

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081616\
 Data File : BG023743.D
 Acq On : 17 Aug 2016 00:57
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

umangi
 8/17/2016 3:10:49 AM

Quant Time: Aug 17 01:55:51 2016
 Quant Method : Z:\HPCHEM1\BNA_G\Methods\8270-BG080116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 19:22:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	114674	20.00	ng	-0.02
21) Naphthalene-d8	10.94	136	534727	20.00	ng	-0.02
38) Acenaphthene-d10	14.78	164	341825	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	737993	20.00	ng	0.00
75) Chrysene-d12	21.87	240	714439	20.00	ng	0.00
86) Perylene-d12	25.28	264	709295	20.00	ng	0.04

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	530264	77.84	ng	-0.02
7) Phenol-d6	7.26	99	821460	77.42	ng	-0.02
23) Nitrobenzene-d5	9.28	82	847256	78.77	ng	-0.02
41) 2,4,6-Tribromophenol	16.28	330	380460	76.09	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	1685533	83.17	ng	0.00
78) Terphenyl-d14	20.16	244	1990390	87.82	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.50	88	95738	32.98	ng	94
3) Pyridine	3.91	79	303500	31.82	ng	95
4) n-Nitrosodimethylamine	3.82	42	133491	37.75	ng	# 84
6) Aniline	7.42	93	508108	33.14	ng	96
8) 2-Chlorophenol	7.67	128	300847	38.44	ng	97
9) Benzaldehyde	7.23	77	247096	41.24	ng	96
10) Phenol	7.29	94	418647	36.88	ng	83
11) bis(2-Chloroethyl)ether	7.51	93	325023	35.90	ng	92
12) 1,3-Dichlorobenzene	7.99	146	346867	38.71	ng	94
13) 1,4-Dichlorobenzene	8.14	146	356278	38.84	ng	95
14) 1,2-Dichlorobenzene	8.46	146	335685	37.31	ng	96
15) Benzyl Alcohol	8.35	79	332854	41.40	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.64	45	492287	36.97	ng	88
17) 2-Methylphenol	8.55	107	287384	37.76	ng	90
18) Hexachloroethane	9.19	117	138498	40.11	ng	97
19) n-Nitroso-di-n-propylamine	8.91	70	325696	35.68	ng	98
20) 3+4-Methylphenols	8.89	107	401472	37.42	ng	92
22) Acetophenone	8.94	105	531629	36.31	ng	# 92
24) Nitrobenzene	9.33	77	485184	40.74	ng	# 90
25) Isophorone	9.85	82	860385	38.40	ng	# 94
26) 2-Nitrophenol	10.05	139	204452	40.67	ng	# 81
27) 2,4-Dimethylphenol	10.10	122	320719	37.47	ng	86
28) bis(2-Chloroethoxy)methane	10.33	93	453388	33.66	ng	99
29) 2,4-Dichlorophenol	10.59	162	345707	40.17	ng	99
30) 1,2,4-Trichlorobenzene	10.80	180	373017	40.19	ng	97
31) Naphthalene	10.99	128	1022024	37.53	ng	100
32) Benzoic acid	10.28	122	265806	35.65	ng	96
33) 4-Chloroaniline	11.10	127	427557	32.85	ng	95
34) Hexachlorobutadiene	11.27	225	257588	42.20	ng	97
35) Caprolactam	11.90	113	115171	31.81	ng	91
36) 4-Chloro-3-methylphenol	12.24	107	415456	36.74	ng	91
37) 2-Methylnaphthalene	12.60	142	760424m	36.73	ng	
39) 1,2,4,5-Tetrachlorobenzene	12.96	216	534423	43.11	ng	98
40) Hexachlorocyclopentadiene	12.94	237	48189	7.71	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.21	196	305934	41.43	ng	96
43) 2,4,5-Trichlorophenol	13.29	196	307127	40.40	ng	92
45) 1,1'-Biphenyl	13.61	154	1043503	39.93	ng	96
46) 2-Chloronaphthalene	13.65	162	801052	39.63	ng	96
47) 2-Nitroaniline	13.86	65	316653	37.73	ng	91
48) Acenaphthylene	14.50	152	1257710	39.36	ng	98
49) Dimethylphthalate	14.23	163	1028803	39.23	ng	98
50) 2,6-Dinitrotoluene	14.36	165	234821	37.81	ng	85
51) Acenaphthene	14.84	154	796395	39.34	ng	99
52) 3-Nitroaniline	14.70	138	238269	34.02	ng	82
53) 2,4-Dinitrophenol	14.92	184	75344	18.92	ng	# 88
54) Dibenzofuran	15.18	168	1125459	38.29	ng	96
55) 4-Nitrophenol	15.02	139	178540	32.46	ng	# 82
56) 2,4-Dinitrotoluene	15.16	165	332114	38.10	ng	94
57) Fluorene	15.84	166	983066	38.33	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.41	232	255451	36.62	ng	95
59) Diethylphthalate	15.58	149	1040116	39.07	ng	100
60) 4-Chlorophenyl-phenylether	15.82	204	528319	38.93	ng	97
61) 4-Nitroaniline	15.87	138	241363	32.37	ng	# 70
62) Azobenzene	16.11	77	1029449	38.15	ng	92
64) 4,6-Dinitro-2-methylphenol	15.93	198	148451	29.59	ng	94
65) n-Nitrosodiphenylamine	16.04	169	862247	39.83	ng	94
66) 4-Bromophenyl-phenylether	16.72	248	345526	40.17	ng	99
67) Hexachlorobenzene	16.85	284	367834	39.09	ng	97
68) Atrazine	16.99	200	330180	39.72	ng	97
69) Pentachlorophenol	17.20	266	185150	33.75	ng	98
70) Phenanthrene	17.59	178	1448296	38.68	ng	97
71) Anthracene	17.68	178	1487905	39.50	ng	99
72) Carbazole	17.96	167	1303133	39.40	ng	100
73) Di-n-butylphthalate	18.49	149	1586707	47.78	ng	97
74) Fluoranthene	19.60	202	1588317	45.13	ng	97
76) Benzidine	19.79	184	865897	43.60	ng	99
77) Pyrene	19.97	202	1599649	38.21	ng	98
79) Butylbenzylphthalate	20.84	149	719628	41.83	ng	98
80) Benzo(a)anthracene	21.85	228	1538427	39.01	ng	97
81) 3,3'-Dichlorobenzidine	21.76	252	644274	39.83	ng	# 97
82) Chrysene	21.92	228	1457479	38.84	ng	99
83) Bis(2-ethylhexyl)phthalate	21.72	149	1002634	42.56	ng	96
84) Di-n-octyl phthalate	22.99	149	1677837	41.73	ng	100
85) Indeno(1,2,3-cd)pyrene	29.19	276	1818373	38.78	ng	# 100
87) Benzo(b)fluoranthene	24.19	252	1556148	38.99	ng	# 96
88) Benzo(k)fluoranthene	24.26	252	1488627	38.84	ng	# 98
89) Benzo(a)pyrene	25.11	252	1500644	39.54	ng	# 98
90) Dibenzo(a,h)anthracene	29.25	278	1560862	40.26	ng	# 96
91) Benzo(g,h,i)perylene	30.41	276	1434220	36.73	ng	# 92

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

