

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
 Data File : BG038778.D
 Acq On : 21 Dec 2018 1:28
 Operator : JU/SJ
 Sample : PB115875BSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115875BSD

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:43:43 PM

Quant Time: Dec 21 04:23:20 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	27165	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	106707	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	69913	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	172531	20.00	ng	0.00
76) Chrysene-d12	21.72	240	173996	20.00	ng	0.00
87) Perylene-d12	25.02	264	188883	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	128317	83.78	ng	0.00
7) Phenol-d6	7.21	99	172908	82.24	ng	0.00
23) Nitrobenzene-d5	9.24	82	152984	73.24	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	84996	88.21	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	398813	78.26	ng	0.00
79) Terphenyl-d14	20.04	244	629230	78.70	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	18582	26.068	ng	# 87
3) Pyridine	3.89	79	54672	27.858	ng	96
4) n-Nitrosodimethylamine	3.82	42	34414	35.788	ng	# 87
6) Aniline	7.38	93	30699	11.438	ng	98
8) 2-Chlorophenol	7.62	128	69590	39.263	ng	99
9) Benzaldehyde	7.20	77	53990	39.583	ng	98
10) Phenol	7.24	94	86844	38.878	ng	96
11) bis(2-Chloroethyl)ether	7.48	93	56865	31.975	ng	99
12) 1,3-Dichlorobenzene	7.94	146	72769	34.724	ng	99
13) 1,4-Dichlorobenzene	8.09	146	74510	34.716	ng	96
14) 1,2-Dichlorobenzene	8.41	146	70400	34.793	ng	97
15) Benzyl Alcohol	8.30	79	61339	37.376	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.57	45	121879	29.917	ng	98
17) 2-Methylphenol	8.50	107	58794	39.285	ng	99
18) Hexachloroethane	9.14	117	25791	32.885	ng	93
19) n-Nitroso-di-n-propylamine	8.86	70	46877	31.744	ng	97
20) 3+4-Methylphenols	8.83	107	80931	39.156	ng	96
22) Acetophenone	8.89	105	99053	37.164	ng	# 97
24) Nitrobenzene	9.28	77	76291	36.105	ng	98
25) Isophorone	9.79	82	134931	35.954	ng	99
26) 2-Nitrophenol	9.99	139	38097	35.227	ng	97
27) 2,4-Dimethylphenol	10.03	122	67929	43.827	ng	98
28) bis(2-Chloroethoxy)methane	10.27	93	77163	36.435	ng	97
29) 2,4-Dichlorophenol	10.52	162	75679	43.890	ng	97
30) 1,2,4-Trichlorobenzene	10.73	180	78909	37.886	ng	98
31) Naphthalene	10.93	128	210651	40.066	ng	99
32) Benzoic acid	10.19	122	28941m	33.336	ng	
33) 4-Chloroaniline	11.04	127	16842	7.378	ng	94
34) Hexachlorobutadiene	11.18	225	52145	37.000	ng	94
35) Caprolactam	11.83	113	23314	32.672	ng	# 88
36) 4-Chloro-3-methylphenol	12.15	107	77035	39.150	ng	94
37) 2-Methylnaphthalene	12.52	142	159391	42.495	ng	96
38) 1-Methylnaphthalene	12.74	142	152314	41.848	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	97708	37.388	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	106677	74.301	nq	97
43) 2,4,6-Trichlorophenol	13.13	196	64641m	40.229	nq	
44) 2,4,5-Trichlorophenol	13.21	196	70241	41.287	nq	97
46) 1,1'-Biphenyl	13.53	154	212605	36.275	nq	100
47) 2-Chloronaphthalene	13.58	162	169838	37.514	nq	97
48) 2-Nitroaniline	13.79	65	52237	36.224	nq	97
49) Acenaphthylene	14.42	152	273715	40.442	nq	98
50) Dimethylphthalate	14.15	163	216972	38.003	nq	98
51) 2,6-Dinitrotoluene	14.27	165	49220	38.042	nq	93
52) Acenaphthene	14.76	154	158127	39.072	nq	100
53) 3-Nitroaniline	14.61	138	17765	13.469	nq	# 84
54) 2,4-Dinitrophenol	14.82	184	34406	46.979	nq	97
55) Dibenzofuran	15.09	168	266661	38.438	nq	97
56) 4-Nitrophenol	14.92	139	72706	72.666	nq	98
57) 2,4-Dinitrotoluene	15.06	165	68246	37.450	nq	97
58) Fluorene	15.74	166	214904	41.812	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.31	232	66366	43.359	nq	95
60) Diethylphthalate	15.50	149	220253	35.142	nq	98
61) 4-Chlorophenyl-phenylether	15.73	204	121886	38.008	nq	98
62) 4-Nitroaniline	15.78	138	50702	37.340	nq	97
63) Azobenzene	16.02	77	185160	35.364	nq	97
65) 4,6-Dinitro-2-methylphenol	15.83	198	35424	28.782	nq	85
66) n-Nitrosodiphenylamine	15.94	169	196474	38.757	nq	96
67) 4-Bromophenyl-phenylether	16.62	248	81645	38.748	nq	97
68) Hexachlorobenzene	16.74	284	85138	38.978	nq	97
69) Atrazine	16.89	200	78000	41.461	nq	98
70) Pentachlorophenol	17.09	266	79268	61.355	nq	97
71) Phenanthrene	17.49	178	369861	41.340	nq	98
72) Anthracene	17.58	178	374807	42.435	nq	99
73) Carbazole	17.85	167	328634	37.622	nq	99
74) Di-n-butylphthalate	18.38	149	377927	37.706	nq	99
75) Fluoranthene	19.49	202	454914	41.095	nq	97
77) Benzidine	19.68	184	165040	32.073	nq	98
78) Pyrene	19.86	202	453308	39.436	nq	100
80) Butylbenzylphthalate	20.72	149	169350	37.710	nq	100
81) Benzo(a)anthracene	21.70	228	443483	41.828	nq	99
82) 3,3'-Dichlorobenzidine	21.62	252	69514	16.531	nq	99
83) Chrysene	21.77	228	414586	40.597	nq	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	238558	37.455	nq	99
85) Di-n-octyl phthalate	22.80	149	406300	38.090	nq	99
86) Indeno(1,2,3-cd)pyrene	28.80	276	494598	41.195	nq	# 81
88) Benzo(b)fluoranthene	23.96	252	432442	41.699	nq	99
89) Benzo(k)fluoranthene	24.03	252	433819	42.009	nq	97
90) Benzo(a)pyrene	24.87	252	421887	42.169	nq	# 98
91) Dibenzo(a,h)anthracene	28.87	278	414717	41.632	nq	98
92) Benzo(g,h,i)perylene	30.00	276	405617	41.317	nq	98

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

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