

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG052720\
 Data File : BG045482.D
 Acq On : 27 May 2020 18:56
 Operator : CG/JU
 Sample : L2745-20MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM62-COMP-2020-0-6MSD

Manual Integrations
 APPROVED

mohammad
 5/29/2020 10:24:39 PM

Quant Time: May 28 01:22:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 27 13:40:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	91275	20.00	ng	0.00
21) Naphthalene-d8	10.77	136	377170	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	274249	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	598789	20.00	ng	0.00
76) Chrysene-d12	21.61	240	518479	20.00	ng	0.00
87) Perylene-d12	24.76	264	575972	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	499197	106.88	ng	0.00
7) Phenol-d6	7.18	99	725598	109.15	ng	0.00
23) Nitrobenzene-d5	9.12	82	475872	68.80	ng	0.00
42) 2,4,6-Tribromophenol	16.10	330	366454	104.48	ng	0.00
45) 2-Fluorobiphenyl	13.23	172	1125366	69.60	ng	0.00
79) Terphenyl-d14	19.96	244	1674517	75.33	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.40	88	82985	35.831 ng	97
3) Pyridine	3.80	79	187679	29.096 ng	98
4) n-Nitrosodimethylamine	3.72	42	88895	42.116 ng	92
6) Aniline	7.29	93	128515	14.810 ng	97
8) 2-Chlorophenol	7.54	128	257571	50.290 ng	99
9) Benzaldehyde	7.09	77	164517	37.986 ng	96
10) Phenol	7.21	94	336013	49.045 ng	98
11) bis(2-Chloroethyl)ether	7.38	93	232274	43.465 ng	96
12) 1,3-Dichlorobenzene	7.85	146	274006	44.518 ng	100
13) 1,4-Dichlorobenzene	7.99	146	274417	45.179 ng	97
14) 1,2-Dichlorobenzene	8.32	146	262349	44.907 ng	97
15) Benzyl Alcohol	8.21	79	257382	46.870 ng	97
16) 2,2'-oxybis(1-Chloropropan	8.49	45	212301	41.853 ng	98
17) 2-Methylphenol	8.44	107	236822	46.182 ng	96
18) Hexachloroethane	9.04	117	108229	46.523 ng	99
19) n-Nitroso-di-n-propylamine	8.77	70	206687	44.963 ng	99
20) 3+4-Methylphenols	8.77	107	307882	44.090 ng	# 63
22) Acetophenone	8.78	105	386661	43.254 ng	# 92
24) Nitrobenzene	9.16	77	318682	43.005 ng	99
25) Isophorone	9.69	82	577689	42.411 ng	97
26) 2-Nitrophenol	9.88	139	149758	47.242 ng	94
27) 2,4-Dimethylphenol	9.97	122	245702	52.259 ng	99
28) bis(2-Chloroethoxy)methane	10.17	93	323054	45.224 ng	96
29) 2,4-Dichlorophenol	10.44	162	273678	49.293 ng	98
30) 1,2,4-Trichlorobenzene	10.63	180	301154	45.904 ng	99
31) Naphthalene	10.82	128	795115	45.239 ng	99
32) Benzoic acid	10.16	122	178873	54.281 ng	89
33) 4-Chloroaniline	10.94	127	61387	8.356 ng	96
34) Hexachlorobutadiene	11.11	225	210569	45.406 ng	98
35) Caprolactam	11.70	113	89669m	44.001 ng	
36) 4-Chloro-3-methylphenol	12.10	107	306175	48.939 ng	98
37) 2-Methylnaphthalene	12.43	142	602590	45.999 ng	97
38) 1-Methylnaphthalene	12.64	142	571554	46.188 ng	95
40) 1,2,4,5-Tetrachlorobenzene	12.80	216	356714	46.567 ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.78	237	297077	67.192	nq	97
43) 2,4,6-Trichlorophenol	13.05	196	246524	46.328	nq	99
44) 2,4,5-Trichlorophenol	13.14	196	271010	44.568	nq	97
46) 1,1'-Biphenyl	13.44	154	776537	46.082	nq	99
47) 2-Chloronaphthalene	13.48	162	596913	43.621	nq	99
48) 2-Nitroaniline	13.68	65	184293	40.943	nq	89
49) Acenaphthylene	14.33	152	975078	44.362	nq	98
50) Dimethylphthalate	14.06	163	855476	45.472	nq	99
51) 2,6-Dinitrotoluene	14.18	165	178240	41.986	nq	98
52) Acenaphthene	14.67	154	703435m	47.811	nq	
53) 3-Nitroaniline	14.51	138	69900	16.198	nq	99
54) 2,4-Dinitrophenol	14.72	184	175984	64.424	nq	99
55) Dibenzofuran	15.00	168	942142	43.121	nq	96
56) 4-Nitrophenol	14.88	139	285989	87.533	nq	95
57) 2,4-Dinitrotoluene	14.97	165	251680	40.935	nq	97
58) Fluorene	15.65	166	770930	42.949	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.24	232	239442	44.758	nq	98
60) Diethylphthalate	15.42	149	793033	40.499	nq	99
61) 4-Chlorophenyl-phenylether	15.65	204	430399	41.902	nq	97
62) 4-Nitroaniline	15.68	138	149768	33.244	nq	98
63) Azobenzene	15.94	77	756456	41.430	nq	96
65) 4,6-Dinitro-2-methylphenol	15.73	198	135259	38.786	nq	98
66) n-Nitrosodiphenylamine	15.86	169	680091	47.304	nq	97
67) 4-Bromophenyl-phenylether	16.54	248	299007	46.115	nq	94
68) Hexachlorobenzene	16.66	284	322990	47.627	nq	99
69) Atrazine	16.82	200	249730	42.172	nq	99
70) Pentachlorophenol	17.02	266	360210	102.295	nq	98
71) Phenanthrene	17.40	178	1238310	47.371	nq	99
72) Anthracene	17.49	178	1222894	46.585	nq	96
73) Carbazole	17.76	167	1094458	42.771	nq	100
74) Di-n-butylphthalate	18.31	149	1288843	41.965	nq	99
75) Fluoranthene	19.41	202	1527202	45.932	nq	98
77) Benzidine	19.58	184	519190	35.654	nq	99
78) Pyrene	19.76	202	1489404	50.394	nq	98
80) Butylbenzylphthalate	20.65	149	554649	43.478	nq	98
81) Benzo(a)anthracene	21.59	228	1397345	48.380	nq	99
82) 3,3'-Dichlorobenzidine	21.51	252	306230	27.029	nq	98
83) Chrysene	21.66	228	1312625	47.264	nq	97
84) Bis(2-ethylhexyl)phthalate	21.50	149	763832	43.557	nq	98
85) Di-n-octyl phthalate	22.70	149	1282712	42.588	nq	100
86) Indeno(1,2,3-cd)pyrene	28.34	276	1732267	47.699	nq	# 91
88) Benzo(b)fluoranthene	23.77	252	1492180	48.306	nq	99
89) Benzo(k)fluoranthene	23.84	252	1374602	47.366	nq	99
90) Benzo(a)pyrene	24.62	252	1352156	47.001	nq	97
91) Dibenzo(a,h)anthracene	28.41	278	1368010	47.320	nq	99
92) Benzo(g,h,i)perylene	29.46	276	1430565	49.478	nq	98

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

