

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
 Data File : BG044019.D
 Acq On : 13 Jan 2020 19:40
 Operator : JU
 Sample : PB126048BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB126048BS

Manual Integrations
 APPROVED

mohammad
 1/14/2020 12:24:53 PM

Quant Time: Jan 14 10:27:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	112555	20.00	ng	-0.02
21) Naphthalene-d8	10.75	136	478390	20.00	ng	-0.02
39) Acenaphthene-d10	14.58	164	316335	20.00	ng	-0.02
64) Phenanthrene-d10	17.32	188	706350	20.00	ng	-0.02
76) Chrysene-d12	21.58	240	625803	20.00	ng	0.00
87) Perylene-d12	24.72	264	583242	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	744105	121.39	ng	0.00
7) Phenol-d6	7.12	99	968527	113.03	ng	0.00
23) Nitrobenzene-d5	9.10	82	596063	66.65	ng	-0.02
42) 2,4,6-Tribromophenol	16.07	330	392776	118.65	ng	-0.02
45) 2-Fluorobiphenyl	13.20	172	1317333	75.45	ng	-0.02
79) Terphenyl-d14	19.94	244	2014705	75.22	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.41	88	101326	37.910	ng	97
3) Pyridine	3.81	79	243452	30.832	ng	94
4) n-Nitrosodimethylamine	3.72	42	127697	36.325	ng	97
6) Aniline	7.27	93	310637	27.601	ng	96
8) 2-Chlorophenol	7.51	128	307707	42.758	ng	95
9) Benzaldehyde	7.08	77	126257m	23.022	ng	
10) Phenol	7.14	94	373263	42.279	ng	98
11) bis(2-Chloroethyl)ether	7.37	93	283130	37.153	ng	99
12) 1,3-Dichlorobenzene	7.84	146	325071	38.600	ng	99
13) 1,4-Dichlorobenzene	7.98	146	325876	39.218	ng	98
14) 1,2-Dichlorobenzene	8.30	146	305837	38.347	ng	99
15) Benzyl Alcohol	8.18	79	252966	37.407	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.48	45	382922	32.478	ng	99
17) 2-Methylphenol	8.39	107	263282	41.210	ng	98
18) Hexachloroethane	9.03	117	121093	37.453	ng	96
19) n-Nitroso-di-n-propylamine	8.75	70	220188	34.506	ng	96
20) 3+4-Methylphenols	8.72	107	358268	40.198	ng	98
22) Acetophenone	8.77	105	417007	35.063	ng	# 98
24) Nitrobenzene	9.14	77	315502	35.622	ng	97
25) Isophorone	9.67	82	604154	36.010	ng	99
26) 2-Nitrophenol	9.86	139	172918	41.266	ng	93
27) 2,4-Dimethylphenol	9.92	122	285646	45.285	ng	98
28) bis(2-Chloroethoxy)methane	10.15	93	382584	34.205	ng	99
29) 2,4-Dichlorophenol	10.39	162	303366	40.320	ng	96
30) 1,2,4-Trichlorobenzene	10.61	180	305512	36.214	ng	98
31) Naphthalene	10.80	128	887656	38.344	ng	99
32) Benzoic acid	10.06	122	164617	34.128	ng	90
33) 4-Chloroaniline	10.90	127	215514	19.518	ng	98
34) Hexachlorobutadiene	11.09	225	176781	34.321	ng	99
35) Caprolactam	11.67	113	114316m	40.813	ng	
36) 4-Chloro-3-methylphenol	12.03	107	321828	40.179	ng	96
37) 2-Methylnaphthalene	12.40	142	668023	38.329	ng	96
38) 1-Methylnaphthalene	12.63	142	629956	38.137	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.78	216	314685	33.940	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.76	237	331670	68.999	nq	98
43) 2,4,6-Trichlorophenol	13.01	196	242454	38.004	nq	99
44) 2,4,5-Trichlorophenol	13.08	196	263837	40.097	nq	99
46) 1,1'-Biphenyl	13.41	154	827251	36.474	nq	98
47) 2-Chloronaphthalene	13.45	162	654309	35.701	nq	98
48) 2-Nitroaniline	13.65	65	204065	34.614	nq	96
49) Acenaphthylene	14.30	152	1027439	39.004	nq	98
50) Dimethylphthalate	14.04	163	851251	37.899	nq	99
51) 2,6-Dinitrotoluene	14.15	165	196033	38.614	nq	95
52) Acenaphthene	14.65	154	649564	35.162	nq	97
53) 3-Nitroaniline	14.48	138	152025	26.370	nq	100
54) 2,4-Dinitrophenol	14.68	184	238314	91.311	nq	98
55) Dibenzofuran	14.98	168	1003248	39.450	nq	98
56) 4-Nitrophenol	14.79	139	387878	87.799	nq	92
57) 2,4-Dinitrotoluene	14.94	165	280084	40.546	nq	98
58) Fluorene	15.63	166	789159	39.152	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.21	232	236810	41.854	nq	96
60) Diethylphthalate	15.40	149	870495	38.748	nq	99
61) 4-Chlorophenyl-phenylether	15.62	204	411515	37.599	nq	95
62) 4-Nitroaniline	15.64	138	205233	36.049	nq	94
63) Azobenzene	15.91	77	783897	37.625	nq	99
65) 4,6-Dinitro-2-methylphenol	15.70	198	171219	47.134	nq	97
66) n-Nitrosodiphenylamine	15.83	169	757637	36.423	nq	99
67) 4-Bromophenyl-phenylether	16.52	248	269035	33.865	nq	96
68) Hexachlorobenzene	16.64	284	293900	34.333	nq	97
69) Atrazine	16.79	200	294466	43.551	nq	99
70) Pentachlorophenol	16.97	266	366473	77.662	nq	99
71) Phenanthrene	17.37	178	1282228	37.846	nq	98
72) Anthracene	17.46	178	1313439	39.227	nq	99
73) Carbazole	17.72	167	1219401	39.426	nq	98
74) Di-n-butylphthalate	18.29	149	1501929	38.034	nq	97
75) Fluoranthene	19.37	202	1509252	38.952	nq	98
77) Benzidine	19.55	184	458258	29.133	nq	96
78) Pyrene	19.74	202	1508334	36.925	nq	97
80) Butylbenzylphthalate	20.63	149	700874	36.039	nq	96
81) Benzo(a)anthracene	21.55	228	1438495	37.071	nq	99
82) 3,3'-Dichlorobenzidine	21.47	252	473374	30.878	nq	100
83) Chrysene	21.63	228	1439556	39.042	nq	95
84) Bis(2-ethylhexyl)phthalate	21.48	149	990469	36.911	nq	99
85) Di-n-octyl phthalate	22.67	149	1732164	38.397	nq	97
86) Indeno(1,2,3-cd)pyrene	28.28	276	1764559	35.215	nq	# 83
88) Benzo(b)fluoranthene	23.71	252	1539029	44.636	nq	98
89) Benzo(k)fluoranthene	23.79	252	1489705	45.434	nq	96
90) Benzo(a)pyrene	24.57	252	1401449	43.065	nq	98
91) Dibenzo(a,h)anthracene	28.38	278	1476501	45.177	nq	96
92) Benzo(g,h,i)perylene	29.48	276	1493537	46.006	nq	98

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

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