

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062117\
 Data File : BG027584.D
 Acq On : 22 Jun 2017 7:05
 Operator : SJ/JU
 Sample : PB99997BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99997BS

Manual Integrations
 APPROVED

mohammad
 6/22/2017 6:04:34 PM

Quant Time: Jun 22 08:02:46 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 18:52:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.51	152	20124	20.00	ng	0.00
21) Naphthalene-d8	11.32	136	98054	20.00	ng	0.00
38) Acenaphthene-d10	15.09	164	66955	20.00	ng	0.00
63) Phenanthrene-d10	17.84	188	182863	20.00	ng	0.00
75) Chrysene-d12	22.21	240	205926	20.00	ng	0.00
86) Perylene-d12	25.84	264	191308	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.09	112	164043	116.02	ng	0.00
7) Phenol-d6	7.64	99	239169	114.26	ng	0.00
23) Nitrobenzene-d5	9.66	82	148211	71.52	ng	0.00
41) 2,4,6-Tribromophenol	16.57	330	124176	124.40	ng	0.00
44) 2-Fluorobiphenyl	13.71	172	349646	71.33	ng	0.00
78) Terphenyl-d14	20.41	244	691746	79.15	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	4.09	88	14847	27.045	ng	96
3) Pyridine	4.47	79	47863	26.932	ng	# 79
4) n-Nitrosodimethylamine	4.38	42	27362	33.736	ng	# 72
6) Aniline	7.83	93	80700	31.590	ng	96
8) 2-Chlorophenol	8.07	128	57109	38.171	ng	96
9) Benzaldehyde	7.65	77	35933	25.263	ng	92
10) Phenol	7.67	94	81040	38.004	ng	94
11) bis(2-Chloroethyl)ether	7.91	93	55201	33.074	ng	90
12) 1,3-Dichlorobenzene	8.40	146	51982	32.914	ng	98
13) 1,4-Dichlorobenzene	8.54	146	52408	32.868	ng	99
14) 1,2-Dichlorobenzene	8.86	146	52721	33.648	ng	94
15) Benzyl Alcohol	8.73	79	65322	39.085	ng	95
16) 2,2'-oxybis(1-Chloropropan	9.01	45	84456	30.498	ng	93
17) 2-Methylphenol	8.92	107	54449	36.287	ng	92
18) Hexachloroethane	9.59	117	20401	32.627	ng	89
19) n-Nitroso-di-n-propylamine	9.29	70	53047	31.905	ng	90
20) 3+4-Methylphenols	9.24	107	75445	36.723	ng	91
22) Acetophenone	9.32	105	93251	34.679	ng	# 89
24) Nitrobenzene	9.71	77	75776	33.361	ng	95
25) Isophorone	10.23	82	147914	34.742	ng	98
26) 2-Nitrophenol	10.42	139	29493	35.104	ng	93
27) 2,4-Dimethylphenol	10.46	122	64359	40.563	ng	98
28) bis(2-Chloroethoxy)methane	10.70	93	82696	34.463	ng	97
29) 2,4-Dichlorophenol	10.96	162	56331	41.104	ng	91
30) 1,2,4-Trichlorobenzene	11.18	180	55404	35.361	ng	96
31) Naphthalene	11.37	128	179452	33.860	ng	98
32) Benzoic acid	10.56	122	41234	38.372	ng	96
33) 4-Chloroaniline	11.47	127	55905	24.872	ng	95
34) Hexachlorobutadiene	11.64	225	33147	35.007	ng	98
35) Caprolactam	12.24	113	26415m	41.356	ng	
36) 4-Chloro-3-methylphenol	12.55	107	88315	42.020	ng	99
37) 2-Methylnaphthalene	12.94	142	139497	36.218	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.30	216	70642	33.645	ng	# 97
40) Hexachlorocyclopentadiene	13.27	237	94552	84.015	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.53	196	53354	39.325	ng	97
43) 2,4,5-Trichlorophenol	13.60	196	60131	37.715	ng	95
45) 1,1'-Biphenyl	13.92	154	191398	34.283	ng	98
46) 2-Chloronaphthalene	13.98	162	143102	34.464	ng	95
47) 2-Nitroaniline	14.17	65	69361	37.210	ng	97
48) Acenaphthylene	14.82	152	242053	33.268	ng	97
49) Dimethylphthalate	14.52	163	220666	37.640	ng	96
50) 2,6-Dinitrotoluene	14.65	165	49481	38.897	ng	95
51) Acenaphthene	15.15	154	197225	38.904	ng	99
52) 3-Nitroaniline	14.99	138	38112	28.384	ng	99
53) 2,4-Dinitrophenol	15.19	184	59388	82.599	ng #	88
54) Dibenzofuran	15.48	168	248242	37.595	ng	98
55) 4-Nitrophenol	15.28	139	89672	79.765	ng #	79
56) 2,4-Dinitrotoluene	15.43	165	74811	41.474	ng #	96
57) Fluorene	16.13	166	218647	34.963	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.70	232	61210	43.599	ng	91
59) Diethylphthalate	15.88	149	242611	37.926	ng	96
60) 4-Chlorophenyl-phenylether	16.11	204	111010	36.073	ng	94
61) 4-Nitroaniline	16.15	138	58574	39.761	ng #	85
62) Azobenzene	16.40	77	252480	38.492	ng	93
64) 4,6-Dinitro-2-methylphenol	16.20	198	42477	36.579	ng	77
65) n-Nitrosodiphenylamine	16.33	169	198314	35.377	ng	100
66) 4-Bromophenyl-phenylether	17.01	248	72449	36.024	ng #	87
67) Hexachlorobenzene	17.14	284	76543	35.614	ng	97
68) Atrazine	17.26	200	79023	38.481	ng	98
69) Pentachlorophenol	17.48	266	101532	76.052	ng	96
70) Phenanthrene	17.88	178	376938	36.234	ng #	92
71) Anthracene	17.97	178	379651	36.247	ng	94
72) Carbazole	18.24	167	337252	33.051	ng	96
73) Di-n-butylphthalate	18.76	149	424778	37.154	ng #	95
74) Fluoranthene	19.87	202	442461	36.507	ng	93
76) Benzidine	20.04	184	224932	44.505	ng	98
77) Pyrene	20.23	202	457873	36.577	ng	94
79) Butylbenzylphthalate	21.11	149	196537	37.617	ng	97
80) Benzo(a)anthracene	22.19	228	445367	36.046	ng	93
81) 3,3'-Dichlorobenzidine	22.08	252	120460	26.966	ng	98
82) Chrysene	22.26	228	417309	35.645	ng	93
83) Bis(2-ethylhexyl)phthalate	22.04	149	280004	36.562	ng	97
84) Di-n-octyl phthalate	23.39	149	481138	37.771	ng #	90
85) Indeno(1,2,3-cd)pyrene	30.03	276	504736	37.832	ng #	94
87) Benzo(b)fluoranthene	24.67	252	439226	35.909	ng #	95
88) Benzo(k)fluoranthene	24.75	252	440093	36.776	ng #	92
89) Benzo(a)pyrene	25.67	252	424738	36.795	ng #	94
90) Dibenzo(a,h)anthracene	30.10	278	433324	36.854	ng #	96
91) Benzo(g,h,i)perylene	31.34	276	416799	36.495	ng #	90

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

