

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090524\
 Data File : BG062667.D
 Acq On : 5 Sep 2024 15:10
 Operator : MA/JU
 Sample : P2403-09
 Misc : MDL-WATER-4 ng
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-WATER-03-QT2-2024

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 09/06/2024
 Supervised By :mohammad ahmed 09/07/2024

Quant Time: Sep 05 16:13:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 31 02:36:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.924	152	53402	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	215385	20.000	ng	0.00
39) Acenaphthene-d10	14.546	164	174723	20.000	ng	0.00
64) Phenanthrene-d10	17.284	188	497380	20.000	ng	0.00
76) Chrysene-d12	21.532	240	585851	20.000	ng	0.00
86) Perylene-d12	24.599	264	664596	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	358924	116.178	ng	0.00
7) Phenol-d6	7.084	99	489410	109.656	ng	0.00
23) Nitrobenzene-d5	9.070	82	333632	83.424	ng	0.00
42) 2,4,6-Tribromophenol	16.032	330	336265	136.754	ng	0.00
45) 2-Fluorobiphenyl	13.165	172	870449	80.696	ng	0.00
79) Terphenyl-d14	19.904	244	2058710	69.678	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.412	88	5451	4.235	ng	97
3) Pyridine	3.811	79	13316	3.591	ng	# 80
4) n-Nitrosodimethylamine	3.717	42	6969	3.866	ng	# 76
6) Aniline	7.243	93	17717	4.249	ng	# 89
8) 2-Chlorophenol	7.489	128	12060	3.636	ng	98
9) Benzaldehyde	7.066	77	13070	5.034	ng	94
10) Phenol	7.113	94	16029	3.548	ng	95
11) bis(2-Chloroethyl)ether	7.342	93	12760	3.689	ng	98
12) 1,3-Dichlorobenzene	7.812	146	14374	3.949	ng	97
13) 1,4-Dichlorobenzene	7.953	146	14370	3.836	ng	92
14) 1,2-Dichlorobenzene	8.271	146	13403	3.640	ng	99
15) Benzyl Alcohol	8.159	79	11000	3.331	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.435	45	23586	3.577	ng	97
17) 2-Methylphenol	8.359	107	10342	3.284	ng	92
18) Hexachloroethane	9.005	117	4832	3.440	ng	# 84
19) n-Nitroso-di-n-propyla...	8.717	70	10843	3.083	ng	96
20) 3+4-Methylphenols	8.682	107	13977	3.008	ng	93
22) Acetophenone	8.735	105	20765	3.574	ng	# 92
24) Nitrobenzene	9.117	77	15794	3.714	ng	# 94
25) Isophorone	9.634	82	31141	3.547	ng	99
26) 2-Nitrophenol	9.828	139	5565	3.464	ng	# 84
27) 2,4-Dimethylphenol	9.886	122	8247	3.503	ng	# 87
28) bis(2-Chloroethoxy)met...	10.122	93	17226	3.665	ng	97
29) 2,4-Dichlorophenol	10.362	162	11959	3.633	ng	95
30) 1,2,4-Trichlorobenzene	10.574	180	15314	3.925	ng	# 91
31) Naphthalene	10.762	128	42308	3.963	ng	97
32) Benzoic acid	9.939	122	5133m	2.027	ng	
33) 4-Chloroaniline	10.868	127	15768	3.753	ng	90
34) Hexachlorobutadiene	11.050	225	10344	3.796	ng	98
35) Caprolactam	11.608	113	4315	3.400	ng	# 90
36) 4-Chloro-3-methylphenol	11.990	107	12530	3.156	ng	# 94
37) 2-Methylnaphthalene	12.372	142	30126	3.638	ng	95
38) 1-Methylnaphthalene	12.589	142	29944	3.601	ng	92
40) 1,2,4,5-Tetrachloroben...	12.736	216	21348	4.063	ng	96
41) Hexachlorocyclopentadiene	12.718	237	4995	3.332	ng	97
43) 2,4,6-Trichlorophenol	12.977	196	12157	3.650	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090524\
 Data File : BG062667.D
 Acq On : 5 Sep 2024 15:10
 Operator : MA/JU
 Sample : P2403-09
 Misc : MDL-WATER-4 ng
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-WATER-03-QT2-2024

Quant Time: Sep 05 16:13:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 31 02:36:01 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 09/06/2024
 Supervised By :mohammad ahmed 09/07/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.048	196	13088	3.470	ng	# 92
46) 1,1'-Biphenyl	13.377	154	46770	4.012	ng	97
47) 2-Chloronaphthalene	13.418	162	35609	3.769	ng	97
48) 2-Nitroaniline	13.623	65	8962	3.607	ng	87
49) Acenaphthylene	14.264	152	56713	4.054	ng	99
50) Dimethylphthalate	13.993	163	48982	3.834	ng	# 99
51) 2,6-Dinitrotoluene	14.111	165	8503	3.519	ng	# 91
52) Acenaphthene	14.610	154	33350	3.813	ng	96
53) 3-Nitroaniline	14.440	138	9079	3.883	ng	86
55) Dibenzofuran	14.939	168	59992	4.000	ng	100
56) 4-Nitrophenol	14.757	139	6268	3.163	ng	87
57) 2,4-Dinitrotoluene	14.898	165	11641	3.531	ng	# 97
58) Fluorene	15.592	166	47512	4.036	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.169	232	14531	4.066	ng	# 92
60) Diethylphthalate	15.357	149	48270	3.812	ng	99
61) 4-Chlorophenyl-phenyle...	15.586	204	27709	4.101	ng	95
62) 4-Nitroaniline	15.597	138	9564	3.709	ng	97
63) Azobenzene	15.874	77	44336	3.729	ng	96
65) 4,6-Dinitro-2-methylph...	15.668	198	5928	2.671	ng	83
66) n-Nitrosodiphenylamine	15.797	169	40725	3.406	ng	97
67) 4-Bromophenyl-phenylether	16.479	248	18578	3.619	ng	93
68) Hexachlorobenzene	16.596	284	23756	3.908	ng	94
69) Atrazine	16.743	200	19669	4.843	ng	94
70) Pentachlorophenol	16.937	266	8542m	2.194	ng	
71) Phenanthrene	17.325	178	92237	3.973	ng	98
72) Anthracene	17.419	178	89550	3.923	ng	99
73) Carbazole	17.689	167	82362	3.925	ng	98
74) Di-n-butylphthalate	18.253	149	88089	3.722	ng	100
75) Fluoranthene	19.340	202	122398	4.169	ng	96
77) Benzidine	19.522	184	47911	4.537	ng	98
78) Pyrene	19.704	202	133570	3.587	ng	95
80) Butylbenzylphthalate	20.597	149	34646	3.167	ng	95
81) Benzo(a)anthracene	21.514	228	143355	3.880	ng	97
82) 3,3'-Dichlorobenzidine	21.432	252	46407	3.855	ng	96
83) Chrysene	21.573	228	140970	3.983	ng	98
84) Bis(2-ethylhexyl)phtha...	21.432	149	53553	3.223	ng	96
85) Di-n-octyl phthalate	22.595	149	88111	3.028	ng	99
87) Indeno(1,2,3-cd)pyrene	28.053	276	161082	3.780	ng	98
88) Benzo(b)fluoranthene	23.629	252	132303	3.604	ng	99
89) Benzo(k)fluoranthene	23.694	252	138489	3.857	ng	97
90) Benzo(a)pyrene	24.446	252	122053	3.764	ng	98
91) Dibenzo(a,h)anthracene	28.106	278	131114	3.735	ng	# 95
92) Benzo(g,h,i)perylene	29.129	276	126081	3.562	ng	97

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

