

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083122\
 Data File : BG054805.D
 Acq On : 31 Aug 2022 13:55
 Operator : CG/JU
 Sample : N4423-04MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-2MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/01/2022
 Supervised By : Jagrut Upadhyay 09/01/2022

Quant Time: Sep 01 05:00:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 25 17:50:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.152	152	36288	20.000	ng	0.00
21) Naphthalene-d8	10.978	136	152690	20.000	ng	0.00
39) Acenaphthene-d10	14.785	164	109657	20.000	ng	0.00
64) Phenanthrene-d10	17.529	188	270494	20.000	ng	0.00
76) Chrysene-d12	21.824	240	257080	20.000	ng	0.00
86) Perylene-d12	25.196	264	275459	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.690	112	248230	119.119	ng	0.00
7) Phenol-d6	7.300	99	353187	123.930	ng	0.00
23) Nitrobenzene-d5	9.321	82	203048	69.810	ng	0.00
42) 2,4,6-Tribromophenol	16.272	330	192318	140.354	ng	0.00
45) 2-Fluorobiphenyl	13.410	172	521751	71.514	ng	0.00
79) Terphenyl-d14	20.132	244	1023305	78.869	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.528	88	39091	39.284	ng	98
3) Pyridine	3.933	79	104223	37.551	ng	99
4) n-Nitrosodimethylamine	3.851	42	53019	44.190	ng	95
6) Aniline	7.464	93	158024	46.669	ng	96
8) 2-Chlorophenol	7.711	128	120383	52.875	ng	97
9) Benzaldehyde	7.282	77	67898	36.351	ng	97
10) Phenol	7.329	94	154053	52.563	ng	96
11) bis(2-Chloroethyl)ether	7.558	93	108480	46.022	ng	98
12) 1,3-Dichlorobenzene	8.034	146	125366	47.581	ng	97
13) 1,4-Dichlorobenzene	8.187	146	126924	47.758	ng	97
14) 1,2-Dichlorobenzene	8.504	146	123551	49.028	ng	99
15) Benzyl Alcohol	8.387	79	107360m	52.145	ng	
16) 2,2'-oxybis(1-Chloropr...	8.675	45	166054	45.022	ng	98
17) 2-Methylphenol	8.586	107	105763	52.413	ng	97
18) Hexachloroethane	9.239	117	44683	46.366	ng	97
19) n-Nitroso-di-n-propyla...	8.957	70	91137	46.504	ng	# 94
20) 3+4-Methylphenols	8.915	107	145710	52.073	ng	98
22) Acetophenone	8.980	105	175799	46.021	ng	# 97
24) Nitrobenzene	9.368	77	129509	44.176	ng	97
25) Isophorone	9.885	82	259500	47.849	ng	99
26) 2-Nitrophenol	10.079	139	66311	51.206	ng	97
27) 2,4-Dimethylphenol	10.132	122	114973	56.035	ng	99
28) bis(2-Chloroethoxy)met...	10.367	93	152641	49.393	ng	98
29) 2,4-Dichlorophenol	10.619	162	121888	52.228	ng	97
30) 1,2,4-Trichlorobenzene	10.831	180	119870	44.946	ng	99
31) Naphthalene	11.025	128	365853	44.659	ng	99
32) Benzoic acid	10.273	122	62007m	43.252	ng	
33) 4-Chloroaniline	11.131	127	131702	39.434	ng	98
34) Hexachlorobutadiene	11.301	225	76675	44.872	ng	98
35) Caprolactam	11.906	113	47107m	57.214	ng	
36) 4-Chloro-3-methylphenol	12.241	107	137012	53.687	ng	99
37) 2-Methylnaphthalene	12.623	142	268236	46.717	ng	97
38) 1-Methylnaphthalene	12.840	142	256285	45.877	ng	98
40) 1,2,4,5-Tetrachloroben...	12.981	216	149167	45.126	ng	99
41) Hexachlorocyclopentadiene	12.958	237	163956	93.527	ng	98
43) 2,4,6-Trichlorophenol	13.222	196	107228	48.872	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083122\
 Data File : BG054805.D
 Acq On : 31 Aug 2022 13:55
 Operator : CG/JU
 Sample : N4423-04MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-2MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/01/2022
 Supervised By : Jagrut Upadhyay 09/01/2022

Quant Time: Sep 01 05:00:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 25 17:50:55 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.293	196	118927	48.266	ng	99
46) 1,1'-Biphenyl	13.622	154	371744	45.651	ng	99
47) 2-Chloronaphthalene	13.669	162	274874	44.446	ng	99
48) 2-Nitroaniline	13.868	65	93981	48.490	ng	93
49) Acenaphthylene	14.509	152	472961	46.344	ng	100
50) Dimethylphthalate	14.233	163	399158	47.248	ng	100
51) 2,6-Dinitrotoluene	14.356	165	86498	50.058	ng	89
52) Acenaphthene	14.850	154	332649m	50.744	ng	
53) 3-Nitroaniline	14.691	138	80624	41.680	ng	95
54) 2,4-Dinitrophenol	14.891	184	96644	97.503	ng	95
55) Dibenzofuran	15.185	168	447255	46.431	ng	100
56) 4-Nitrophenol	14.991	139	159010	106.224	ng	97
57) 2,4-Dinitrotoluene	15.143	165	125795	52.632	ng	84
58) Fluorene	15.831	166	381815	47.051	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.402	232	109134	49.134	ng	99
60) Diethylphthalate	15.590	149	402309	47.690	ng	98
61) 4-Chlorophenyl-phenylether	15.813	204	196037	48.970	ng	100
62) 4-Nitroaniline	15.854	138	98167	49.706	ng	96
63) Azobenzene	16.113	77	357960	44.649	ng	96
65) 4,6-Dinitro-2-methylph...	15.901	198	68058	49.429	ng	98
66) n-Nitrosodiphenylamine	16.031	169	349738	44.835	ng	100
67) 4-Bromophenyl-phenylether	16.712	248	139532	46.984	ng	98
68) Hexachlorobenzene	16.830	284	156437	46.858	ng	97
69) Atrazine	16.971	200	142282	51.702	ng	99
70) Pentachlorophenol	17.176	266	174012	95.930	ng	99
71) Phenanthrene	17.576	178	647320	44.923	ng	99
72) Anthracene	17.664	178	656955	46.894	ng	99
73) Carbazole	17.934	167	613298	47.515	ng	99
74) Di-n-butylphthalate	18.475	149	748623	45.835	ng	99
75) Fluoranthene	19.579	202	801685	45.735	ng	99
77) Benzidine	19.756	184	437515	68.710	ng	99
78) Pyrene	19.944	202	805654	44.943	ng	99
80) Butylbenzylphthalate	20.813	149	330742	44.233	ng	95
81) Benzo(a)anthracene	21.806	228	782752	44.906	ng	99
82) 3,3'-Dichlorobenzidine	21.718	252	247297	40.465	ng	97
83) Chrysene	21.877	228	731325	43.881	ng	100
84) Bis(2-ethylhexyl)phtha...	21.689	149	477838	44.356	ng	98
85) Di-n-octyl phthalate	22.952	149	802376	43.920	ng	95
87) Indeno(1,2,3-cd)pyrene	29.086	276	932866	44.609	ng	99
88) Benzo(b)fluoranthene	24.121	252	817774	47.325	ng	99
89) Benzo(k)fluoranthene	24.192	252	799274	45.961	ng	99
90) Benzo(a)pyrene	25.038	252	725430	49.138	ng	98
91) Dibenzo(a,h)anthracene	29.151	278	805040	47.253	ng	99
92) Benzo(g,h,i)perylene	30.302	276	804306	46.731	ng	98

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

Manual Integrations

Quant Time: Sep 01 05:00:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 25 17:50:55 2022
Response via : Initial Calibration

