

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041322\
 Data File : BG053139.D
 Acq On : 13 Apr 2022 15:27
 Operator : CG/JU
 Sample : N2368-03MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E14ST-EB3-041122-5-10MS

Quant Time: Apr 14 02:56:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 15:31:24 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian
 Giraldo
 04/14/2022

Supervised By :Jagrut
 Upadhyay
 04/14/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							04/14/2022
1) 1,4-Dichlorobenzene-d4	8.123	152	33109	20.000	ng	0.00	Supervised By :Jagrut Upadhyay
21) Naphthalene-d8	10.931	136	135722	20.000	ng	0.00	
39) Acenaphthene-d10	14.743	164	102999	20.000	ng	0.00	
64) Phenanthrene-d10	17.486	188	237250	20.000	ng	0.00	
76) Chrysene-d12	21.775	240	233117	20.000	ng	0.00	
86) Perylene-d12	25.070	264	247470	20.000	ng	# 0.00	04/14/2022
System Monitoring Compounds							
5) 2-Fluorophenol	5.685	112	219734	113.765	ng	0.00	
7) Phenol-d6	7.283	99	328390	110.252	ng	0.00	
23) Nitrobenzene-d5	9.286	82	230697	69.017	ng	0.00	
42) 2,4,6-Tribromophenol	16.229	330	168977	103.118	ng	0.00	
45) 2-Fluorobiphenyl	13.368	172	497840	66.649	ng	0.00	
79) Terphenyl-d14	20.089	244	863880	62.587	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	3.564	88	34457	37.356	ng	#	86
3) Pyridine	3.964	79	89633	36.844	ng	#	79
4) n-Nitrosodimethylamine	3.875	42	56149	46.608	ng		88
6) Aniline	7.441	93	114836	33.267	ng		98
8) 2-Chlorophenol	7.688	128	108470	52.019	ng		94
9) Benzaldehyde	7.253	77	84715m	47.582	ng		
10) Phenol	7.312	94	148998	50.255	ng		89
11) bis(2-Chloroethyl)ether	7.529	93	101175	47.340	ng		95
12) 1,3-Dichlorobenzene	8.011	146	111664m	46.084	ng		
13) 1,4-Dichlorobenzene	8.158	146	112202	45.422	ng		94
14) 1,2-Dichlorobenzene	8.475	146	111262	46.707	ng		96
15) Benzyl Alcohol	8.352	79	124099	46.874	ng		97
16) 2,2'-oxybis(1-Chloropr...	8.645	45	161372	45.226	ng		99
17) 2-Methylphenol	8.563	107	99093	49.391	ng		97
18) Hexachloroethane	9.209	117	44216	47.733	ng		98
19) n-Nitroso-di-n-propyla...	8.922	70	97816	44.857	ng	#	96
20) 3+4-Methylphenols	8.886	107	139835	49.373	ng		96
22) Acetophenone	8.939	105	169545	45.028	ng		97
24) Nitrobenzene	9.327	77	148284	45.608	ng		90
25) Isophorone	9.850	82	264704	46.591	ng		98
26) 2-Nitrophenol	10.038	139	62541	46.001	ng		91
27) 2,4-Dimethylphenol	10.096	122	106486	54.689	ng		94
28) bis(2-Chloroethoxy)met...	10.326	93	144416	45.245	ng		99
29) 2,4-Dichlorophenol	10.578	162	118610	48.456	ng		98
30) 1,2,4-Trichlorobenzene	10.796	180	124107	44.797	ng		96
31) Naphthalene	10.983	128	325601	44.966	ng		99
32) Benzoic acid	10.255	122	69641	54.430	ng		97
33) 4-Chloroaniline	11.083	127	55172	17.792	ng		96
34) Hexachlorobutadiene	11.271	225	89332	43.409	ng		97
35) Caprolactam	11.853	113	40627m	47.087	ng		
36) 4-Chloro-3-methylphenol	12.205	107	134449	47.179	ng		97
37) 2-Methylnaphthalene	12.581	142	240788	44.940	ng		95
38) 1-Methylnaphthalene	12.799	142	231283m	44.222	ng		
40) 1,2,4,5-Tetrachloroben...	12.951	216	156599	44.765	ng		99
41) Hexachlorocyclopentadiene	12.928	237	189019	104.671	ng		95
43) 2,4,6-Trichlorophenol	13.181	196	111121	46.072	ng		96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.257	196	121595	47.310	ng	97	04/14/2022
46) 1,1'-Biphenyl	13.580	154	332959	44.889	ng	98	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.621	162	258666	43.710	ng	94	
48) 2-Nitroaniline	13.821	65	101497	45.195	ng	91	
49) Acenaphthylene	14.467	152	441538	47.664	ng	99	
50) Dimethylphthalate	14.191	163	358819	43.555	ng	100	
51) 2,6-Dinitrotoluene	14.308	165	78471	45.658	ng	# 81	04/14/2022
52) Acenaphthene	14.808	154	327530m	52.137	ng		
53) 3-Nitroaniline	14.637	138	45023	24.775	ng	87	
54) 2,4-Dinitrophenol	14.849	184	106762	97.629	ng	# 85	
55) Dibenzofuran	15.143	168	424658	43.171	ng	98	
56) 4-Nitrophenol	14.955	139	128979	93.058	ng	# 77	
57) 2,4-Dinitrotoluene	15.096	165	111948	45.752	ng	# 82	
58) Fluorene	15.789	166	351840	44.285	ng	97	
59) 2,3,4,6-Tetrachlorophenol	15.366	232	113629	45.525	ng	99	
60) Diethylphthalate	15.548	149	371102	42.993	ng	99	
61) 4-Chlorophenyl-phenyle...	15.777	204	195714	43.456	ng	97	
62) 4-Nitroaniline	15.801	138	78659	41.018	ng	# 78	
63) Azobenzene	16.071	77	369157	42.419	ng	98	
65) 4,6-Dinitro-2-methylph...	15.865	198	71167	42.995	ng	97	
66) n-Nitrosodiphenylamine	15.988	169	313990	46.023	ng	99	
67) 4-Bromophenyl-phenylether	16.670	248	137798	45.328	ng	98	
68) Hexachlorobenzene	16.799	284	153720	46.692	ng	96	
69) Atrazine	16.934	200	133561	50.116	ng	98	
70) Pentachlorophenol	17.140	266	183187	101.812	ng	89	
71) Phenanthrene	17.528	178	614630	47.429	ng	98	
72) Anthracene	17.622	178	590794	46.018	ng	99	
73) Carbazole	17.886	167	538929	43.184	ng	99	
74) Di-n-butylphthalate	18.438	149	618390	43.691	ng	99	
75) Fluoranthene	19.537	202	770993	46.414	ng	98	
77) Benzidine	19.707	184	352762	53.037	ng	100	
78) Pyrene	19.895	202	762586	45.932	ng	99	
80) Butylbenzylphthalate	20.770	149	272071	43.955	ng	98	
81) Benzo(a)anthracene	21.751	228	730735	45.846	ng	98	
82) 3,3'-Dichlorobenzidine	21.657	252	140502	23.969	ng	98	
83) Chrysene	21.822	228	667880	45.150	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.646	149	377428	45.111	ng	99	
85) Di-n-octyl phthalate	22.879	149	635709	45.431	ng	96	
87) Indeno(1,2,3-cd)pyrene	28.853	276	824782	44.860	ng	99	
88) Benzo(b)fluoranthene	24.025	252	773946	49.568	ng	98	
89) Benzo(k)fluoranthene	24.095	252	694653	45.267	ng	97	
90) Benzo(a)pyrene	24.918	252	726279	54.601	ng	98	
91) Dibenzo(a,h)anthracene	28.918	278	706940	46.717	ng	98	
92) Benzo(g,h,i)perylene	30.040	276	712212	47.752	ng	99	

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

