

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG062217\
 Data File : BG027627.D
 Acq On : 23 Jun 2017 11:53
 Operator : SJ/JU
 Sample : PB100013BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100013BS

Manual Integrations
 APPROVED

mohammad
 6/26/2017 8:33:08 AM

Quant Time: Jun 23 16:07:21 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	15035	20.00	ng	0.00
21) Naphthalene-d8	11.33	136	70737	20.00	ng	0.00
38) Acenaphthene-d10	15.09	164	48876	20.00	ng	0.00
63) Phenanthrene-d10	17.84	188	146694	20.00	ng	0.00
75) Chrysene-d12	22.21	240	175852	20.00	ng	0.00
86) Perylene-d12	25.84	264	163703	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	6.09	112	136132	128.86	ng	0.00
7) Phenol-d6	7.64	99	187960	120.19	ng	0.00
23) Nitrobenzene-d5	9.67	82	119255	79.77	ng	0.00
41) 2,4,6-Tribromophenol	16.58	330	116702	160.16	ng	0.00
44) 2-Fluorobiphenyl	13.71	172	288622	80.66	ng	0.00
78) Terphenyl-d14	20.41	244	645985	86.55	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	4.09	88	14450	35.23	ng	97
3) Pyridine	4.47	79	41549	31.29	ng	# 83
4) n-Nitrosodimethylamine	4.38	42	24876	41.05	ng	# 87
6) Aniline	7.83	93	33717	17.67	ng	95
8) 2-Chlorophenol	8.07	128	49571	44.35	ng	94
9) Benzaldehyde	7.65	77	11698	11.01	ng	90
10) Phenol	7.67	94	67583	42.42	ng	96
11) bis(2-Chloroethyl)ether	7.91	93	43639	35.00	ng	88
12) 1,3-Dichlorobenzene	8.39	146	46565	39.46	ng	94
13) 1,4-Dichlorobenzene	8.54	146	47101	39.54	ng	96
14) 1,2-Dichlorobenzene	8.86	146	45876	39.19	ng	98
15) Benzyl Alcohol	8.73	79	51032	40.87	ng	96
16) 2,2'-oxybis(1-Chloropropan	9.01	45	63822	30.85	ng	88
17) 2-Methylphenol	8.92	107	44347	39.56	ng	94
18) Hexachloroethane	9.59	117	18377	39.34	ng	85
19) n-Nitroso-di-n-propylamine	9.29	70	42643	34.33	ng	89
20) 3+4-Methylphenols	9.25	107	62640	40.81	ng	92
22) Acetophenone	9.32	105	76808	39.59	ng	# 90
24) Nitrobenzene	9.71	77	64176	39.16	ng	# 90
25) Isophorone	10.23	82	114179	37.18	ng	99
26) 2-Nitrophenol	10.42	139	24956	41.18	ng	95
27) 2,4-Dimethylphenol	10.46	122	52691	46.03	ng	93
28) bis(2-Chloroethoxy)methane	10.70	93	66339	38.32	ng	97
29) 2,4-Dichlorophenol	10.96	162	49073	49.64	ng	95
30) 1,2,4-Trichlorobenzene	11.18	180	48092	42.55	ng	95
31) Naphthalene	11.37	128	152676	39.93	ng	98
32) Benzoic acid	10.57	122	32091	40.78	ng	95
33) 4-Chloroaniline	11.48	127	9810	6.05	ng	# 89
34) Hexachlorobutadiene	11.64	225	30908	45.25	ng	90
35) Caprolactam	12.24	113	21676m	47.04	ng	
36) 4-Chloro-3-methylphenol	12.55	107	74779	49.32	ng	99
37) 2-Methylnaphthalene	12.94	142	120531	43.38	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.30	216	64119	41.83	ng	96
40) Hexachlorocyclopentadiene	13.28	237	82014	99.83	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.53	196	46721	47.17	nq	99
43) 2,4,5-Trichlorophenol	13.61	196	54415	46.75	nq	94
45) 1,1'-Biphenyl	13.92	154	164476	40.36	nq	96
46) 2-Chloronaphthalene	13.98	162	124582	41.10	nq	94
47) 2-Nitroaniline	14.17	65	57629	42.35	nq	95
48) Acenaphthylene	14.82	152	208170	39.19	nq	99
49) Dimethylphthalate	14.53	163	192072	44.88	nq	# 95
50) 2,6-Dinitrotoluene	14.65	165	42508	45.78	nq	98
51) Acenaphthene	15.15	154	171885	46.45	nq	99
52) 3-Nitroaniline	14.99	138	14880	15.18	nq	# 80
53) 2,4-Dinitrophenol	15.19	184	53031	101.04	nq	# 83
54) Dibenzofuran	15.49	168	218191	45.27	nq	98
55) 4-Nitrophenol	15.28	139	76148	92.79	nq	# 87
56) 2,4-Dinitrotoluene	15.44	165	67742	51.45	nq	# 98
57) Fluorene	16.13	166	192271	42.12	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.70	232	57278	55.89	nq	92
59) Diethylphthalate	15.88	149	217915	46.67	nq	97
60) 4-Chlorophenyl-phenylether	16.11	204	100850	44.89	nq	95
61) 4-Nitroaniline	16.15	138	49586	46.11	nq	88
62) Azobenzene	16.41	77	217657	45.46	nq	95
64) 4,6-Dinitro-2-methylphenol	16.20	198	39425	42.32	nq	86
65) n-Nitrosodiphenylamine	16.33	169	179376	39.89	nq	98
66) 4-Bromophenyl-phenylether	17.01	248	68592	42.51	nq	# 89
67) Hexachlorobenzene	17.14	284	72847	42.25	nq	96
68) Atrazine	17.27	200	73620	44.69	nq	98
69) Pentachlorophenol	17.48	266	96747	90.34	nq	97
70) Phenanthrene	17.88	178	346356	41.50	nq	94
71) Anthracene	17.97	178	345683	41.14	nq	97
72) Carbazole	18.24	167	309263	37.78	nq	97
73) Di-n-butylphthalate	18.76	149	394332	42.99	nq	96
74) Fluoranthene	19.87	202	415661	42.75	nq	94
76) Benzidine	20.04	184	98212	22.76	nq	98
77) Pyrene	20.23	202	430040	40.23	nq	95
79) Butylbenzylphthalate	21.11	149	184507	41.35	nq	94
80) Benzo(a)anthracene	22.18	228	434667	41.20	nq	93
81) 3,3'-Dichlorobenzidine	22.08	252	81012	21.24	nq	# 98
82) Chrysene	22.26	228	403739	40.38	nq	95
83) Bis(2-ethylhexyl)phthalate	22.04	149	264095	40.38	nq	97
84) Di-n-octyl phthalate	23.39	149	455868	41.91	nq	# 90
85) Indeno(1,2,3-cd)pyrene	30.03	276	498828	43.78	nq	# 94
87) Benzo(b)fluoranthene	24.68	252	433465	41.41	nq	# 97
88) Benzo(k)fluoranthene	24.76	252	431889	42.18	nq	# 93
89) Benzo(a)pyrene	25.67	252	416358	42.15	nq	# 96
90) Dibenzo(a,h)anthracene	30.10	278	424480	42.19	nq	# 97
91) Benzo(g,h,i)perylene	31.35	276	411544	42.11	nq	# 94

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

