

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012319\
 Data File : BG039260.D
 Acq On : 23 Jan 2019 12:54
 Operator : JU/SJ
 Sample : IDOC-S-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 IDOC-S-01

Manual Integrations
 APPROVED

Sohil
 1/24/2019 12:24:51 PM

Quant Time: Jan 23 13:43:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 04:20:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	15444	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	61107	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	43409	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	116147	20.00	ng	0.00
76) Chrysene-d12	21.67	240	122727	20.00	ng	-0.01
87) Perylene-d12	24.93	264	141879	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	120463	147.96	ng	0.00
7) Phenol-d6	7.17	99	172194	146.97	ng	0.00
23) Nitrobenzene-d5	9.19	82	84017	75.23	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	97391	133.84	ng	-0.01
45) 2-Fluorobiphenyl	13.27	172	245663	77.45	ng	0.00
79) Terphenyl-d14	20.00	244	509314	76.77	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.43	88	11985	37.592	ng	86
3) Pyridine	3.84	79	30725	35.467	ng	98
4) n-Nitrosodimethylamine	3.76	42	17355	44.520	ng	93
6) Aniline	7.33	93	38294	27.215	ng	97
8) 2-Chlorophenol	7.57	128	39615	41.279	ng	96
9) Benzaldehyde	7.15	77	10878	16.099	ng	87
10) Phenol	7.20	94	51909	44.190	ng	97
11) bis(2-Chloroethyl)ether	7.43	93	31843	35.468	ng	95
12) 1,3-Dichlorobenzene	7.89	146	41084	35.730	ng	96
13) 1,4-Dichlorobenzene	8.05	146	41526	35.659	ng	98
14) 1,2-Dichlorobenzene	8.36	146	40709	36.664	ng	96
15) Benzyl Alcohol	8.25	79	34713	39.342	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.53	45	62804	36.482	ng	99
17) 2-Methylphenol	8.45	107	34236	41.973	ng	97
18) Hexachloroethane	9.09	117	14372	36.327	ng	98
19) n-Nitroso-di-n-propylamine	8.81	70	26171	34.015	ng	97
20) 3+4-Methylphenols	8.78	107	45721	39.316	ng	98
22) Acetophenone	8.84	105	55394	34.712	ng	# 96
24) Nitrobenzene	9.23	77	40192	35.607	ng	99
25) Isophorone	9.74	82	76311	39.670	ng	100
26) 2-Nitrophenol	9.94	139	24398	39.961	ng	98
27) 2,4-Dimethylphenol	9.99	122	42182	49.109	ng	93
28) bis(2-Chloroethoxy)methane	10.23	93	48431	41.891	ng	98
29) 2,4-Dichlorophenol	10.47	162	49518	46.404	ng	97
30) 1,2,4-Trichlorobenzene	10.68	180	50018	39.011	ng	95
31) Naphthalene	10.87	128	124373	40.581	ng	98
32) Benzoic acid	10.16	122	12051	27.769	ng	91
33) 4-Chloroaniline	11.00	127	26503	19.326	ng	99
34) Hexachlorobutadiene	11.13	225	34126	38.682	ng	98
35) Caprolactam	11.78	113	16260m	37.999	ng	
36) 4-Chloro-3-methylphenol	12.11	107	50413	44.171	ng	90
37) 2-Methylnaphthalene	12.48	142	93923	40.875	ng	99
38) 1-Methylnaphthalene	12.70	142	88559	40.154	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.84	216	60655	37.206	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.80	237	68283	74.772	nq	99
43) 2,4,6-Trichlorophenol	13.09	196	40873	39.835	nq	95
44) 2,4,5-Trichlorophenol	13.16	196	45296	41.769	nq	98
46) 1,1'-Biphenyl	13.48	154	126844	36.877	nq	97
47) 2-Chloronaphthalene	13.53	162	103480	37.177	nq	99
48) 2-Nitroaniline	13.75	65	30681	39.345	nq	98
49) Acenaphthylene	14.37	152	170481	40.749	nq	99
50) Dimethylphthalate	14.10	163	140080	38.187	nq	99
51) 2,6-Dinitrotoluene	14.24	165	31094	37.914	nq	96
52) Acenaphthene	14.71	154	97308	38.712	nq	96
53) 3-Nitroaniline	14.57	138	20561	24.714	nq	96
54) 2,4-Dinitrophenol	14.79	184	29770	62.013	nq	98
55) Dibenzofuran	15.05	168	170534	38.413	nq	97
56) 4-Nitrophenol	14.88	139	46746	78.089	nq	97
57) 2,4-Dinitrotoluene	15.03	165	45452	38.629	nq	97
58) Fluorene	15.70	166	139734	41.816	nq	96
59) 2,3,4,6-Tetrachlorophenol	15.28	232	44392m	43.365	nq	
60) Diethylphthalate	15.45	149	145349	38.458	nq	98
61) 4-Chlorophenyl-phenylether	15.68	204	81071	38.509	nq	99
62) 4-Nitroaniline	15.74	138	33418	38.558	nq	94
63) Azobenzene	15.98	77	113216	38.306	nq	97
65) 4,6-Dinitro-2-methylphenol	15.79	198	28263	32.845	nq	96
66) n-Nitrosodiphenylamine	15.90	169	125437	36.431	nq	97
67) 4-Bromophenyl-phenylether	16.58	248	54461	36.477	nq	95
68) Hexachlorobenzene	16.69	284	57845	35.506	nq	98
69) Atrazine	16.85	200	53368	36.486	nq	97
70) Pentachlorophenol	17.05	266	52795	54.447	nq	94
71) Phenanthrene	17.44	178	243921	39.732	nq	99
72) Anthracene	17.53	178	254682	41.957	nq	99
73) Carbazole	17.81	167	216334	36.103	nq	99
74) Di-n-butylphthalate	18.34	149	260548	39.085	nq	98
75) Fluoranthene	19.45	202	312020	40.998	nq	99
77) Benzidine	19.64	184	100863	26.870	nq	98
78) Pyrene	19.81	202	305800	39.099	nq	98
80) Butylbenzylphthalate	20.68	149	118108	39.704	nq	95
81) Benzo(a)anthracene	21.65	228	304392	40.743	nq	99
82) 3,3'-Dichlorobenzidine	21.57	252	72154	23.704	nq	97
83) Chrysene	21.72	228	286754	40.356	nq	99
84) Bis(2-ethylhexyl)phthalate	21.52	149	172433	41.747	nq	96
85) Di-n-octyl phthalate	22.73	149	274756	39.340	nq	99
86) Indeno(1,2,3-cd)pyrene	28.67	276	356664	42.008	nq	99
88) Benzo(b)fluoranthene	23.89	252	298933	38.211	nq	99
89) Benzo(k)fluoranthene	23.96	252	292089	37.762	nq	97
90) Benzo(a)pyrene	24.78	252	342092	45.240	nq	# 97
91) Dibenzo(a,h)anthracene	28.72	278	296171	39.377	nq	99
92) Benzo(g,h,i)perylene	29.85	276	310049	41.639	nq	96

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

