

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042222\
 Data File : BG053273.D
 Acq On : 22 Apr 2022 22:51
 Operator : CG/JU
 Sample : N2487-10MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

UST-1-WMS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 04/25/2022
 Supervised By : Jagrut Upadhyay 04/25/2022

Quant Time: Apr 24 01:16:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
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Internal Standards

1) 1,4-Dichlorobenzene-d4	8.108	152	22744	20.000	ng	0.00
21) Naphthalene-d8	10.921	136	99028	20.000	ng	0.00
39) Acenaphthene-d10	14.734	164	81354	20.000	ng	0.00
64) Phenanthrene-d10	17.477	188	209281	20.000	ng	0.00
76) Chrysene-d12	21.765	240	199369	20.000	ng	0.00
86) Perylene-d12	25.061	264	209139	20.000	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.676	112	149070	112.352	ng	0.00
7) Phenol-d6	7.267	99	158076	77.258	ng	0.00
23) Nitrobenzene-d5	9.271	82	235716	96.648	ng	0.00
42) 2,4,6-Tribromophenol	16.226	330	198878	153.656	ng	0.00
45) 2-Fluorobiphenyl	13.359	172	557727	94.532	ng	0.00
79) Terphenyl-d14	20.080	244	1113567	94.333	ng	0.00

Target Compounds

Compound	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.549	88	18070	28.518	ng	# 88
3) Pyridine	3.954	79	37815	22.628	ng	# 85
4) n-Nitrosodimethylamine	3.860	42	29539	35.694	ng	80
6) Aniline	7.426	93	81852	34.518	ng	99
8) 2-Chlorophenol	7.673	128	78960	55.124	ng	89
9) Benzaldehyde	7.238	77	67602	55.274	ng	97
10) Phenol	7.297	94	59927	29.424	ng	92
11) bis(2-Chloroethyl)ether	7.520	93	82994	56.530	ng	94
12) 1,3-Dichlorobenzene	7.996	146	81545	48.991	ng	94
13) 1,4-Dichlorobenzene	8.143	146	84749	49.943	ng	95
14) 1,2-Dichlorobenzene	8.466	146	83645	51.116	ng	98
15) Benzyl Alcohol	8.342	79	85025	46.751	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.630	45	131142	53.503	ng	94
17) 2-Methylphenol	8.548	107	68288	49.548	ng	95
18) Hexachloroethane	9.200	117	30386	47.752	ng	97
19) n-Nitroso-di-n-propyla...	8.906	70	81619	54.487	ng	# 93
20) 3+4-Methylphenols	8.877	107	89906	46.210	ng	92
22) Acetophenone	8.930	105	145235	52.864	ng	# 97
24) Nitrobenzene	9.312	77	122794	51.763	ng	92
25) Isophorone	9.835	82	235534	56.818	ng	99
26) 2-Nitrophenol	10.028	139	51924	52.344	ng	87
27) 2,4-Dimethylphenol	10.087	122	86002	60.535	ng	96
28) bis(2-Chloroethoxy)met...	10.316	93	124131	53.300	ng	99
29) 2,4-Dichlorophenol	10.569	162	97126	54.381	ng	98
30) 1,2,4-Trichlorobenzene	10.780	180	96665	47.821	ng	94
31) Naphthalene	10.974	128	263590	49.891	ng	99
32) Benzoic acid	10.222	122	24701m	28.325	ng	
33) 4-Chloroaniline	11.074	127	57674	25.491	ng	92
34) Hexachlorobutadiene	11.256	225	65306	43.492	ng	99
35) Caprolactam	11.850	113	10191m	16.188	ng	
36) 4-Chloro-3-methylphenol	12.196	107	115652	55.621	ng	95
37) 2-Methylnaphthalene	12.566	142	204224	52.239	ng	97
38) 1-Methylnaphthalene	12.789	142	198914	52.126	ng	99
40) 1,2,4,5-Tetrachloroben...	12.936	216	136703	49.475	ng	98
41) Hexachlorocyclopentadiene	12.919	237	124722	87.441	ng	94
43) 2,4,6-Trichlorophenol	13.171	196	98589	51.752	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.248	196	107133	52.774	ng	97	
46) 1,1'-Biphenyl	13.571	154	294116	50.202	ng	99	
47) 2-Chloronaphthalene	13.612	162	232849	49.817	ng	97	
48) 2-Nitroaniline	13.812	65	97109	54.746	ng	93	
49) Acenaphthylene	14.458	152	409209	55.927	ng	98	
50) Dimethylphthalate	14.182	163	345088	53.033	ng	99	
51) 2,6-Dinitrotoluene	14.299	165	76098	56.058	ng	# 84	04/25/2022
52) Acenaphthene	14.798	154	302355m	60.935	ng		
53) 3-Nitroaniline	14.634	138	43911	30.591	ng	85	
54) 2,4-Dinitrophenol	14.845	184	97484	112.863	ng	# 88	
55) Dibenzofuran	15.133	168	403236	51.899	ng	97	
56) 4-Nitrophenol	14.945	139	72012	65.780	ng	# 77	
57) 2,4-Dinitrotoluene	15.092	165	110692	57.275	ng	# 86	
58) Fluorene	15.779	166	336751	53.663	ng	97	
59) 2,3,4,6-Tetrachlorophenol	15.362	232	105192	53.358	ng	98	
60) Diethylphthalate	15.539	149	362158	53.120	ng	100	
61) 4-Chlorophenyl-phenyle...	15.768	204	187156	52.612	ng	97	
62) 4-Nitroaniline	15.797	138	78262	51.669	ng	# 81	
63) Azobenzene	16.061	77	359668	52.325	ng	99	
65) 4,6-Dinitro-2-methylph...	15.862	198	68499	46.914	ng	97	
66) n-Nitrosodiphenylamine	15.979	169	303093	50.362	ng	98	
67) 4-Bromophenyl-phenylether	16.661	248	135588	50.562	ng	98	
68) Hexachlorobenzene	16.790	284	151352	52.117	ng	94	
69) Atrazine	16.925	200	134692	57.295	ng	95	
70) Pentachlorophenol	17.136	266	176529	111.224	ng	89	
71) Phenanthrene	17.524	178	597176	52.240	ng	98	
72) Anthracene	17.612	178	589736	52.074	ng	99	
73) Carbazole	17.883	167	541759	49.213	ng	99	
74) Di-n-butylphthalate	18.429	149	635362	50.889	ng	99	
75) Fluoranthene	19.527	202	734010	50.093	ng	99	
77) Benzidine	19.704	184	210476	37.001	ng	98	
78) Pyrene	19.892	202	740145	52.126	ng	99	
80) Butylbenzylphthalate	20.761	149	278295	52.572	ng	98	
81) Benzo(a)anthracene	21.742	228	725069	53.191	ng	98	
82) 3,3'-Dichlorobenzidine	21.648	252	175025	34.913	ng	99	
83) Chrysene	21.813	228	661126	52.259	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.636	149	386314	53.989	ng	99	
85) Di-n-octyl phthalate	22.864	149	656894	54.891	ng	97	
87) Indeno(1,2,3-cd)pyrene	28.838	276	821422	52.866	ng	# 94	
88) Benzo(b)fluoranthene	24.015	252	771864	58.495	ng	98	
89) Benzo(k)fluoranthene	24.086	252	703944	54.280	ng	98	
90) Benzo(a)pyrene	24.908	252	719427	63.998	ng	99	
91) Dibenzo(a,h)anthracene	28.903	278	712657	55.727	ng	98	
92) Benzo(g,h,i)perylene	30.025	276	711364	56.436	ng	98	

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

