

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG113018\
 Data File : BG038264.D
 Acq On : 30 Nov 2018 19:20
 Operator : JU/SJ
 Sample : J5761-02MSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AU-06-112918MSD

Manual Integrations
 APPROVED

mohammad
 12/3/2018 3:33:19 PM

Quant Time: Dec 03 04:41:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	32121	20.00	ng	0.00
21) Naphthalene-d8	10.90	136	136110	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	83146	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	192559	20.00	ng	0.00
76) Chrysene-d12	21.75	240	180194	20.00	ng	0.00
87) Perylene-d12	25.07	264	174035	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	215595	115.89	ng	0.00
7) Phenol-d6	7.23	99	268523	101.12	ng	0.00
23) Nitrobenzene-d5	9.26	82	219103	86.17	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	116107	122.44	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	458146	80.27	ng	0.00
79) Terphenyl-d14	20.06	244	752761	98.30	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.52	88	29206	33.236	ng	# 1
3) Pyridine	3.93	79	75772	30.228	ng	92
4) n-Nitrosodimethylamine	3.84	42	47329	52.044	ng	81
6) Aniline	7.41	93	26613	7.765	ng	# 96
8) 2-Chlorophenol	7.64	128	106020	49.113	ng	99
9) Benzaldehyde	7.22	77	69461	36.169	ng	99
10) Phenol	7.26	94	107671	38.278	ng	93
11) bis(2-Chloroethyl)ether	7.50	93	100824	43.905	ng	99
12) 1,3-Dichlorobenzene	7.97	146	107171	43.095	ng	99
13) 1,4-Dichlorobenzene	8.12	146	107863	42.937	ng	99
14) 1,2-Dichlorobenzene	8.44	146	104044	43.381	ng	98
15) Benzyl Alcohol	8.33	79	94751	49.309	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.61	45	219222	51.410	ng	98
17) 2-Methylphenol	8.52	107	91133	47.921	ng	96
18) Hexachloroethane	9.17	117	39766	43.496	ng	92
19) n-Nitroso-di-n-propylamine	8.88	70	84891	44.179	ng	94
20) 3+4-Methylphenols	8.85	107	122618	45.483	ng	98
22) Acetophenone	8.91	105	159924	47.224	ng	# 97
24) Nitrobenzene	9.31	77	129047	49.613	ng	95
25) Isophorone	9.82	82	237820	51.965	ng	96
26) 2-Nitrophenol	10.01	139	60938	46.929	ng	94
27) 2,4-Dimethylphenol	10.06	122	105723	54.038	ng	95
28) bis(2-Chloroethoxy)methane	10.30	93	140496	48.009	ng	97
29) 2,4-Dichlorophenol	10.54	162	110425	50.179	ng	97
30) 1,2,4-Trichlorobenzene	10.76	180	112938	45.798	ng	96
31) Naphthalene	10.95	128	333748	49.853	ng	98
32) Benzoic acid	10.16	122	18234m	23.088	ng	
33) 4-Chloroaniline	11.07	127	7100	2.280	ng	# 89
34) Hexachlorobutadiene	11.22	225	68391	45.062	ng	100
35) Caprolactam	11.84	113	24519	27.836	ng	94
36) 4-Chloro-3-methylphenol	12.17	107	118216	49.171	ng	99
37) 2-Methylnaphthalene	12.55	142	245616	51.077	ng	98
38) 1-Methylnaphthalene	12.77	142	234861	50.740	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.91	216	128155	46.126	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	132635	89.152	nq	99
43) 2,4,6-Trichlorophenol	13.15	196	91065	48.099	nq	98
44) 2,4,5-Trichlorophenol	13.23	196	93618	46.995	nq	97
46) 1,1'-Biphenyl	13.55	154	314534	45.029	nq	99
47) 2-Chloronaphthalene	13.60	162	249658	46.838	nq	99
48) 2-Nitroaniline	13.81	65	87823	52.351	nq	97
49) Acenaphthylene	14.45	152	408688	50.987	nq	99
50) Dimethylphthalate	14.17	163	319650	48.494	nq	100
51) 2,6-Dinitrotoluene	14.30	165	72759	48.771	nq	95
52) Acenaphthene	14.78	154	234667	48.630	nq	99
53) 3-Nitroaniline	14.64	138	10700	6.500	nq	95
54) 2,4-Dinitrophenol	14.84	184	7632	17.663	nq	98
55) Dibenzofuran	15.12	168	382794	47.973	nq	96
56) 4-Nitrophenol	14.93	139	68908	54.274	nq	88
57) 2,4-Dinitrotoluene	15.08	165	100445	49.485	nq	97
58) Fluorene	15.76	166	307563	51.661	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.33	232	79400	47.633	nq	99
60) Diethylphthalate	15.52	149	329785	47.104	nq	98
61) 4-Chlorophenyl-phenylether	15.75	204	162267	46.580	nq	98
62) 4-Nitroaniline	15.80	138	74831	45.758	nq	95
63) Azobenzene	16.05	77	313289	50.047	nq	97
65) 4,6-Dinitro-2-methylphenol	15.84	198	13802	11.332	nq	89
66) n-Nitrosodiphenylamine	15.97	169	278336	47.779	nq	99
67) 4-Bromophenyl-phenylether	16.64	248	102587	48.539	nq	98
68) Hexachlorobenzene	16.76	284	103543	47.463	nq	98
69) Atrazine	16.91	200	101304	47.174	nq	99
70) Pentachlorophenol	17.11	266	54260	36.622	nq	96
71) Phenanthrene	17.51	178	503322	51.053	nq	99
72) Anthracene	17.60	178	515265	53.021	nq	100
73) Carbazole	17.87	167	463098	47.496	nq	99
74) Di-n-butylphthalate	18.40	149	560567	51.217	nq	98
75) Fluoranthene	19.51	202	591495	52.586	nq	98
77) Benzidine	19.70	184	42593	6.736	nq	96
78) Pyrene	19.88	202	602875	51.173	nq	98
80) Butylbenzylphthalate	20.75	149	248776	48.286	nq	99
81) Benzo(a)anthracene	21.73	228	564236	51.816	nq	99
82) 3,3'-Dichlorobenzidine	21.64	252	43967	10.365	nq	96
83) Chrysene	21.80	228	531700	51.033	nq	99
84) Bis(2-ethylhexyl)phthalate	21.60	149	354004	48.181	nq	97
85) Di-n-octyl phthalate	22.83	149	579102	48.719	nq	99
86) Indeno(1,2,3-cd)pyrene	28.87	276	398213	34.413	nq	# 84
88) Benzo(b)fluoranthene	24.00	252	521602	54.463	nq	99
89) Benzo(k)fluoranthene	24.08	252	532530	56.350	nq	99
90) Benzo(a)pyrene	24.91	252	499646	54.590	nq	98
91) Dibenzo(a,h)anthracene	28.94	278	298770	33.926	nq	99
92) Benzo(g,h,i)perylene	30.06	276	316174	36.107	nq	96

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

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