

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062518\
 Data File : BG035371.D
 Acq On : 25 Jun 2018 10:28
 Operator : JU/SJ
 Sample : SSTDC0040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDC0040

Manual Integrations
 APPROVED

Sohil
 6/27/2018 10:48:19 AM

Quant Time: Jun 25 11:09:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 08 17:16:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	48872	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	197487	20.00	ng	-0.02
38) Acenaphthene-d10	14.67	164	134487	20.00	ng	-0.02
63) Phenanthrene-d10	17.42	188	354427	20.00	ng	-0.02
75) Chrysene-d12	21.69	240	397590	20.00	ng	-0.03
86) Perylene-d12	24.91	264	391056	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	226678	78.27	ng	-0.02
7) Phenol-d6	7.21	99	335859	78.58	ng	-0.01
23) Nitrobenzene-d5	9.21	82	334445	80.50	ng	-0.02
41) 2,4,6-Tribromophenol	16.16	330	147293	85.56	ng	-0.02
44) 2-Fluorobiphenyl	13.30	172	770945	74.52	ng	-0.03
78) Terphenyl-d14	20.02	244	1316623	69.37	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	53483	38.708	ng	# 69
3) Pyridine	3.95	79	144710	38.816	ng	# 73
4) n-Nitrosodimethylamine	3.86	42	55750	34.809	ng	# 83
6) Aniline	7.38	93	208035	37.654	ng	95
8) 2-Chlorophenol	7.62	128	114709	37.789	ng	95
9) Benzaldehyde	7.19	77	115862	35.192	ng	94
10) Phenol	7.24	94	168153	38.015	ng	97
11) bis(2-Chloroethyl)ether	7.47	93	126963	36.981	ng	88
12) 1,3-Dichlorobenzene	7.95	146	138444	38.225	ng	96
13) 1,4-Dichlorobenzene	8.09	146	142210	38.036	ng	97
14) 1,2-Dichlorobenzene	8.41	146	134689	38.352	ng	99
15) Benzyl Alcohol	8.29	79	140228	34.624	ng	91
16) 2,2'-oxybis(1-Chloropropan	8.57	45	112704m	32.982	ng	
17) 2-Methylphenol	8.49	107	112938	38.727	ng	# 92
18) Hexachloroethane	9.14	117	52894	39.516	ng	93
19) n-Nitroso-di-n-propylamine	8.86	70	122410	33.710	ng	# 85
20) 3+4-Methylphenols	8.82	107	158023	37.804	ng	95
22) Acetophenone	8.87	105	222224	36.326	ng	# 88
24) Nitrobenzene	9.26	77	167594	39.393	ng	97
25) Isophorone	9.78	82	297766	33.946	ng	96
26) 2-Nitrophenol	9.97	139	66892	45.615	ng	# 91
27) 2,4-Dimethylphenol	10.02	122	110743	35.928	ng	98
28) bis(2-Chloroethoxy)methane	10.26	93	169207	35.896	ng	99
29) 2,4-Dichlorophenol	10.50	162	132188	39.413	ng	95
30) 1,2,4-Trichlorobenzene	10.72	180	157258	39.103	ng	98
31) Naphthalene	10.91	128	375747	36.625	ng	98
32) Benzoic acid	10.16	122	82606	40.011	ng	95
33) 4-Chloroaniline	11.01	127	164811	36.997	ng	94
34) Hexachlorobutadiene	11.19	225	122311	39.106	ng	95
35) Caprolactam	11.78	113	46024	37.130	ng	# 66
36) 4-Chloro-3-methylphenol	12.13	107	162663	38.907	ng	92
37) 2-Methylnaphthalene	12.51	142	287589	36.974	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	200182	39.962	ng	99
40) Hexachlorocyclopentadiene	12.85	237	112806	35.910	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	127057	42.359	nq	99
43) 2,4,5-Trichlorophenol	13.18	196	138423	42.046	nq	96
45) 1,1'-Biphenyl	13.51	154	403838	37.344	nq	96
46) 2-Chloronaphthalene	13.56	162	309574	37.863	nq	98
47) 2-Nitroaniline	13.75	65	103379	42.819	nq	99
48) Acenaphthylene	14.40	152	508952	37.124	nq	98
49) Dimethylphthalate	14.13	163	425664	36.296	nq	99
50) 2,6-Dinitrotoluene	14.25	165	93441	43.674	nq #	86
51) Acenaphthene	14.74	154	347415	38.785	nq	97
52) 3-Nitroaniline	14.58	138	94516	45.020	nq #	93
53) 2,4-Dinitrophenol	14.78	184	40883	47.216	nq #	67
54) Dibenzofuran	15.07	168	504532	37.608	nq	94
55) 4-Nitrophenol	14.88	139	67219	37.937	nq #	89
56) 2,4-Dinitrotoluene	15.03	165	134176	47.210	nq #	87
57) Fluorene	15.72	166	422373	37.672	nq	96
58) 2,3,4,6-Tetrachlorophenol	15.29	232	134730	42.126	nq #	93
59) Diethylphthalate	15.49	149	435592	36.405	nq	97
60) 4-Chlorophenyl-phenylether	15.71	204	251798	39.385	nq	95
61) 4-Nitroaniline	15.74	138	98458	43.730	nq #	84
62) Azobenzene	16.01	77	398795	34.012	nq	93
64) 4,6-Dinitro-2-methylphenol	15.79	198	76798	47.459	nq	85
65) n-Nitrosodiphenylamine	15.92	169	389909	36.632	nq	98
66) 4-Bromophenyl-phenylether	16.60	248	164918	38.639	nq	98
67) Hexachlorobenzene	16.72	284	166811	38.398	nq	94
68) Atrazine	16.87	200	165080	36.369	nq	94
69) Pentachlorophenol	17.06	266	115438	39.517	nq	98
70) Phenanthrene	17.46	178	680142	37.777	nq	99
71) Anthracene	17.55	178	673389	37.339	nq	98
72) Carbazole	17.82	167	630881	37.703	nq	99
73) Di-n-butylphthalate	18.37	149	703610	35.279	nq #	98
74) Fluoranthene	19.47	202	896558	38.268	nq	96
76) Benzidine	19.64	184	355631	24.042	nq #	96
77) Pyrene	19.83	202	912660	34.820	nq	99
79) Butylbenzylphthalate	20.71	149	317872	33.398	nq #	96
80) Benzo(a)anthracene	21.66	228	906969	35.697	nq	98
81) 3,3'-Dichlorobenzidine	21.58	252	338709	35.434	nq #	99
82) Chrysene	21.73	228	845165	36.332	nq	98
83) Bis(2-ethylhexyl)phthalate	21.57	149	443540	32.981	nq	99
84) Di-n-octyl phthalate	22.78	149	753803	32.672	nq #	92
85) Indeno(1,2,3-cd)pyrene	28.56	276	998161	36.637	nq #	88
87) Benzo(b)fluoranthene	23.88	252	878769	36.491	nq #	95
88) Benzo(k)fluoranthene	23.95	252	873433	37.077	nq #	96
89) Benzo(a)pyrene	24.75	252	827154	36.502	nq #	97
90) Dibenzo(a,h)anthracene	28.64	278	826413	36.943	nq #	96
91) Benzo(g,h,i)perylene	29.72	276	812960	37.135	nq #	91

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

