

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083118\
 Data File : BG036510.D
 Acq On : 31 Aug 2018 17:49
 Operator : JU/SJ
 Sample : J4270-06MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BU-02-083018MS

Quant Time: Sep 01 01:01:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 13:03:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	56321	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	233646	20.00	ng	0.00
38) Acenaphthene-d10	14.68	164	154999	20.00	ng	0.00
63) Phenanthrene-d10	17.43	188	410454	20.00	ng	0.00
75) Chrysene-d12	21.69	240	431409	20.00	ng	0.00
86) Perylene-d12	24.91	264	450055	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.63	112	382297	107.67	ng	0.00
7) Phenol-d6	7.22	99	511279	100.58	ng	0.00
23) Nitrobenzene-d5	9.22	82	344403	79.98	ng	0.00
41) 2,4,6-Tribromophenol	16.17	330	237769	128.56	ng	0.00
44) 2-Fluorobiphenyl	13.31	172	777947	71.66	ng	0.00
78) Terphenyl-d14	20.04	244	1412825	76.55	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.55	88	55335	33.395	ng	# 93
3) Pyridine	3.94	79	152324	32.751	ng	95
4) n-Nitrosodimethylamine	3.85	42	82293	39.725	ng	96
6) Aniline	7.39	93	141520	21.933	ng	97
8) 2-Chlorophenol	7.63	128	160889	41.786	ng	96
9) Benzaldehyde	7.20	77	95276	27.217	ng	97
10) Phenol	7.24	94	201289	37.838	ng	98
11) bis(2-Chloroethyl)ether	7.48	93	155430	36.854	ng	97
12) 1,3-Dichlorobenzene	7.95	146	167005	37.986	ng	99
13) 1,4-Dichlorobenzene	8.10	146	167137	37.654	ng	99
14) 1,2-Dichlorobenzene	8.41	146	160815	37.936	ng	99
15) Benzyl Alcohol	8.30	79	160236	39.102	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.58	45	295755	34.774	ng	98
17) 2-Methylphenol	8.50	107	145619	41.225	ng	97
18) Hexachloroethane	9.15	117	60408	37.958	ng	90
19) n-Nitroso-di-n-propylamine	8.87	70	132943	34.993	ng	97
20) 3+4-Methylphenols	8.83	107	200687	40.694	ng	96
22) Acetophenone	8.88	105	251650	41.016	ng	# 98
24) Nitrobenzene	9.27	77	196318	42.277	ng	98
25) Isophorone	9.78	82	378825	44.283	ng	99
26) 2-Nitrophenol	9.98	139	84970	46.177	ng	93
27) 2,4-Dimethylphenol	10.03	122	161349	47.361	ng	97
28) bis(2-Chloroethoxy)methane	10.27	93	219996	41.902	ng	98
29) 2,4-Dichlorophenol	10.51	162	173816	46.554	ng	94
30) 1,2,4-Trichlorobenzene	10.73	180	175466	40.173	ng	98
31) Naphthalene	10.92	128	513584	43.972	ng	99
32) Benzoic acid	10.15	122	74541	31.238	ng	98
33) 4-Chloroaniline	11.02	127	30814	5.757	ng	96
34) Hexachlorobutadiene	11.20	225	120082	40.919	ng	97
35) Caprolactam	11.79	113	58162	39.105	ng	88
36) 4-Chloro-3-methylphenol	12.13	107	192261	44.317	ng	97
37) 2-Methylnaphthalene	12.51	142	391922	45.860	ng	97
39) 1,2,4,5-Tetrachlorobenzene	12.88	216	224601	40.947	ng	99
40) Hexachlorocyclopentadiene	12.86	237	264463	92.665	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083118\
 Data File : BG036510.D
 Acq On : 31 Aug 2018 17:49
 Operator : JU/SJ
 Sample : J4270-06MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BU-02-083018MS

Quant Time: Sep 01 01:01:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 13:03:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	146959	43.982	ng	100
43) 2,4,5-Trichlorophenol	13.19	196	164434	46.139	ng	98
45) 1,1'-Biphenyl	13.52	154	514259	39.970	ng	98
46) 2-Chloronaphthalene	13.56	162	393706	40.083	ng	100
47) 2-Nitroaniline	13.76	65	140327	45.496	ng	98
48) Acenaphthylene	14.41	152	673489	43.874	ng	100
49) Dimethylphthalate	14.14	163	536420	42.383	ng	99
50) 2,6-Dinitrotoluene	14.25	165	118558	45.523	ng	95
51) Acenaphthene	14.75	154	423347	43.062	ng	99
52) 3-Nitroaniline	14.58	138	59013	21.027	ng	95
53) 2,4-Dinitrophenol	14.79	184	55426	56.186	ng	95
54) Dibenzofuran	15.08	168	645369	41.901	ng	98
55) 4-Nitrophenol	14.88	139	179680	78.302	ng	96
56) 2,4-Dinitrotoluene	15.04	165	169394	49.906	ng	90
57) Fluorene	15.73	166	516516	44.860	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.31	232	166291	49.662	ng	97
59) Diethylphthalate	15.50	149	535331	41.970	ng	99
60) 4-Chlorophenyl-phenylether	15.72	204	294909	41.479	ng	99
61) 4-Nitroaniline	15.75	138	131966	44.934	ng	97
62) Azobenzene	16.02	77	530117	40.847	ng	97
64) 4,6-Dinitro-2-methylphenol	15.80	198	59592	33.051	ng	97
65) n-Nitrosodiphenylamine	15.93	169	485801	39.833	ng	97
66) 4-Bromophenyl-phenylether	16.62	248	199984	41.615	ng	99
67) Hexachlorobenzene	16.73	284	197550	40.202	ng	98
68) Atrazine	16.88	200	203359	44.423	ng	98
69) Pentachlorophenol	17.07	266	213988	64.075	ng	97
70) Phenanthrene	17.47	178	907022	43.489	ng	99
71) Anthracene	17.56	178	917774	44.145	ng	99
72) Carbazole	17.83	167	833169	40.128	ng	98
73) Di-n-butylphthalate	18.39	149	954026	41.263	ng	100
74) Fluoranthene	19.47	202	1142880	44.945	ng	99
76) Benzidine	19.65	184	190737	13.858	ng	98
77) Pyrene	19.84	202	1126081	42.447	ng	98
79) Butylbenzylphthalate	20.72	149	441408	41.809	ng	95
80) Benzo(a)anthracene	21.67	228	1145125	43.815	ng	99
81) 3,3'-Dichlorobenzidine	21.59	252	182656	17.536	ng	99
82) Chrysene	21.74	228	1050052	42.387	ng	99
83) Bis(2-ethylhexyl)phthalate	21.59	149	610035	41.291	ng	100
84) Di-n-octyl phthalate	22.80	149	1039209	42.112	ng	97
85) Indeno(1,2,3-cd)pyrene	28.56	276	1244215	43.242	ng	97
87) Benzo(b)fluoranthene	23.89	252	1139895	45.611	ng	98
88) Benzo(k)fluoranthene	23.95	252	1119373	45.847	ng	98
89) Benzo(a)pyrene	24.75	252	1108141	46.143	ng	100
90) Dibenzo(a,h)anthracene	28.63	278	981052	42.316	ng	97
91) Benzo(g,h,i)perylene	29.69	276	995191	42.715	ng	95

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

