

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\
 Data File : BG032582.D
 Acq On : 7 Feb 2018 3:41
 Operator : SJ/JU
 Sample : J1458-01MSD
 Misc : MDL-W-8 ppm
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-01MSD

Manual Integrations
 APPROVED

Sohil
 2/7/2018 4:15:04 PM

Quant Time: Feb 07 05:45:40 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 18:57:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.71	152	18891	20.00	ng	0.00
21) Naphthalene-d8	11.58	136	71151	20.00	ng	0.00
38) Acenaphthene-d10	15.34	164	49450	20.00	ng	0.00
63) Phenanthrene-d10	18.07	188	161673	20.00	ng	0.00
75) Chrysene-d12	22.40	240	200535	20.00	ng	0.00
86) Perylene-d12	26.07	264	245005	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.14	112	142476	151.49	ng	0.00
7) Phenol-d6	7.83	99	220366	175.58	ng	0.00
23) Nitrobenzene-d5	9.90	82	134685	118.23	ng	0.00
41) 2,4,6-Tribromophenol	16.82	330	184298	195.79	ng	0.00
44) 2-Fluorobiphenyl	13.97	172	355379	85.42	ng	0.00
78) Terphenyl-d14	20.62	244	884201	94.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.80	88	14832	46.871	ng	# 70
3) Pyridine	4.26	79	39363	46.068	ng	# 78
4) n-Nitrosodimethylamine	4.16	42	22379	50.313	ng	# 72
6) Aniline	8.00	93	36625	23.405	ng	96
8) 2-Chlorophenol	8.26	128	71322	64.152	ng	# 81
9) Benzaldehyde	7.81	77	14240	21.968	ng	95
10) Phenol	7.86	94	83454	70.020	ng	91
11) bis(2-Chloroethyl)ether	8.09	93	50971	56.127	ng	# 76
12) 1,3-Dichlorobenzene	8.59	146	80434m	53.498	ng	
13) 1,4-Dichlorobenzene	8.74	146	69080	45.841	ng	89
14) 1,2-Dichlorobenzene	9.08	146	77149	52.261	ng	97
15) Benzyl Alcohol	8.95	79	55651	53.802	ng	97
16) 2,2'-oxybis(1-Chloropropan	9.23	45	46735	54.310	ng	65
17) 2-Methylphenol	9.15	107	51856	53.570	ng	97
18) Hexachloroethane	9.83	117	23226	45.906	ng	92
19) n-Nitroso-di-n-propylamine	9.52	70	34622	41.605	ng	# 83
20) 3+4-Methylphenols	9.48	107	58789	43.297	ng	99
22) Acetophenone	9.55	105	77330	46.365	ng	# 90
24) Nitrobenzene	9.95	77	64682	58.721	ng	# 85
25) Isophorone	10.47	82	122660	63.238	ng	# 86
26) 2-Nitrophenol	10.67	139	34564	51.908	ng	# 81
27) 2,4-Dimethylphenol	10.72	122	54137	56.645	ng	90
28) bis(2-Chloroethoxy)methane	10.96	93	55923	52.577	ng	# 95
29) 2,4-Dichlorophenol	11.22	162	65045	51.375	ng	92
30) 1,2,4-Trichlorobenzene	11.44	180	72395	42.902	ng	96
31) Naphthalene	11.64	128	171165	49.255	ng	99
32) Benzoic acid	10.77	122	3859m	11.875	ng	
33) 4-Chloroaniline	11.74	127	37815	24.398	ng	# 91
34) Hexachlorobutadiene	11.90	225	55576	42.331	ng	93
35) Caprolactam	12.48	113	16296	44.836	ng	# 48
36) 4-Chloro-3-methylphenol	12.82	107	58888	49.035	ng	# 89
37) 2-Methylnaphthalene	13.21	142	141139	48.344	ng	96
39) 1,2,4,5-Tetrachlorobenzene	13.56	216	96405	41.672	ng	98
40) Hexachlorocyclopentadiene	13.53	237	137896	95.401	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.79	196	57442	45.648	nq	97
43) 2,4,5-Trichlorophenol	13.79	196	57442	39.159	nq	95
45) 1,1'-Biphenyl	14.18	154	194005	46.437	nq	97
46) 2-Chloronaphthalene	14.23	162	149602	45.687	nq	97
47) 2-Nitroaniline	14.43	65	31864	44.004	nq #	69
48) Acenaphthylene	15.07	152	230591	46.582	nq	99
49) Dimethylphthalate	14.78	163	229705	50.379	nq	98
50) 2,6-Dinitrotoluene	14.90	165	44358	46.352	nq #	75
51) Acenaphthene	15.40	154	161457	47.030	nq	97
52) 3-Nitroaniline	15.24	138	26960	32.253	nq #	67
53) 2,4-Dinitrophenol	15.44	184	25359	44.378	nq #	62
54) Dibenzofuran	15.73	168	241708	47.568	nq	100
55) 4-Nitrophenol	15.53	139	67081	107.194	nq #	78
56) 2,4-Dinitrotoluene	15.68	165	65653	48.183	nq #	71
57) Fluorene	16.38	166	237900	56.342	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.95	232	72904	50.390	nq #	87
59) Diethylphthalate	16.11	149	209538	47.185	nq	98
60) 4-Chlorophenyl-phenylether	16.36	204	145415	54.614	nq	93
61) 4-Nitroaniline	16.40	138	48169	53.765	nq	83
62) Azobenzene	16.65	77	153512	57.846	nq #	82
64) 4,6-Dinitro-2-methylphenol	16.44	198	49532	41.010	nq #	69
65) n-Nitrosodiphenylamine	16.57	169	222524	46.120	nq	97
66) 4-Bromophenyl-phenylether	17.25	248	97655	40.772	nq	96
67) Hexachlorobenzene	17.38	284	110399	41.651	nq #	87
68) Atrazine	17.50	200	101666	49.085	nq	94
69) Pentachlorophenol	17.72	266	124506	74.983	nq	93
70) Phenanthrene	18.12	178	414004	45.919	nq	99
71) Anthracene	18.21	178	428019	47.678	nq	98
72) Carbazole	18.47	167	376006	47.404	nq	100
73) Di-n-butylphthalate	18.98	149	432481	49.246	nq	99
74) Fluoranthene	20.08	202	552438	46.722	nq	94
76) Benzidine	20.25	184	154146	39.146	nq	97
77) Pyrene	20.44	202	539400	43.855	nq	98
79) Butylbenzylphthalate	21.29	149	192797	49.695	nq #	81
80) Benzo(a)anthracene	22.38	228	580235	46.672	nq	99
81) 3,3'-Dichlorobenzidine	22.27	252	137199	28.484	nq #	96
82) Chrysene	22.45	228	547203	45.577	nq	98
83) Bis(2-ethylhexyl)phthalate	22.22	149	270688	49.208	nq #	98
84) Di-n-octyl phthalate	23.58	149	385391	44.171	nq #	88
85) Indeno(1,2,3-cd)pyrene	30.32	276	835306	54.997	nq #	79
87) Benzo(b)fluoranthene	24.89	252	641082	43.305	nq #	96
88) Benzo(k)fluoranthene	24.97	252	610181	41.608	nq #	97
89) Benzo(a)pyrene	25.90	252	667356	47.286	nq #	94
90) Dibenzo(a,h)anthracene	30.40	278	715238	49.540	nq #	92
91) Benzo(g,h,i)perylene	31.66	276	684175	47.743	nq #	91

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

