

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071822\
 Data File : BG054311.D
 Acq On : 18 Jul 2022 14:29
 Operator : CG/JU
 Sample : N3214-08
 Misc : LOQ-water-10ppm
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER-02-QT2-2022

Quant Time: Jul 18 15:30:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 18 14:19:07 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 07/19/2022
 Supervised By :Jagrut Upadhyay 07/19/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.997	152	17472	20.000	ng	0.00
21) Naphthalene-d8	10.806	136	62519	20.000	ng	# 0.00
39) Acenaphthene-d10	14.637	164	48769	20.000	ng	0.00
64) Phenanthrene-d10	17.381	188	135324	20.000	ng	# 0.00
76) Chrysene-d12	21.646	240	177659	20.000	ng	# 0.00
86) Perylene-d12	24.842	264	191711	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.565	112	126239	150.854	ng	0.00
7) Phenol-d6	7.163	99	168947	147.338	ng	0.00
23) Nitrobenzene-d5	9.161	82	93059	81.059	ng	0.00
42) 2,4,6-Tribromophenol	16.123	330	158812	155.095	ng	0.00
45) 2-Fluorobiphenyl	13.256	172	290523	84.037	ng	0.00
79) Terphenyl-d14	19.989	244	739169	82.538	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.432	88	3102m	9.318	ng	
3) Pyridine	3.838	79	7050	8.348	ng	# 85
4) n-Nitrosodimethylamine	3.755	42	4045	8.566	ng	# 72
6) Aniline	7.322	93	12885	9.735	ng	98
8) 2-Chlorophenol	7.563	128	8585	8.927	ng	93
9) Benzaldehyde	7.134	77	7398	11.485	ng	# 72
10) Phenol	7.193	94	10308	9.048	ng	91
11) bis(2-Chloroethyl)ether	7.416	93	7558	9.099	ng	89
12) 1,3-Dichlorobenzene	7.892	146	11204m	9.562	ng	
13) 1,4-Dichlorobenzene	8.039	146	11377	9.402	ng	94
14) 1,2-Dichlorobenzene	8.356	146	10982	9.624	ng	92
15) Benzyl Alcohol	8.232	79	8185	8.679	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.520	45	8889	8.497	ng	91
17) 2-Methylphenol	8.444	107	7231	8.806	ng	90
18) Hexachloroethane	9.079	117	3967	9.901	ng	# 70
19) n-Nitroso-di-n-propyla...	8.797	70	6516	8.510	ng	# 87
20) 3+4-Methylphenols	8.773	107	9305	8.042	ng	96
22) Acetophenone	8.814	105	13941	9.271	ng	# 97
24) Nitrobenzene	9.202	77	10352	9.227	ng	# 89
25) Isophorone	9.719	82	18252	9.141	ng	# 94
26) 2-Nitrophenol	9.919	139	5095	9.009	ng	# 85
27) 2,4-Dimethylphenol	9.977	122	8261	10.574	ng	93
28) bis(2-Chloroethoxy)met...	10.207	93	9913	9.844	ng	# 98
29) 2,4-Dichlorophenol	10.459	162	10697	9.071	ng	87
30) 1,2,4-Trichlorobenzene	10.671	180	14093	9.650	ng	98
31) Naphthalene	10.859	128	28518	9.087	ng	96
32) Benzoic acid	10.148	122	2682m	10.762	ng	
33) 4-Chloroaniline	10.965	127	11784	9.233	ng	92
34) Hexachlorobutadiene	11.141	225	12133	9.400	ng	93
35) Caprolactam	11.711	113	2451	8.255	ng	# 70
36) 4-Chloro-3-methylphenol	12.093	107	9189	8.565	ng	92
37) 2-Methylnaphthalene	12.463	142	21191	8.793	ng	99
38) 1-Methylnaphthalene	12.680	142	21035	8.957	ng	93
40) 1,2,4,5-Tetrachloroben...	12.833	216	21604	10.111	ng	# 93
41) Hexachlorocyclopentadiene	12.815	237	6674	13.314	ng	93
43) 2,4,6-Trichlorophenol	13.074	196	10361	8.906	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071822\
 Data File : BG054311.D
 Acq On : 18 Jul 2022 14:29
 Operator : CG/JU
 Sample : N3214-08
 Misc : LOQ-water-10ppm
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER-02-QT2-2022

Quant Time: Jul 18 15:30:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 18 14:19:07 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 07/19/2022
 Supervised By :Jagrut Upadhyay 07/19/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.150	196	11704	8.904	ng	# 89
46) 1,1'-Biphenyl	13.468	154	32366	9.701	ng	97
47) 2-Chloronaphthalene	13.509	162	25877	9.268	ng	94
48) 2-Nitroaniline	13.708	65	6458	8.862	ng	89
49) Acenaphthylene	14.355	152	40546	9.773	ng	98
50) Dimethylphthalate	14.079	163	34632	9.091	ng	# 97
51) 2,6-Dinitrotoluene	14.202	165	7521	9.001	ng	# 70
52) Acenaphthene	14.695	154	23004	9.156	ng	98
53) 3-Nitroaniline	14.531	138	5794	8.291	ng	# 90
54) 2,4-Dinitrophenol	14.754	184	4076	10.040	ng	# 76
55) Dibenzofuran	15.030	168	41940	9.382	ng	# 95
56) 4-Nitrophenol	14.860	139	3550	7.614	ng	90
57) 2,4-Dinitrotoluene	14.995	165	10600	8.850	ng	# 84
58) Fluorene	15.677	166	33779	9.191	ng	92
59) 2,3,4,6-Tetrachlorophenol	15.265	232	11407	8.913	ng	# 79
60) Diethylphthalate	15.442	149	33270	9.104	ng	95
61) 4-Chlorophenyl-phenyle...	15.671	204	23899	9.906	ng	# 89
62) 4-Nitroaniline	15.694	138	6662	9.074	ng	# 85
63) Azobenzene	15.965	77	24229	9.176	ng	# 79
65) 4,6-Dinitro-2-methylph...	15.765	198	7630	9.126	ng	77
66) n-Nitrosodiphenylamine	15.882	169	30292	8.966	ng	99
67) 4-Bromophenyl-phenylether	16.564	248	16793	8.836	ng	90
68) Hexachlorobenzene	16.693	284	20419	9.395	ng	# 89
69) Atrazine	16.828	200	15901	10.576	ng	95
70) Pentachlorophenol	17.040	266	7906	9.296	ng	94
71) Phenanthrene	17.422	178	62458	9.268	ng	98
72) Anthracene	17.510	178	62713	9.483	ng	99
73) Carbazole	17.780	167	52212	9.590	ng	97
74) Di-n-butylphthalate	18.332	149	58077	9.039	ng	97
75) Fluoranthene	19.431	202	86471	9.413	ng	97
77) Benzidine	19.607	184	39766	12.359	ng	98
78) Pyrene	19.790	202	90265	8.855	ng	99
80) Butylbenzylphthalate	20.671	149	26229	8.257	ng	# 83
81) Benzo(a)anthracene	21.623	228	102649	9.102	ng	99
82) 3,3'-Dichlorobenzidine	21.535	252	38894	11.089	ng	# 95
83) Chrysene	21.687	228	95489	8.962	ng	99
84) Bis(2-ethylhexyl)phtha...	21.529	149	38987	8.673	ng	# 90
85) Di-n-octyl phthalate	22.721	149	66151	8.626	ng	# 92
87) Indeno(1,2,3-cd)pyrene	28.509	276	125658	8.616	ng	# 98
88) Benzo(b)fluoranthene	23.826	252	107126	9.449	ng	# 96
89) Benzo(k)fluoranthene	23.896	252	103841	9.203	ng	# 95
90) Benzo(a)pyrene	24.696	252	103109	10.649	ng	# 97
91) Dibenzo(a,h)anthracene	28.562	278	110407	9.206	ng	# 97
92) Benzo(g,h,i)perylene	29.643	276	109010	9.210	ng	# 90

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

