

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010919\
 Data File : BG039044.D
 Acq On : 10 Jan 2019 7:35
 Operator : JU/SJ
 Sample : K1006-02MS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JC-03-010819-AMS

Manual Integrations
 APPROVED

Sohil
 1/10/2019 3:15:28 PM

Quant Time: Jan 10 09:48:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 09 14:23:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	21345	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	97512	20.00	ng	0.00
39) Acenaphthene-d10	14.67	164	77154	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	212399	20.00	ng	0.00
76) Chrysene-d12	21.70	240	171637	20.00	ng	0.00
87) Perylene-d12	24.99	264	192088	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	154755	137.53	ng	0.00
7) Phenol-d6	7.19	99	203635	125.75	ng	0.00
23) Nitrobenzene-d5	9.21	82	148039	83.07	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	179094	138.48	ng	0.00
45) 2-Fluorobiphenyl	13.29	172	433507	76.90	ng	0.00
79) Terphenyl-d14	20.03	244	847206	91.31	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.46	88	12351	28.030	ng	# 1
3) Pyridine	3.88	79	38590	32.231	ng	93
4) n-Nitrosodimethylamine	3.79	42	23851	44.269	ng	86
6) Aniline	7.35	93	65557	33.710	ng	97
8) 2-Chlorophenol	7.59	128	61746	46.552	ng	97
9) Benzaldehyde	7.17	77	20101	21.525	ng	98
10) Phenol	7.22	94	66854	41.179	ng	97
11) bis(2-Chloroethyl)ether	7.45	93	51617	41.599	ng	99
12) 1,3-Dichlorobenzene	7.92	146	53488	33.658	ng	98
13) 1,4-Dichlorobenzene	8.06	146	56119	34.868	ng	99
14) 1,2-Dichlorobenzene	8.38	146	55572	36.213	ng	97
15) Benzyl Alcohol	8.28	79	59041	48.415	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.55	45	99711	41.908	ng	98
17) 2-Methylphenol	8.48	107	55412	49.154	ng	95
18) Hexachloroethane	9.10	117	18460	33.761	ng	96
19) n-Nitroso-di-n-propylamine	8.83	70	47349	44.527	ng	96
20) 3+4-Methylphenols	8.80	107	75671	47.081	ng	98
22) Acetophenone	8.86	105	96925	38.062	ng	# 99
24) Nitrobenzene	9.25	77	71510	39.701	ng	99
25) Isophorone	9.76	82	142968	46.575	ng	98
26) 2-Nitrophenol	9.96	139	38923	39.951	ng	97
27) 2,4-Dimethylphenol	10.01	122	66003	48.153	ng	92
28) bis(2-Chloroethoxy)methane	10.24	93	79276	42.971	ng	97
29) 2,4-Dichlorophenol	10.49	162	79719	46.816	ng	99
30) 1,2,4-Trichlorobenzene	10.70	180	69316	33.879	ng	100
31) Naphthalene	10.90	128	189671	38.782	ng	100
32) Benzoic acid	10.17	122	18444	27.020	ng	97
33) 4-Chloroaniline	11.02	127	41133	18.796	ng	98
34) Hexachlorobutadiene	11.16	225	45081	32.022	ng	95
35) Caprolactam	11.81	113	20716m	30.338	ng	
36) 4-Chloro-3-methylphenol	12.13	107	83881	46.056	ng	92
37) 2-Methylnaphthalene	12.50	142	162848	44.412	ng	97
38) 1-Methylnaphthalene	12.72	142	154793	43.982	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.86	216	103827	35.832	ng	99

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41) Hexachlorocyclopentadiene	12.82	237	96065	60.381	nq	99
43) 2,4,6-Trichlorophenol	13.11	196	78334	42.954	nq	99
44) 2,4,5-Trichlorophenol	13.18	196	86552	44.905	nq	96
46) 1,1'-Biphenyl	13.51	154	229880	37.602	nq	100
47) 2-Chloronaphthalene	13.55	162	183900	37.172	nq	99
48) 2-Nitroaniline	13.77	65	59623	43.019	nq	96
49) Acenaphthylene	14.40	152	316336	42.541	nq	99
50) Dimethylphthalate	14.12	163	278773	42.758	nq	99
51) 2,6-Dinitrotoluene	14.26	165	62571	42.925	nq	97
52) Acenaphthene	14.74	154	182798	40.915	nq	97
53) 3-Nitroaniline	14.59	138	41722	28.215	nq	99
54) 2,4-Dinitrophenol	14.80	184	60101	69.706	nq	94
55) Dibenzofuran	15.07	168	321380	40.730	nq	99
56) 4-Nitrophenol	14.90	139	80972	76.103	nq	99
57) 2,4-Dinitrotoluene	15.04	165	91025	43.526	nq	95
58) Fluorene	15.72	166	266975	44.950	nq	94
59) 2,3,4,6-Tetrachlorophenol	15.30	232	82903	45.565	nq	98
60) Diethylphthalate	15.47	149	292896	43.602	nq	99
61) 4-Chlorophenyl-phenylether	15.70	204	152115	40.652	nq	100
62) 4-Nitroaniline	15.76	138	61486	39.914	nq	97
63) Azobenzene	16.00	77	218529	41.599	nq	96
65) 4,6-Dinitro-2-methylphenol	15.81	198	51866	32.961	nq	91
66) n-Nitrosodiphenylamine	15.93	169	249322	39.597	nq	96
67) 4-Bromophenyl-phenylether	16.60	248	106863	39.140	nq	94
68) Hexachlorobenzene	16.72	284	114457	38.418	nq	98
69) Atrazine	16.87	200	107533	40.201	nq	98
70) Pentachlorophenol	17.07	266	109284	60.865	nq	95
71) Phenanthrene	17.47	178	477413	42.525	nq	98
72) Anthracene	17.56	178	484492	43.647	nq	99
73) Carbazole	17.83	167	414831	37.857	nq	99
74) Di-n-butylphthalate	18.36	149	492139	40.371	nq	99
75) Fluoranthene	19.47	202	545254	39.177	nq	99
77) Benzidine	19.66	184	125693	23.943	nq	99
78) Pyrene	19.83	202	532515	48.684	nq	99
80) Butylbenzylphthalate	20.70	149	180379	43.358	nq	97
81) Benzo(a)anthracene	21.68	228	466636	44.661	nq	100
82) 3,3'-Dichlorobenzidine	21.60	252	118481	27.831	nq	97
83) Chrysene	21.75	228	433001	43.573	nq	99
84) Bis(2-ethylhexyl)phthalate	21.54	149	248660	43.047	nq	96
85) Di-n-octyl phthalate	22.76	149	429931	44.016	nq	100
86) Indeno(1,2,3-cd)pyrene	28.74	276	553971	46.653	nq	# 96
88) Benzo(b)fluoranthene	23.93	252	471864	44.550	nq	99
89) Benzo(k)fluoranthene	24.00	252	466413	44.538	nq	100
90) Benzo(a)pyrene	24.83	252	463200	45.245	nq	97
91) Dibenzo(a,h)anthracene	28.81	278	460172	45.190	nq	98
92) Benzo(g,h,i)perylene	29.93	276	461013	45.730	nq	98

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

