

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121719\
 Data File : BG043815.D
 Acq On : 17 Dec 2019 11:43
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 12/18/2019 2:40:33 PM

Quant Time: Dec 18 07:08:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 09 18:16:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	149920	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	668890	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	473864	20.00	ng	0.00
64) Phenanthrene-d10	17.37	188	1042792	20.00	ng	0.00
76) Chrysene-d12	21.63	240	846473	20.00	ng	0.00
87) Perylene-d12	24.79	264	962340	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.59	112	632134	79.89	ng	0.00
7) Phenol-d6	7.17	99	922491	80.84	ng	0.01
23) Nitrobenzene-d5	9.16	82	940235	78.92	ng	0.00
42) 2,4,6-Tribromophenol	16.12	330	386603	85.58	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	2049685	75.98	ng	0.00
79) Terphenyl-d14	19.99	244	2689246	76.43	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	138797	37.880	ng	100
3) Pyridine	3.89	79	430066	42.475	ng	97
4) n-Nitrosodimethylamine	3.80	42	178874	36.846	ng	94
6) Aniline	7.32	93	601231	39.819	ng	97
8) 2-Chlorophenol	7.57	128	368736	39.989	ng	99
9) Benzaldehyde	7.14	77	263406	42.722	ng	89
10) Phenol	7.20	94	473626	40.202	ng	97
11) bis(2-Chloroethyl)ether	7.42	93	388384	38.851	ng	96
12) 1,3-Dichlorobenzene	7.89	146	430152	38.927	ng	99
13) 1,4-Dichlorobenzene	8.04	146	425465	38.520	ng	97
14) 1,2-Dichlorobenzene	8.36	146	413884	38.891	ng	98
15) Benzyl Alcohol	8.24	79	373726	41.154	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.53	45	634799	39.621	ng	98
17) 2-Methylphenol	8.45	107	343971	40.901	ng	98
18) Hexachloroethane	9.09	117	163402	40.712	ng	98
19) n-Nitroso-di-n-propylamine	8.81	70	345895	39.649	ng	100
20) 3+4-Methylphenols	8.78	107	493861	42.174	ng	97
22) Acetophenone	8.82	105	632120	37.893	ng	100
24) Nitrobenzene	9.20	77	463461	38.193	ng	99
25) Isophorone	9.73	82	912467	38.351	ng	99
26) 2-Nitrophenol	9.92	139	228615	48.152	ng	95
27) 2,4-Dimethylphenol	9.98	122	343869	40.900	ng	98
28) bis(2-Chloroethoxy)methane	10.21	93	584946	38.409	ng	97
29) 2,4-Dichlorophenol	10.45	162	420422	40.720	ng	97
30) 1,2,4-Trichlorobenzene	10.67	180	467798	38.275	ng	97
31) Naphthalene	10.86	128	1238509	38.588	ng	99
32) Benzoic acid	10.12	122	270067	46.333	ng	95
33) 4-Chloroaniline	10.96	127	611368	39.287	ng	99
34) Hexachlorobutadiene	11.16	225	283423	38.272	ng	97
35) Caprolactam	11.74	113	166021	40.973	ng	94
36) 4-Chloro-3-methylphenol	12.08	107	455707	40.998	ng	99
37) 2-Methylnaphthalene	12.46	142	972395	39.089	ng	98
38) 1-Methylnaphthalene	12.68	142	919450	38.869	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	537624	37.757	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	278554	41.678	nq	99
43) 2,4,6-Trichlorophenol	13.06	196	381611	42.165	nq	99
44) 2,4,5-Trichlorophenol	13.14	196	393422	41.492	nq	98
46) 1,1'-Biphenyl	13.46	154	1249433	38.246	nq	97
47) 2-Chloronaphthalene	13.51	162	1018868	38.150	nq	99
48) 2-Nitroaniline	13.70	65	336658	42.743	nq	97
49) Acenaphthylene	14.35	152	1519430	38.192	nq	98
50) Dimethylphthalate	14.09	163	1290675	38.161	nq	99
51) 2,6-Dinitrotoluene	14.20	165	301737	41.398	nq	99
52) Acenaphthene	14.70	154	1027640	39.265	nq	98
53) 3-Nitroaniline	14.53	138	343809	41.220	nq	94
54) 2,4-Dinitrophenol	14.73	184	142296	55.078	nq	94
55) Dibenzofuran	15.03	168	1474718	38.066	nq	99
56) 4-Nitrophenol	14.83	139	265947	43.705	nq	96
57) 2,4-Dinitrotoluene	14.98	165	421203	42.768	nq	96
58) Fluorene	15.68	166	1207784	38.678	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.26	232	350722m	41.225	nq	
60) Diethylphthalate	15.46	149	1296077	38.364	nq	99
61) 4-Chlorophenyl-phenylether	15.67	204	672128	38.433	nq	99
62) 4-Nitroaniline	15.69	138	347866	40.252	nq	98
63) Azobenzene	15.97	77	1182159	38.089	nq	98
65) 4,6-Dinitro-2-methylphenol	15.75	198	197467	48.868	nq	99
66) n-Nitrosodiphenylamine	15.88	169	1116921	38.537	nq	98
67) 4-Bromophenyl-phenylether	16.57	248	427433	38.339	nq	95
68) Hexachlorobenzene	16.68	284	448965	37.969	nq	98
69) Atrazine	16.84	200	325706	35.251	nq	98
70) Pentachlorophenol	17.02	266	266120	42.309	nq	96
71) Phenanthrene	17.42	178	1876927	38.020	nq	99
72) Anthracene	17.51	178	1875640	38.382	nq	97
73) Carbazole	17.77	167	1736289	38.010	nq	98
74) Di-n-butylphthalate	18.35	149	2048299	39.114	nq	98
75) Fluoranthene	19.42	202	2104186	37.242	nq	97
77) Benzidine	19.60	184	815042	44.135	nq	97
78) Pyrene	19.78	202	2132016	38.541	nq	98
80) Butylbenzylphthalate	20.68	149	998990	43.361	nq	96
81) Benzo(a)anthracene	21.61	228	2063929	38.633	nq	97
82) 3,3'-Dichlorobenzidine	21.53	252	773056	39.170	nq	99
83) Chrysene	21.68	228	1974509	38.792	nq	97
84) Bis(2-ethylhexyl)phthalate	21.54	149	1337579	40.476	nq	99
85) Di-n-octyl phthalate	22.75	149	2295711	42.494	nq	99
86) Indeno(1,2,3-cd)pyrene	28.38	276	2093423	32.134	nq	# 94
88) Benzo(b)fluoranthene	23.79	252	2100021	37.901	nq	99
89) Benzo(k)fluoranthene	23.86	252	2056803	38.428	nq	98
90) Benzo(a)pyrene	24.65	252	2021613	38.286	nq	98
91) Dibenzo(a,h)anthracene	28.45	278	1829760	34.693	nq	98
92) Benzo(g,h,i)perylene	29.50	276	1858897m	35.716	nq	

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

