

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073018\
 Data File : BG035981.D
 Acq On : 30 Jul 2018 18:16
 Operator : JU/SJ
 Sample : PB111568BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111568BS

Manual Integrations
 APPROVED

Sohil
 7/31/2018 12:20:47 PM

Quant Time: Jul 31 01:49:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 25 11:04:14 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	32238	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	123176	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	92034	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	267708	20.00	ng	0.00
75) Chrysene-d12	21.66	240	286657	20.00	ng	0.00
86) Perylene-d12	24.85	264	275040	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	166307	85.65	ng	0.00
7) Phenol-d6	7.19	99	243564	84.07	ng	0.00
23) Nitrobenzene-d5	9.19	82	238633	80.93	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	118742	97.89	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	580635	84.52	ng	0.00
78) Terphenyl-d14	19.99	244	1042231	80.14	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	28719	29.059	ng	# 70
3) Pyridine	3.91	79	78818	30.549	ng	# 72
4) n-Nitrosodimethylamine	3.82	42	35996	32.493	ng	# 83
6) Aniline	7.35	93	69325	18.462	ng	90
8) 2-Chlorophenol	7.59	128	87004	44.055	ng	98
9) Benzaldehyde	7.16	77	58695	33.019	ng	98
10) Phenol	7.22	94	128584	43.931	ng	93
11) bis(2-Chloroethyl)ether	7.44	93	83897	36.601	ng	92
12) 1,3-Dichlorobenzene	7.91	146	93932	39.387	ng	# 89
13) 1,4-Dichlorobenzene	8.06	146	95495	39.410	ng	98
14) 1,2-Dichlorobenzene	8.37	146	92544	39.756	ng	90
15) Benzyl Alcohol	8.26	79	101232	38.479	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.55	45	96788	50.365	ng	72
17) 2-Methylphenol	8.47	107	86828	43.632	ng	93
18) Hexachloroethane	9.10	117	36203	39.075	ng	92
19) n-Nitroso-di-n-propylamine	8.82	70	81790	33.526	ng	# 81
20) 3+4-Methylphenols	8.79	107	116839	42.322	ng	93
22) Acetophenone	8.84	105	148533	38.777	ng	# 92
24) Nitrobenzene	9.23	77	124260	41.904	ng	# 93
25) Isophorone	9.74	82	236899	43.281	ng	99
26) 2-Nitrophenol	9.94	139	51203	48.619	ng	# 91
27) 2,4-Dimethylphenol	10.00	122	94208	52.846	ng	84
28) bis(2-Chloroethoxy)methane	10.23	93	123959	41.099	ng	100
29) 2,4-Dichlorophenol	10.48	162	107621	52.886	ng	95
30) 1,2,4-Trichlorobenzene	10.69	180	109491	43.878	ng	95
31) Naphthalene	10.88	128	274188	42.733	ng	99
32) Benzoic acid	10.17	122	28930m	24.106	ng	
33) 4-Chloroaniline	10.99	127	25957	9.282	ng	95
34) Hexachlorobutadiene	11.15	225	83570	42.949	ng	98
35) Caprolactam	11.76	113	36831m	42.175	ng	
36) 4-Chloro-3-methylphenol	12.11	107	127871	46.879	ng	92
37) 2-Methylnaphthalene	12.48	142	217046	45.404	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.85	216	151004	43.471	ng	99
40) Hexachlorocyclopentadiene	12.82	237	170709	93.706	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.09	196	94494	46.516	nq	99
43) 2,4,5-Trichlorophenol	13.16	196	107066	47.113	nq	97
45) 1,1'-Biphenyl	13.48	154	307499	43.095	nq	98
46) 2-Chloronaphthalene	13.53	162	243483	43.869	nq	100
47) 2-Nitroaniline	13.73	65	88491	45.099	nq	97
48) Acenaphthylene	14.37	152	410685	44.173	nq	98
49) Dimethylphthalate	14.10	163	348480	43.217	nq	98
50) 2,6-Dinitrotoluene	14.22	165	80486	49.016	nq	91
51) Acenaphthene	14.71	154	240350	41.500	nq	97
52) 3-Nitroaniline	14.56	138	29431	17.536	nq #	80
53) 2,4-Dinitrophenol	14.77	184	55891	67.326	nq #	67
54) Dibenzofuran	15.04	168	410228	44.538	nq	94
55) 4-Nitrophenol	14.87	139	112707	94.299	nq #	87
56) 2,4-Dinitrotoluene	15.01	165	120943	51.340	nq #	85
57) Fluorene	15.69	166	347891	45.064	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.28	232	110881	50.888	nq #	92
59) Diethylphthalate	15.46	149	359568	43.366	nq	98
60) 4-Chlorophenyl-phenylether	15.68	204	201505	44.410	nq	97
61) 4-Nitroaniline	15.72	138	80269	45.723	nq #	83
62) Azobenzene	15.98	77	358274	43.807	nq	95
64) 4,6-Dinitro-2-methylphenol	15.78	198	66608	44.696	nq	84
65) n-Nitrosodiphenylamine	15.90	169	314906	41.326	nq	96
66) 4-Bromophenyl-phenylether	16.58	248	137427	44.248	nq	92
67) Hexachlorobenzene	16.70	284	142333	44.501	nq	96
68) Atrazine	16.85	200	147230	46.928	nq	96
69) Pentachlorophenol	17.05	266	128359	65.887	nq	94
70) Phenanthrene	17.43	178	590489	44.204	nq	97
71) Anthracene	17.53	178	616001	46.312	nq	97
72) Carbazole	17.79	167	556273	44.038	nq	98
73) Di-n-butylphthalate	18.34	149	642228	43.024	nq #	97
74) Fluoranthene	19.44	202	794933	44.179	nq	96
76) Benzidine	19.62	184	195925	25.998	nq	97
77) Pyrene	19.80	202	797564	44.002	nq	97
79) Butylbenzylphthalate	20.68	149	289630	44.684	nq #	96
80) Benzo(a)anthracene	21.63	228	795377	44.525	nq	99
81) 3,3'-Dichlorobenzidine	21.55	252	147800	23.729	nq #	98
82) Chrysene	21.70	228	741786	44.811	nq	98
83) Bis(2-ethylhexyl)phthalate	21.53	149	404620	45.917	nq #	98
84) Di-n-octyl phthalate	22.73	149	688735	46.679	nq #	93
85) Indeno(1,2,3-cd)pyrene	28.49	276	837560	45.700	nq #	86
87) Benzo(b)fluoranthene	23.84	252	765756	45.807	nq #	97
88) Benzo(k)fluoranthene	23.91	252	728452	44.573	nq #	97
89) Benzo(a)pyrene	24.70	252	738759	47.502	nq #	96
90) Dibenzo(a,h)anthracene	28.55	278	695238	45.535	nq #	97
91) Benzo(g,h,i)perylene	29.64	276	702741	47.036	nq #	94

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

