

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG073116\
 Data File : BG023359.D
 Acq On : 30 Jul 2016 16:07
 Operator : UM/SJ
 Sample : PB92558BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB92558BS

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:12:30 PM

Quant Time: Aug 01 02:00:20 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG072616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 27 17:30:33 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	151231	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	732046	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	527167	20.00	ng	0.00
63) Phenanthrene-d10	17.57	188	1166856	20.00	ng	0.00
75) Chrysene-d12	21.89	240	1050367	20.00	ng	0.00
86) Perylene-d12	25.30	264	1078730	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.68	112	1336407	150.86	ng	0.00
7) Phenol-d6	7.29	99	2000601	147.10	ng	0.00
23) Nitrobenzene-d5	9.31	82	1314527	87.81	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	915693	167.82	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	2626492	93.25	ng	0.00
78) Terphenyl-d14	20.17	244	2924063	90.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	124226	32.70	ng	# 90
3) Pyridine	3.93	79	432572	32.53	ng	93
4) n-Nitrosodimethylamine	3.85	42	201353	33.51	ng	# 67
6) Aniline	7.46	93	648603	35.70	ng	94
8) 2-Chlorophenol	7.70	128	519980	52.61	ng	96
9) Benzaldehyde	7.26	77	256538	26.26	ng	95
10) Phenol	7.32	94	764636	53.28	ng	80
11) bis(2-Chloroethyl)ether	7.55	93	516733	42.36	ng	92
12) 1,3-Dichlorobenzene	8.03	146	488061m	43.82	ng	
13) 1,4-Dichlorobenzene	8.17	146	497414	43.77	ng	95
14) 1,2-Dichlorobenzene	8.50	146	486861	44.42	ng	94
15) Benzyl Alcohol	8.38	79	563204	61.53	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.67	45	748122	38.54	ng	89
17) 2-Methylphenol	8.58	107	518085	52.82	ng	# 88
18) Hexachloroethane	9.23	117	193316	40.83	ng	93
19) n-Nitroso-di-n-propylamine	8.94	70	521873	42.66	ng	97
20) 3+4-Methylphenols	8.92	107	734094	55.04	ng	92
22) Acetophenone	8.97	105	753335	39.42	ng	# 93
24) Nitrobenzene	9.36	77	721453	41.50	ng	# 86
25) Isophorone	9.88	82	1404603	45.43	ng	# 93
26) 2-Nitrophenol	10.08	139	316663	48.23	ng	# 77
27) 2,4-Dimethylphenol	10.13	122	586063	52.96	ng	85
28) bis(2-Chloroethoxy)methane	10.36	93	807322	46.54	ng	99
29) 2,4-Dichlorophenol	10.62	162	614070	57.31	ng	97
30) 1,2,4-Trichlorobenzene	10.83	180	537516	45.40	ng	99
31) Naphthalene	11.02	128	1566387	46.05	ng	98
32) Benzoic acid	10.25	122	205433	27.06	ng	# 90
33) 4-Chloroaniline	11.14	127	175690	11.18	ng	# 87
34) Hexachlorobutadiene	11.30	225	336454	43.36	ng	97
35) Caprolactam	11.92	113	219324m	48.14	ng	
36) 4-Chloro-3-methylphenol	12.26	107	740993	52.64	ng	93
37) 2-Methylnaphthalene	12.63	142	1228442	48.71	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	13.00	216	591523	34.66	ng	98
40) Hexachlorocyclopentadiene	12.97	237	813100	100.54	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.24	196	523729	52.25	ng	97
43) 2,4,5-Trichlorophenol	13.32	196	570214	53.62	ng	91
45) 1,1'-Biphenyl	13.64	154	1520224	41.35	ng	96
46) 2-Chloronaphthalene	13.68	162	1329219	45.12	ng	94
47) 2-Nitroaniline	13.89	65	572537	45.94	ng	# 86
48) Acenaphthylene	14.53	152	2095905	46.47	ng	95
49) Dimethylphthalate	14.25	163	1798194	46.22	ng	97
50) 2,6-Dinitrotoluene	14.38	165	430931	47.30	ng	# 82
51) Acenaphthene	14.87	154	1340399	45.81	ng	98
52) 3-Nitroaniline	14.72	138	211683	23.07	ng	# 82
53) 2,4-Dinitrophenol	14.92	184	443231	83.09	ng	91
54) Dibenzofuran	15.20	168	2067546	49.94	ng	98
55) 4-Nitrophenol	15.03	139	603687	75.85	ng	# 81
56) 2,4-Dinitrotoluene	15.17	165	598913	46.69	ng	# 86
57) Fluorene	15.86	166	1706583	46.68	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.43	232	509780	56.62	ng	94
59) Diethylphthalate	15.62	149	1771922	44.47	ng	94
60) 4-Chlorophenyl-phenylether	15.84	204	902062	45.58	ng	99
61) 4-Nitroaniline	15.89	138	427094	44.39	ng	# 65
62) Azobenzene	16.14	77	1974548	47.79	ng	91
64) 4,6-Dinitro-2-methylphenol	15.94	198	348561	46.51	ng	94
65) n-Nitrosodiphenylamine	16.06	169	1585305	50.89	ng	96
66) 4-Bromophenyl-phenylether	16.74	248	631117	53.49	ng	97
67) Hexachlorobenzene	16.87	284	646361	54.49	ng	95
68) Atrazine	17.01	200	487510	40.80	ng	97
69) Pentachlorophenol	17.22	266	660536	90.22	ng	98
70) Phenanthrene	17.61	178	2502029	49.02	ng	94
71) Anthracene	17.71	178	2546776	49.39	ng	92
72) Carbazole	17.98	167	2179799	46.15	ng	97
73) Di-n-butylphthalate	18.52	149	2269047	43.06	ng	96
74) Fluoranthene	19.62	202	2511537	43.12	ng	94
76) Benzidine	19.80	184	929558	31.47	ng	98
77) Pyrene	19.99	202	2491043	41.79	ng	96
79) Butylbenzylphthalate	20.86	149	1204853	45.95	ng	99
80) Benzo(a)anthracene	21.87	228	2536294	46.12	ng	99
81) 3,3'-Dichlorobenzidine	21.78	252	710029	31.84	ng	# 96
82) Chrysene	21.94	228	2357371	44.37	ng	95
83) Bis(2-ethylhexyl)phthalate	21.75	149	1669025	45.49	ng	98
84) Di-n-octyl phthalate	23.02	149	2762625	46.03	ng	100
85) Indeno(1,2,3-cd)pyrene	29.19	276	3241235	52.30	ng	# 100
87) Benzo(b)fluoranthene	24.21	252	2719745	47.46	ng	# 95
88) Benzo(k)fluoranthene	24.27	252	2653536	46.32	ng	# 94
89) Benzo(a)pyrene	25.13	252	2656981	47.31	ng	# 96
90) Dibenzo(a,h)anthracene	29.26	278	2697190	50.77	ng	# 90
91) Benzo(g,h,i)perylene	30.40	276	2700594	48.29	ng	# 92

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

