

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110519\
 Data File : BG043506.D
 Acq On : 5 Nov 2019 13:58
 Operator : JU
 Sample : PB124540BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124540BS

Manual Integrations
 APPROVED

mohammad
 11/6/2019 8:41:04 AM

Quant Time: Nov 06 02:13:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	36289	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	149975	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	124171	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	296025	20.00	ng	0.00
76) Chrysene-d12	21.66	240	347448	20.00	ng	0.00
87) Perylene-d12	24.85	264	354515	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	258536	130.55	ng	0.00
7) Phenol-d6	7.20	99	344069	120.76	ng	0.00
23) Nitrobenzene-d5	9.17	82	171371	58.91	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	230639	97.61	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	470633	52.37	ng	0.00
79) Terphenyl-d14	20.00	244	989800	53.44	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.47	88	29147	37.768	ng	96
3) Pyridine	3.87	79	65394	33.592	ng	98
4) n-Nitrosodimethylamine	3.78	42	42109	42.866	ng	97
6) Aniline	7.34	93	109666	33.412	ng	99
8) 2-Chlorophenol	7.59	128	102314	46.802	ng	96
9) Benzaldehyde	7.15	77	50626	36.155	ng	96
10) Phenol	7.23	94	124582	47.849	ng	96
11) bis(2-Chloroethyl)ether	7.43	93	82176	40.823	ng	99
12) 1,3-Dichlorobenzene	7.90	146	111261	40.621	ng	99
13) 1,4-Dichlorobenzene	8.05	146	114442	41.297	ng	98
14) 1,2-Dichlorobenzene	8.37	146	107991	39.912	ng	98
15) Benzyl Alcohol	8.25	79	91849	45.900	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.54	45	113062	38.627	ng	96
17) 2-Methylphenol	8.47	107	86339	45.488	ng	97
18) Hexachloroethane	9.10	117	40699	40.707	ng	97
19) n-Nitroso-di-n-propylamine	8.81	70	75373	40.938	ng	92
20) 3+4-Methylphenols	8.80	107	116385	44.717	ng	95
22) Acetophenone	8.83	105	151216	39.372	ng	# 98
24) Nitrobenzene	9.22	77	112145	40.468	ng	# 96
25) Isophorone	9.74	82	216575	40.721	ng	98
26) 2-Nitrophenol	9.93	139	59785	44.686	ng	92
27) 2,4-Dimethylphenol	9.99	122	95571	49.840	ng	96
28) bis(2-Chloroethoxy)methane	10.22	93	115127	39.220	ng	99
29) 2,4-Dichlorophenol	10.48	162	111551	45.403	ng	88
30) 1,2,4-Trichlorobenzene	10.68	180	120700	38.978	ng	97
31) Naphthalene	10.87	128	319476	41.966	ng	98
32) Benzoic acid	10.18	122	48089	35.701	ng	93
33) 4-Chloroaniline	10.98	127	83700	26.822	ng	98
34) Hexachlorobutadiene	11.15	225	85660	37.421	ng	96
35) Caprolactam	11.75	113	35777m	42.281	ng	
36) 4-Chloro-3-methylphenol	12.12	107	114867	42.861	ng	92
37) 2-Methylnaphthalene	12.47	142	244520	41.862	ng	98
38) 1-Methylnaphthalene	12.69	142	230883	41.544	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.84	216	156523	34.061	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	140035	58.010	nq	97
43) 2,4,6-Trichlorophenol	13.08	196	100815	37.252	nq	96
44) 2,4,5-Trichlorophenol	13.17	196	103346	37.773	nq	96
46) 1,1'-Biphenyl	13.47	154	323057	34.199	nq	97
47) 2-Chloronaphthalene	13.52	162	253304	35.243	nq	99
48) 2-Nitroaniline	13.72	65	77047	35.867	nq	92
49) Acenaphthylene	14.37	152	412105	36.841	nq	98
50) Dimethylphthalate	14.10	163	333662	34.692	nq	99
51) 2,6-Dinitrotoluene	14.21	165	75754	36.255	nq	94
52) Acenaphthene	14.71	154	242244	35.511	nq	95
53) 3-Nitroaniline	14.55	138	56553	28.033	nq	100
54) 2,4-Dinitrophenol	14.76	184	88910	69.222	nq	95
55) Dibenzofuran	15.04	168	401890	36.329	nq	99
56) 4-Nitrophenol	14.89	139	114489	84.688	nq	91
57) 2,4-Dinitrotoluene	15.01	165	112790	37.772	nq	99
58) Fluorene	15.69	166	336883	36.309	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.28	232	110251	37.928	nq	96
60) Diethylphthalate	15.46	149	341219	35.309	nq	98
61) 4-Chlorophenyl-phenylether	15.68	204	191223	35.329	nq	98
62) 4-Nitroaniline	15.72	138	78243	37.555	nq	99
63) Azobenzene	15.98	77	280049	35.806	nq	97
65) 4,6-Dinitro-2-methylphenol	15.78	198	68259	39.182	nq	90
66) n-Nitrosodiphenylamine	15.89	169	303733	36.342	nq	99
67) 4-Bromophenyl-phenylether	16.57	248	137213	34.442	nq	96
68) Hexachlorobenzene	16.70	284	155765	34.062	nq	94
69) Atrazine	16.84	200	135311	46.594	nq	97
70) Pentachlorophenol	17.05	266	165045	79.207	nq	99
71) Phenanthrene	17.43	178	551449	36.378	nq	99
72) Anthracene	17.52	178	571804	37.702	nq	98
73) Carbazole	17.80	167	531779	38.719	nq	97
74) Di-n-butylphthalate	18.34	149	629275	37.761	nq	100
75) Fluoranthene	19.44	202	750397	37.971	nq	100
77) Benzidine	19.62	184	202470	28.955	nq	100
78) Pyrene	19.80	202	762512	35.455	nq	100
80) Butylbenzylphthalate	20.68	149	291294	35.750	nq	94
81) Benzo(a)anthracene	21.64	228	789880	36.293	nq	99
82) 3,3'-Dichlorobenzidine	21.56	252	268488	31.895	nq	96
83) Chrysene	21.70	228	760162	35.153	nq	98
84) Bis(2-ethylhexyl)phthalate	21.54	149	420999	36.373	nq	99
85) Di-n-octyl phthalate	22.74	149	725793	36.961	nq	99
86) Indeno(1,2,3-cd)pyrene	28.50	276	1013128	37.325	nq	99
88) Benzo(b)fluoranthene	23.84	252	817894	40.735	nq	98
89) Benzo(k)fluoranthene	23.91	252	846741	41.890	nq	97
90) Benzo(a)pyrene	24.71	252	768879	40.101	nq	99
91) Dibenzo(a,h)anthracene	28.55	278	858546	43.777	nq	98
92) Benzo(g,h,i)perylene	29.65	276	851782	44.031	nq	98

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

