

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080918\
 Data File : BG036253.D
 Acq On : 10 Aug 2018 4:46
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 8/10/2018 4:01:53 PM

Quant Time: Aug 10 13:18:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 12:30:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	39881	20.00	ng	-0.02
21) Naphthalene-d8	10.81	136	158315	20.00	ng	-0.03
38) Acenaphthene-d10	14.64	164	111502	20.00	ng	-0.01
63) Phenanthrene-d10	17.38	188	305696	20.00	ng	-0.01
75) Chrysene-d12	21.64	240	321745	20.00	ng	-0.01
86) Perylene-d12	24.83	264	315030	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	178181	74.18	ng	0.00
7) Phenol-d6	7.19	99	266121	74.25	ng	0.00
23) Nitrobenzene-d5	9.17	82	290529	76.66	ng	-0.01
41) 2,4,6-Tribromophenol	16.13	330	130876	89.05	ng	-0.01
44) 2-Fluorobiphenyl	13.26	172	664633	79.86	ng	-0.02
78) Terphenyl-d14	19.99	244	1150083	78.79	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	39569m	32.365	ng	
3) Pyridine	3.92	79	110741	34.696	ng	# 67
4) n-Nitrosodimethylamine	3.83	42	46223	33.728	ng	# 81
6) Aniline	7.34	93	160151	34.476	ng	98
8) 2-Chlorophenol	7.58	128	93643	38.330	ng	97
9) Benzaldehyde	7.15	77	79603	36.199	ng	98
10) Phenol	7.21	94	136762	37.771	ng	91
11) bis(2-Chloroethyl)ether	7.43	93	103203	36.395	ng	88
12) 1,3-Dichlorobenzene	7.90	146	114429	38.786	ng	# 93
13) 1,4-Dichlorobenzene	8.04	146	117966	39.353	ng	96
14) 1,2-Dichlorobenzene	8.36	146	112328	39.007	ng	95
15) Benzyl Alcohol	8.25	79	118999	36.564	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.53	45	76995	32.387	ng	71
17) 2-Methylphenol	8.46	107	93484	37.974	ng	# 87
18) Hexachloroethane	9.09	117	44595	38.909	ng	# 85
19) n-Nitroso-di-n-propylamine	8.82	70	101684	33.693	ng	# 85
20) 3+4-Methylphenols	8.79	107	129005	37.774	ng	91
22) Acetophenone	8.83	105	180098	36.582	ng	# 94
24) Nitrobenzene	9.21	77	142693	37.440	ng	96
25) Isophorone	9.74	82	252041	35.827	ng	97
26) 2-Nitrophenol	9.93	139	60871	44.970	ng	# 85
27) 2,4-Dimethylphenol	9.98	122	89120	38.896	ng	87
28) bis(2-Chloroethoxy)methane	10.21	93	139701	36.038	ng	95
29) 2,4-Dichlorophenol	10.47	162	110223	42.142	ng	94
30) 1,2,4-Trichlorobenzene	10.67	180	133116	41.506	ng	97
31) Naphthalene	10.86	128	315647	38.275	ng	98
32) Benzoic acid	10.17	122	36548	23.695	ng	94
33) 4-Chloroaniline	10.97	127	136844	38.072	ng	97
34) Hexachlorobutadiene	11.14	225	111650	44.644	ng	97
35) Caprolactam	11.78	113	36610	32.617	ng	# 65
36) 4-Chloro-3-methylphenol	12.10	107	135291	38.590	ng	93
37) 2-Methylnaphthalene	12.46	142	238742	38.858	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	12.83	216	175728	41.756	ng	98
40) Hexachlorocyclopentadiene	12.80	237	83436	37.803	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.07	196	103607	42.097	nq	96
43) 2,4,5-Trichlorophenol	13.16	196	112690	40.930	nq	99
45) 1,1'-Biphenyl	13.47	154	335587	38.820	nq	98
46) 2-Chloronaphthalene	13.51	162	264786	39.378	nq	98
47) 2-Nitroaniline	13.72	65	92241	38.802	nq	97
48) Acenaphthylene	14.35	152	437217	38.816	nq	96
49) Dimethylphthalate	14.09	163	371240	38.001	nq	100
50) 2,6-Dinitrotoluene	14.21	165	82049	41.244	nq	93
51) Acenaphthene	14.70	154	265927	37.900	nq	99
52) 3-Nitroaniline	14.55	138	75748	37.254	nq #	93
53) 2,4-Dinitrophenol	14.75	184	36255	39.106	nq #	71
54) Dibenzofuran	15.03	168	432523	38.760	nq	93
55) 4-Nitrophenol	14.87	139	48792	33.695	nq #	87
56) 2,4-Dinitrotoluene	15.00	165	118358	41.470	nq #	88
57) Fluorene	15.68	166	359805	38.470	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.27	232	106235	40.243	nq #	93
59) Diethylphthalate	15.45	149	370031	36.836	nq	98
60) 4-Chlorophenyl-phenylether	15.67	204	220620	40.133	nq	97
61) 4-Nitroaniline	15.71	138	72380	34.031	nq #	73
62) Azobenzene	15.96	77	348254	35.147	nq	95
64) 4,6-Dinitro-2-methylphenol	15.77	198	69244	40.691	nq	87
65) n-Nitrosodiphenylamine	15.89	169	331783	38.131	nq	98
66) 4-Bromophenyl-phenylether	16.56	248	144199	40.658	nq	97
67) Hexachlorobenzene	16.69	284	149499	40.933	nq	94
68) Atrazine	16.83	200	123013	34.337	nq	95
69) Pentachlorophenol	17.04	266	71912m	32.326	nq	
70) Phenanthrene	17.42	178	578011	37.893	nq	98
71) Anthracene	17.52	178	569496	37.495	nq	98
72) Carbazole	17.79	167	526598	36.508	nq	99
73) Di-n-butylphthalate	18.33	149	626483	36.754	nq #	97
74) Fluoranthene	19.43	202	760350	37.006	nq	97
76) Benzidine	19.61	184	305815	36.154	nq	98
77) Pyrene	19.79	202	764563	37.581	nq	97
79) Butylbenzylphthalate	20.67	149	290198	39.889	nq #	96
80) Benzo(a)anthracene	21.62	228	756205	37.715	nq	99
81) 3,3'-Dichlorobenzidine	21.54	252	271879	38.889	nq #	96
82) Chrysene	21.69	228	712357	38.340	nq	99
83) Bis(2-ethylhexyl)phthalate	21.52	149	402267	40.671	nq	99
84) Di-n-octyl phthalate	22.71	149	667916	40.332	nq #	93
85) Indeno(1,2,3-cd)pyrene	28.47	276	775950	37.721	nq #	84
87) Benzo(b)fluoranthene	23.82	252	711492	37.159	nq #	96
88) Benzo(k)fluoranthene	23.89	252	726104	38.789	nq #	97
89) Benzo(a)pyrene	24.68	252	685820	38.500	nq #	95
90) Dibenzo(a,h)anthracene	28.52	278	657057	37.572	nq #	93
91) Benzo(g,h,i)perylene	29.60	276	648585	37.900	nq #	93

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

