

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG073116\
 Data File : BG023379.D
 Acq On : 31 Jul 2016 22:20
 Operator : UM/SJ
 Sample : H4223-04MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-32-16-17MSD

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:13:32 PM

Quant Time: Aug 01 03:06:31 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	8.12	152	147242	20.00	ng	-0.02
21) Naphthalene-d8	10.95	136	751491	20.00	ng	-0.03
38) Acenaphthene-d10	14.79	164	561127	20.00	ng	-0.02
63) Phenanthrene-d10	17.54	188	1374780	20.00	ng	-0.03
75) Chrysene-d12	21.86	240	1133038	20.00	ng	-0.03
86) Perylene-d12	25.24	264	1127117	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	1134947	131.59	ng	0.00
7) Phenol-d6	7.28	99	1745446	131.81	ng	0.00
23) Nitrobenzene-d5	9.30	82	1135518	73.89	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	894983	154.10	ng	-0.02
44) 2-Fluorobiphenyl	13.40	172	2317111	77.29	ng	-0.02
78) Terphenyl-d14	20.15	244	2898644	83.16	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	100607	27.20	ng	# 92
3) Pyridine	3.93	79	348318	26.90	ng	93
4) n-Nitrosodimethylamine	3.84	42	160412	27.42	ng	# 72
6) Aniline	7.44	93	762508	43.11	ng	94
8) 2-Chlorophenol	7.69	128	449428	46.70	ng	99
9) Benzaldehyde	7.25	77	305157	32.08	ng	94
10) Phenol	7.31	94	705363	50.48	ng	# 80
11) bis(2-Chloroethyl)ether	7.53	93	450648	37.95	ng	92
12) 1,3-Dichlorobenzene	8.01	146	402322	37.10	ng	95
13) 1,4-Dichlorobenzene	8.16	146	415941	37.60	ng	93
14) 1,2-Dichlorobenzene	8.48	146	411054	38.52	ng	92
15) Benzyl Alcohol	8.36	79	507218	56.92	ng	# 90
16) 2,2'-oxybis(1-Chloropropan	8.66	45	651354	34.47	ng	89
17) 2-Methylphenol	8.57	107	466104	48.81	ng	# 83
18) Hexachloroethane	9.21	117	160351	34.79	ng	93
19) n-Nitroso-di-n-propylamine	8.93	70	471402	39.58	ng	98
20) 3+4-Methylphenols	8.90	107	660002	50.83	ng	87
22) Acetophenone	8.95	105	679644	34.64	ng	# 88
24) Nitrobenzene	9.34	77	642385	36.00	ng	# 86
25) Isophorone	9.87	82	1288194	40.59	ng	# 93
26) 2-Nitrophenol	10.06	139	289227	42.91	ng	# 75
27) 2,4-Dimethylphenol	10.11	122	521588	45.91	ng	84
28) bis(2-Chloroethoxy)methane	10.35	93	744603	41.81	ng	99
29) 2,4-Dichlorophenol	10.60	162	532744	48.44	ng	97
30) 1,2,4-Trichlorobenzene	10.82	180	478061	39.33	ng	98
31) Naphthalene	11.01	128	1432447	41.02	ng	98
32) Benzoic acid	10.11	122	521588	54.97	ng	# 7
33) 4-Chloroaniline	11.12	127	577715	35.81	ng	93
34) Hexachlorobutadiene	11.29	225	283783	35.62	ng	93
35) Caprolactam	11.90	113	224054m	47.90	ng	
36) 4-Chloro-3-methylphenol	12.24	107	666882	46.15	ng	# 88
37) 2-Methylnaphthalene	12.61	142	1143067	44.15	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	12.97	216	535386	29.47	ng	98
40) Hexachlorocyclopentadiene	12.95	237	669210	77.74	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.22	196	473895	44.41	ng	98
43) 2,4,5-Trichlorophenol	13.30	196	497330	43.94	ng	91
45) 1,1'-Biphenyl	13.61	154	1378912	35.23	ng	92
46) 2-Chloronaphthalene	13.66	162	1213005	38.68	ng	95
47) 2-Nitroaniline	13.87	65	548162	41.32	ng	88
48) Acenaphthylene	14.51	152	1954144	40.70	ng	97
49) Dimethylphthalate	14.24	163	2257531	54.52	ng	# 94
50) 2,6-Dinitrotoluene	14.36	165	413861	42.68	ng	# 77
51) Acenaphthene	14.85	154	1260171	40.46	ng	99
52) 3-Nitroaniline	14.69	138	424740	43.49	ng	# 78
53) 2,4-Dinitrophenol	14.90	184	186289	32.81	ng	# 86
54) Dibenzofuran	15.19	168	1954538	44.35	ng	95
55) 4-Nitrophenol	15.01	139	716956	84.63	ng	# 74
56) 2,4-Dinitrotoluene	15.15	165	604870	44.30	ng	# 89
57) Fluorene	15.84	166	1647393	42.34	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.41	232	499607	52.14	ng	95
59) Diethylphthalate	15.59	149	1757039	41.43	ng	96
60) 4-Chlorophenyl-phenylether	15.82	204	849170	40.31	ng	98
61) 4-Nitroaniline	15.87	138	477526	46.63	ng	# 61
62) Azobenzene	16.12	77	1890478	42.99	ng	90
64) 4,6-Dinitro-2-methylphenol	15.92	198	371513	42.07	ng	94
65) n-Nitrosodiphenylamine	16.04	169	1549196	42.21	ng	93
66) 4-Bromophenyl-phenylether	16.72	248	608083	43.74	ng	97
67) Hexachlorobenzene	16.84	284	640453	45.82	ng	96
68) Atrazine	16.99	200	638427	45.35	ng	95
69) Pentachlorophenol	17.20	266	749225	86.86	ng	98
70) Phenanthrene	17.59	178	2579027m	42.89	ng	
71) Anthracene	17.68	178	2623107	43.17	ng	93
72) Carbazole	17.95	167	2471277	44.41	ng	95
73) Di-n-butylphthalate	18.49	149	2675363	43.09	ng	# 95
74) Fluoranthene	19.60	202	2791714	40.68	ng	90
76) Benzidine	19.78	184	1882429	59.07	ng	97
77) Pyrene	19.96	202	2759828	42.92	ng	93
79) Butylbenzylphthalate	20.84	149	1144284	40.45	ng	96
80) Benzo(a)anthracene	21.84	228	2394244	40.36	ng	98
81) 3,3'-Dichlorobenzidine	21.75	252	845885	35.17	ng	# 98
82) Chrysene	21.91	228	2265742	39.53	ng	96
83) Bis(2-ethylhexyl)phthalate	21.72	149	1522393	38.46	ng	98
84) Di-n-octyl phthalate	22.98	149	2534777	39.15	ng	99
85) Indeno(1,2,3-cd)pyrene	29.10	276	2974499	44.49	ng	# 100
87) Benzo(b)fluoranthene	24.16	252	2537075	42.37	ng	# 97
88) Benzo(k)fluoranthene	24.23	252	2390242	39.93	ng	# 96
89) Benzo(a)pyrene	25.08	252	2441007	41.60	ng	# 97
90) Dibenzo(a,h)anthracene	29.18	278	2450783	44.16	ng	# 91
91) Benzo(g,h,i)perylene	30.33	276	2450584	41.94	ng	# 91

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

