

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091418\
 Data File : BG036710.D
 Acq On : 14 Sep 2018 18:03
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 9/17/2018 2:36:05 PM

Quant Time: Sep 15 02:21:33 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	8.05	152	59847	20.00	ng	-0.01
21) Naphthalene-d8	10.86	136	268816	20.00	ng	-0.01
38) Acenaphthene-d10	14.67	164	179775	20.00	ng	-0.01
63) Phenanthrene-d10	17.42	188	444511	20.00	ng	-0.01
75) Chrysene-d12	21.69	240	423976	20.00	ng	0.00
86) Perylene-d12	24.89	264	444309	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	290684	77.05	ng	0.00
7) Phenol-d6	7.21	99	421957	78.12	ng	-0.01
23) Nitrobenzene-d5	9.21	82	424894	85.76	ng	-0.01
41) 2,4,6-Tribromophenol	16.17	330	196554	91.63	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	986926	78.38	ng	-0.01
78) Terphenyl-d14	20.03	244	1450562	79.97	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.53	88	59664	33.887	ng	99
3) Pyridine	3.93	79	189713	38.386	ng	98
4) n-Nitrosodimethylamine	3.84	42	78392	35.612	ng	97
6) Aniline	7.38	93	267472	39.011	ng	97
8) 2-Chlorophenol	7.61	128	158685	38.786	ng	97
9) Benzaldehyde	7.19	77	133879	35.992	ng	98
10) Phenol	7.24	94	219339	38.802	ng	98
11) bis(2-Chloroethyl)ether	7.47	93	171063	38.171	ng	98
12) 1,3-Dichlorobenzene	7.94	146	177305	37.953	ng	98
13) 1,4-Dichlorobenzene	8.09	146	182454	38.683	ng	98
14) 1,2-Dichlorobenzene	8.41	146	172764	38.354	ng	99
15) Benzyl Alcohol	8.29	79	169552	38.937	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.58	45	326061	36.078	ng	98
17) 2-Methylphenol	8.49	107	146007	38.899	ng	99
18) Hexachloroethane	9.13	117	67543	39.941	ng	95
19) n-Nitroso-di-n-propylamine	8.86	70	158818	39.341	ng	92
20) 3+4-Methylphenols	8.82	107	213195	40.684	ng	96
22) Acetophenone	8.87	105	266383	37.737	ng	# 98
24) Nitrobenzene	9.26	77	218012	40.806	ng	99
25) Isophorone	9.77	82	389525	39.576	ng	99
26) 2-Nitrophenol	9.97	139	105760	49.955	ng	95
27) 2,4-Dimethylphenol	10.02	122	145282	37.066	ng	98
28) bis(2-Chloroethoxy)methane	10.26	93	233643	38.679	ng	96
29) 2,4-Dichlorophenol	10.50	162	182083	42.388	ng	94
30) 1,2,4-Trichlorobenzene	10.72	180	200479	39.895	ng	98
31) Naphthalene	10.91	128	519340	38.648	ng	99
32) Benzoic acid	10.09	122	23694m	11.872	ng	
33) 4-Chloroaniline	11.01	127	247120	40.131	ng	99
34) Hexachlorobutadiene	11.19	225	139300	41.258	ng	95
35) Caprolactam	11.79	113	72791	42.537	ng	100
36) 4-Chloro-3-methylphenol	12.12	107	209206	41.914	ng	99
37) 2-Methylnaphthalene	12.51	142	394164	40.088	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	252617	39.707	ng	99
40) Hexachlorocyclopentadiene	12.85	237	142984	43.195	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	169047	43.620	nq	99
43) 2,4,5-Trichlorophenol	13.18	196	177691	42.987	nq	98
45) 1,1'-Biphenyl	13.51	154	569014	38.131	nq	99
46) 2-Chloronaphthalene	13.56	162	440285	38.648	nq	100
47) 2-Nitroaniline	13.75	65	157268	43.962	nq	97
48) Acenaphthylene	14.40	152	683895	38.412	nq	99
49) Dimethylphthalate	14.13	163	584103	39.790	nq	99
50) 2,6-Dinitrotoluene	14.25	165	129133	42.750	nq	95
51) Acenaphthene	14.74	154	420715	36.896	nq	97
52) 3-Nitroaniline	14.58	138	145277	44.630	nq	95
53) 2,4-Dinitrophenol	14.78	184	43261	37.810	nq	96
54) Dibenzofuran	15.07	168	688367	38.533	nq	99
55) 4-Nitrophenol	14.87	139	118166	44.398	nq	99
56) 2,4-Dinitrotoluene	15.03	165	182339	46.316	nq	91
57) Fluorene	15.72	166	513563	38.456	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.30	232	172315	44.369	nq	96
59) Diethylphthalate	15.49	149	572160	38.675	nq	99
60) 4-Chlorophenyl-phenylether	15.71	204	332394	40.308	nq	98
61) 4-Nitroaniline	15.74	138	144547	42.435	nq	96
62) Azobenzene	16.01	77	558611	37.111	nq	97
64) 4,6-Dinitro-2-methylphenol	15.80	198	93028	47.642	nq	99
65) n-Nitrosodiphenylamine	15.93	169	515024	38.993	nq	97
66) 4-Bromophenyl-phenylether	16.61	248	214243	41.166	nq	100
67) Hexachlorobenzene	16.73	284	218746	41.105	nq	97
68) Atrazine	16.87	200	169951	34.281	nq	99
69) Pentachlorophenol	17.07	266	132630	36.671	nq	97
70) Phenanthrene	17.47	178	863907	38.248	nq	99
71) Anthracene	17.55	178	862244	38.296	nq	99
72) Carbazole	17.82	167	851431	37.866	nq	99
73) Di-n-butylphthalate	18.38	149	925097	36.946	nq	99
74) Fluoranthene	19.47	202	1022424	37.128	nq	100
76) Benzidine	19.65	184	552548	40.849	nq	99
77) Pyrene	19.83	202	1024502	39.295	nq	98
79) Butylbenzylphthalate	20.71	149	417698	40.257	nq	97
80) Benzo(a)anthracene	21.67	228	991645	38.608	nq	99
81) 3,3'-Dichlorobenzidine	21.58	252	401429	39.215	nq	98
82) Chrysene	21.73	228	948648	38.965	nq	99
83) Bis(2-ethylhexyl)phthalate	21.57	149	563833	38.833	nq	97
84) Di-n-octyl phthalate	22.78	149	945900	39.003	nq	99
85) Indeno(1,2,3-cd)pyrene	28.54	276	1104619	39.063	nq	# 95
87) Benzo(b)fluoranthene	23.88	252	940568	38.122	nq	97
88) Benzo(k)fluoranthene	23.95	252	952345	39.510	nq	99
89) Benzo(a)pyrene	24.74	252	934896	39.433	nq	99
90) Dibenzo(a,h)anthracene	28.61	278	901561	39.390	nq	98
91) Benzo(g,h,i)perylene	29.67	276	909607	39.547	nq	97

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

