

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110716\
 Data File : BG024756.D
 Acq On : 8 Nov 2016 12:53
 Operator : UM/SJ
 Sample : PB94569BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB94569BSD

Manual Integrations
 APPROVED

Umangi
 11/8/2016 5:59:47 PM

Quant Time: Nov 08 15:11:27 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 26 11:14:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	83814	20.00	ng	0.00
21) Naphthalene-d8	11.09	136	322805	20.00	ng	0.00
38) Acenaphthene-d10	14.88	164	274150	20.00	ng	-0.02
63) Phenanthrene-d10	17.62	188	675295	20.00	ng	0.00
75) Chrysene-d12	21.92	240	924061	20.00	ng	0.00
86) Perylene-d12	25.35	264	966173	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.79	112	345375	67.54	ng	0.00
7) Phenol-d6	7.40	99	497766	70.96	ng	0.00
23) Nitrobenzene-d5	9.44	82	605922	85.46	ng	-0.02
41) 2,4,6-Tribromophenol	16.36	330	311503	76.06	ng	0.00
44) 2-Fluorobiphenyl	13.50	172	1331730	80.74	ng	-0.02
78) Terphenyl-d14	20.20	244	2674902	86.50	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.65	88	71089	34.03	ng	92
3) Pyridine	4.06	79	216608	34.63	ng	88
4) n-Nitrosodimethylamine	3.98	42	112811	34.84	ng	82
6) Aniline	7.58	93	114955	12.94	ng	98
8) 2-Chlorophenol	7.82	128	208023	40.01	ng	98
9) Benzaldehyde	7.40	77	31441	7.25	ng	87
10) Phenol	7.43	94	309510	42.80	ng	96
11) bis(2-Chloroethyl)ether	7.67	93	213049	39.22	ng	88
12) 1,3-Dichlorobenzene	8.14	146	239866	37.27	ng	94
13) 1,4-Dichlorobenzene	8.30	146	247549	38.94	ng	94
14) 1,2-Dichlorobenzene	8.62	146	242011	39.59	ng	97
15) Benzyl Alcohol	8.50	79	259063	39.70	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.78	45	355230	35.24	ng	93
17) 2-Methylphenol	8.69	107	208616	43.54	ng	95
18) Hexachloroethane	9.35	117	91474	37.43	ng	92
19) n-Nitroso-di-n-propylamine	9.06	70	199319	38.56	ng	93
20) 3+4-Methylphenols	9.01	107	274805	42.72	ng	89
22) Acetophenone	9.09	105	342573	41.31	ng	# 89
24) Nitrobenzene	9.48	77	306142	38.81	ng	98
25) Isophorone	9.99	82	537682	40.77	ng	98
26) 2-Nitrophenol	10.19	139	123072	42.04	ng	94
27) 2,4-Dimethylphenol	10.23	122	232263	45.32	ng	97
28) bis(2-Chloroethoxy)methane	10.48	93	296241	43.41	ng	100
29) 2,4-Dichlorophenol	10.72	162	251014	46.66	ng	96
30) 1,2,4-Trichlorobenzene	10.95	180	261367	40.96	ng	96
31) Naphthalene	11.14	128	645689	40.10	ng	98
32) Benzoic acid	10.33	122	112453	26.35	ng	94
33) 4-Chloroaniline	11.24	127	76212	10.90	ng	# 90
34) Hexachlorobutadiene	11.40	225	206379	41.32	ng	98
35) Caprolactam	12.00	113	94408m	47.94	ng	
36) 4-Chloro-3-methylphenol	12.34	107	295439	45.49	ng	# 92
37) 2-Methylnaphthalene	12.73	142	496905	42.29	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.08	216	346480	37.47	ng	97
40) Hexachlorocyclopentadiene	13.05	237	488359	93.79	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.32	196	235288	42.33	nq	95
43) 2,4,5-Trichlorophenol	13.40	196	254419	42.34	nq #	91
45) 1,1'-Biphenyl	13.71	154	711488	37.40	nq	94
46) 2-Chloronaphthalene	13.77	162	566706	39.12	nq	99
47) 2-Nitroaniline	13.97	65	233220	39.32	nq	93
48) Acenaphthylene	14.61	152	858867	38.66	nq	99
49) Dimethylphthalate	14.32	163	789308	40.55	nq	97
50) 2,6-Dinitrotoluene	14.45	165	173065	41.65	nq	93
51) Acenaphthene	14.95	154	572044	34.54	nq	98
52) 3-Nitroaniline	14.78	138	69849	16.25	nq #	96
53) 2,4-Dinitrophenol	14.99	184	250792	83.30	nq	99
54) Dibenzofuran	15.28	168	942167	41.15	nq	99
55) 4-Nitrophenol	15.08	139	318363	84.99	nq #	89
56) 2,4-Dinitrotoluene	15.23	165	264785	43.86	nq #	95
57) Fluorene	15.92	166	761654	40.12	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.49	232	290934	46.70	nq #	91
59) Diethylphthalate	15.67	149	833421	40.84	nq	98
60) 4-Chlorophenyl-phenylether	15.90	204	437497	41.07	nq	97
61) 4-Nitroaniline	15.95	138	185760	39.55	nq #	79
62) Azobenzene	16.20	77	826147	40.60	nq	93
64) 4,6-Dinitro-2-methylphenol	15.99	198	185165	39.80	nq	98
65) n-Nitrosodiphenylamine	16.12	169	717208	38.85	nq	98
66) 4-Bromophenyl-phenylether	16.80	248	324889	39.63	nq	93
67) Hexachlorobenzene	16.92	284	362486	40.02	nq #	87
68) Atrazine	17.06	200	345365	43.61	nq	99
69) Pentachlorophenol	17.26	266	419954	70.26	nq	95
70) Phenanthrene	17.67	178	1379272	40.07	nq	96
71) Anthracene	17.76	178	1411603	40.65	nq	100
72) Carbazole	18.03	167	1275518	40.28	nq	98
73) Di-n-butylphthalate	18.55	149	1534191	42.01	nq	97
74) Fluoranthene	19.66	202	1857393	42.69	nq	97
76) Benzidine	19.83	184	547076	25.13	nq	96
77) Pyrene	20.02	202	1857721	41.13	nq	99
79) Butylbenzylphthalate	20.88	149	768952	40.83	nq	96
80) Benzo(a)anthracene	21.90	228	2043600	40.26	nq	98
81) 3,3'-Dichlorobenzidine	21.80	252	350747	17.13	nq	99
82) Chrysene	21.97	228	1917090	39.65	nq	98
83) Bis(2-ethylhexyl)phthalate	21.76	149	1041170	38.65	nq	100
84) Di-n-octyl phthalate	23.03	149	1774974	39.70	nq #	93
85) Indeno(1,2,3-cd)pyrene	29.29	276	2531792	37.94	nq #	99
87) Benzo(b)fluoranthene	24.25	252	2078662m	39.80	nq	
88) Benzo(k)fluoranthene	24.32	252	2021554	40.08	nq	97
89) Benzo(a)pyrene	25.18	252	2024036	39.66	nq	100
90) Dibenzo(a,h)anthracene	29.36	278	2128566	38.00	nq	98
91) Benzo(g,h,i)perylene	30.53	276	2123029	37.94	nq	98

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

