23) Nitrobenzene-d5	9.292	82	329339	58.25	ng	0.00
42) 2,4,6-Tribromophenol	16.231	330	233655	87.25	ng	0.00
45) 2-Fluorobiphenyl	13.376	172	674363	54.23	ng	0.00
79) Terphenyl-d14	20.091	244	1134507	48.17	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.564	88	48655	29.23	ng	# 96
3) Pyridine	3.963	79	152936	30.61	ng	90
4) n-Nitrosodimethylamine	3.875	42	91163	39.49	ng	80
6) Aniline	7.441	93	156025	22.75	ng	94
8) 2-Chlorophenol	7.688	128	177639	45.45	ng	90
9) Benzaldehyde	7.253	77	119904	41.18	ng	89
10) Phenol	7.300	94	251642	46.25	ng	75
11) bis(2-Chloroethyl)ether	7.541	93	173653	40.75	ng	97
12) 1,3-Dichlorobenzene	8.017	146	186334	39.50	ng	# 93
13) 1,4-Dichlorobenzene	8.158	146	191197	40.43	ng	96
14) 1,2-Dichlorobenzene	8.481	146	186587	41.34	ng	97
15) Benzyl Alcohol	8.358	79	196662	40.95	ng	# 88
16) 2,2'-oxybis(1-Chloropr...	8.652	45	220574	36.88	ng	89
17) 2-Methylphenol	8.558	107	159085	43.26	ng	# 88
18) Hexachloroethane	9.216	117	72138	39.10	ng	100
19) n-Nitroso-di-n-propyla...	8.928	70	153808	37.12	ng	97
20) 3+4-Methylphenols	8.887	107	214648	42.19	ng	89
22) Acetophenone	8.946	105	277753	40.29	ng	93
24) Nitrobenzene	9.333	77	234888	41.48	ng	# 83
25) Isophorone	9.856	82	410941	40.50	ng	94
26) 2-Nitrophenol	10.044	139	104407	44.81	ng	# 75
27) 2,4-Dimethylphenol	10.097	122	160306	47.35	ng	87
28) bis(2-Chloroethoxy)met...	10.338	93	236338	38.71	ng	97
29) 2,4-Dichlorophenol	10.579	162	196088	45.02	ng	97
30) 1,2,4-Trichlorobenzene	10.802	180	219951	41.77	ng	98
31) Naphthalene	10.990	128	530804	41.03	ng	99
32) Benzoic acid	10.244	122	131213	43.28	ng	# 77
33) 4-Chloroaniline	11.090	127	49716	8.41	ng	# 91
34) Hexachlorobutadiene	11.278	225	156063	41.73	ng	98
35) Caprolactam	11.866	113	63448m	38.27	ng	
36) 4-Chloro-3-methylphenol	12.201	107	204431	41.83	ng	86
37) 2-Methylnaphthalene	12.588	142	395672	40.50	ng	# 92
38) 1-Methylnaphthalene	12.806	142	375818	40.32	ng	# 100
40) 1,2,4,5-Tetrachloroben...	12.953	216	270006	44.05	ng	# 97
41) Hexachlorocyclopentadiene	12.935	237	335983	96.77	ng	100
43) 2,4,6-Trichlorophenol	13.182	196	186748	43.68	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.252	196	203341	45.90	ng	# 95
46) 1,1'-Biphenyl	13.587	154	523140	42.09	ng	99
47) 2-Chloronaphthalene	13.628	162	429970	39.75	ng	99
48) 2-Nitroaniline	13.822	65	148559	37.84	ng	# 82
49) Acenaphthylene	14.474	152	646062	40.46	ng	99
50) Dimethylphthalate	14.198	163	604451	41.53	ng	99
51) 2,6-Dinitrotoluene	14.316	165	127654	42.20	ng	# 72
52) Acenaphthene	14.815	154	507153m	46.92	ng	
53) 3-Nitroaniline	14.645	138	61938	18.66	ng	# 80
54) 2,4-Dinitrophenol	14.850	184	179385	94.83	ng	# 71
55) Dibenzofuran	15.144	168	660261	40.28	ng	96
56) 4-Nitrophenol	14.944	139	219030	84.59	ng	# 60
57) 2,4-Dinitrotoluene	15.103	165	177411	40.60	ng	# 80
58) Fluorene	15.796	166	525749	39.89	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.368	232	191651	43.26	ng	# 97
60) Diethylphthalate	15.561	149	567051	37.90	ng	95
61) 4-Chlorophenyl-phenyle...	15.785	204	324548	40.21	ng	98
62) 4-Nitroaniline	15.808	138	122306	33.83	ng	# 61
63) Azobenzene	16.078	77	540440	38.37	ng	92
65) 4,6-Dinitro-2-methylph...	15.861	198	123370	50.19	ng	83
66) n-Nitrosodiphenylamine	15.996	169	495997	43.33	ng	99
67) 4-Bromophenyl-phenylether	16.678	248	216623	42.56	ng	96
68) Hexachlorobenzene	16.795	284	228442	41.27	ng	96
69) Atrazine	16.942	200	179190	41.21	ng	96
70) Pentachlorophenol	17.136	266	308493	91.79	ng	98
71) Phenanthrene	17.536	178	891352	40.56	ng	97
72) Anthracene	17.624	178	906635	41.95	ng	98
73) Carbazole	17.888	167	811941	40.25	ng	98
74) Di-n-butylphthalate	18.446	149	932380	37.86	ng	# 97
75) Fluoranthene	19.533	202	1151091	39.95	ng	97
77) Benzidine	19.709	184	317373	26.94	ng	99
78) Pyrene	19.897	202	1151422	38.51	ng	100
80) Butylbenzylphthalate	20.779	149	428690	37.53	ng	90
81) Benzo(a)anthracene	21.748	228	1201871	40.50	ng	99
82) 3,3'-Dichlorobenzidine	21.660	252	236198	23.62	ng	# 97
83) Chrysene	21.819	228	1164079	41.28	ng	99
84) Bis(2-ethylhexyl)phtha...	21.654	149	619937	39.63	ng	99
85) Di-n-octyl phthalate	22.888	149	1056846	40.39	ng	96
87) Indeno(1,2,3-cd)pyrene	28.805	276	1446355	45.07	ng	# 90
88) Benzo(b)fluoranthene	24.010	252	1280347	45.15	ng	# 98
89) Benzo(k)fluoranthene	24.081	252	1206613	43.79	ng	# 96
90) Benzo(a)pyrene	24.898	252	1148151	43.12	ng	# 98
91) Dibenzo(a,h)anthracene	28.881	278	1210268	45.70	ng	# 95
92) Benzo(g,h,i)perylene	29.992	276	1187087	46.53	ng	# 94

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

