

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011733.D
 Acq On : 19 Sep 2022 13:42
 Operator : CG/JU
 Sample : N3980-09
 Misc : MDL-MDL-WATER-4ppm
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MDL-WATER-03-QT3-2022

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/20/2022
 Supervised By : Jagrut Upadhyay 09/22/2022

Quant Time: Sep 20 05:06:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 05:06:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	475526	20.000	ng	0.00
21) Naphthalene-d8	10.752	136	1898723	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	1066587	20.000	ng	0.00
64) Phenanthrene-d10	17.340	188	2083317	20.000	ng	# 0.01
76) Chrysene-d12	21.416	240	1909502	20.000	ng	# 0.01
86) Perylene-d12	23.874	264	1851589	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	2713405	103.863	ng	0.00
7) Phenol-d6	7.099	99	3572984	102.599	ng	0.00
23) Nitrobenzene-d5	9.105	82	2431157	75.938	ng	0.00
42) 2,4,6-Tribromophenol	16.075	330	817984	98.154	ng	0.01
45) 2-Fluorobiphenyl	13.210	172	5235079	76.296	ng	0.01
79) Terphenyl-d14	19.892	244	6724785	76.583	ng	0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.376	88	45402	3.946	ng	# 86
3) Pyridine	3.799	79	93044m	2.805	ng	
4) n-Nitrosodimethylamine	3.693	42	47064	3.701	ng	# 45
6) Aniline	7.264	93	164714	3.840	ng	91
8) 2-Chlorophenol	7.505	128	114540	3.698	ng	91
9) Benzaldehyde	7.075	77	105428	4.730	ng	96
10) Phenol	7.122	94	142556	3.640	ng	89
11) bis(2-Chloroethyl)ether	7.364	93	115900	3.714	ng	94
12) 1,3-Dichlorobenzene	7.828	146	127054	3.698	ng	# 93
13) 1,4-Dichlorobenzene	7.975	146	130842	3.745	ng	97
14) 1,2-Dichlorobenzene	8.293	146	123133	3.697	ng	98
15) Benzyl Alcohol	8.181	79	84089	3.382	ng	90
16) 2,2'-oxybis(1-Chloropr...	8.469	45	167191	4.012	ng	94
17) 2-Methylphenol	8.381	107	82104	3.228	ng	93
18) Hexachloroethane	9.028	117	42940	3.553	ng	93
19) n-Nitroso-di-n-propyla...	8.752	70	76315	3.385	ng	# 87
20) 3+4-Methylphenols	8.711	107	107748	3.085	ng	93
22) Acetophenone	8.769	105	166186	3.734	ng	# 90
24) Nitrobenzene	9.152	77	130540	3.919	ng	90
25) Isophorone	9.681	82	192344	3.396	ng	96
26) 2-Nitrophenol	9.863	139	34057	2.492	ng	# 88
27) 2,4-Dimethylphenol	9.922	122	85015	3.590	ng	91
28) bis(2-Chloroethoxy)met...	10.163	93	143825	3.971	ng	97
29) 2,4-Dichlorophenol	10.399	162	75636	2.884	ng	96
30) 1,2,4-Trichlorobenzene	10.616	180	101717	3.503	ng	99
31) Naphthalene	10.805	128	356917	3.622	ng	100
33) 4-Chloroaniline	10.916	127	132404	3.425	ng	# 91
34) Hexachlorobutadiene	11.093	225	54282	3.393	ng	97
35) Caprolactam	11.687	113	19242	2.264	ng	# 87
36) 4-Chloro-3-methylphenol	12.034	107	82852	2.937	ng	92
37) 2-Methylnaphthalene	12.410	142	229751	3.484	ng	97
38) 1-Methylnaphthalene	12.628	142	220744	3.424	ng	98
40) 1,2,4,5-Tetrachloroben...	12.775	216	102640	3.646	ng	97
41) Hexachlorocyclopentadiene	12.757	237	49444	3.519	ng	96
43) 2,4,6-Trichlorophenol	13.016	196	50181	2.621	ng	99
44) 2,4,5-Trichlorophenol	13.081	196	58481	2.732	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.416	154	307045	3.957	ng	99
47) 2-Chloronaphthalene	13.457	162	222161	3.639	ng	99
48) 2-Nitroaniline	13.663	65	44946	2.625	ng	92
49) Acenaphthylene	14.304	152	303922	3.217	ng	99
50) Dimethylphthalate	14.040	163	251819	3.384	ng	100
51) 2,6-Dinitrotoluene	14.151	165	38661	2.562	ng	94
52) Acenaphthene	14.646	154	216222m	3.441	ng	
53) 3-Nitroaniline	14.487	138	41305	2.374	ng	# 97
55) Dibenzofuran	14.981	168	330981	3.658	ng	98
57) 2,4-Dinitrotoluene	14.940	165	43339	2.217	ng	94
58) Fluorene	15.634	166	257353	3.507	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.204	232	45257	2.573	ng	91
60) Diethylphthalate	15.404	149	241553	3.261	ng	98
61) 4-Chlorophenyl-phenyle...	15.628	204	119874	3.533	ng	99
62) 4-Nitroaniline	15.645	138	40030	2.289	ng	# 81
63) Azobenzene	15.922	77	222041	3.119	ng	91
66) n-Nitrosodiphenylamine	15.840	169	206060	3.324	ng	99
67) 4-Bromophenyl-phenylether	16.522	248	61008	3.128	ng	# 91
68) Hexachlorobenzene	16.634	284	68839	3.277	ng	# 88
69) Atrazine	16.792	200	58053	3.068	ng	95
70) Pentachlorophenol	16.987	266	28709	2.155	ng	91
71) Phenanthrene	17.381	178	403089	3.556	ng	99
72) Anthracene	17.469	178	354293	3.305	ng	99
73) Carbazole	17.739	167	334100	3.303	ng	100
74) Di-n-butylphthalate	18.292	149	343741	2.820	ng	99
75) Fluoranthene	19.339	202	402642	3.293	ng	94
77) Benzidine	19.522	184	155519	3.425	ng	98
78) Pyrene	19.692	202	421261	3.271	ng	99
80) Butylbenzylphthalate	20.563	149	130713	2.422	ng	96
81) Benzo(a)anthracene	21.404	228	414024	3.364	ng	98
82) 3,3'-Dichlorobenzidine	21.333	252	122233	3.270	ng	# 98
83) Chrysene	21.457	228	406056	3.453	ng	98
84) Bis(2-ethylhexyl)phtha...	21.322	149	195154	2.514	ng	# 96
85) Di-n-octyl phthalate	22.269	149	300903	2.259	ng	98
87) Indeno(1,2,3-cd)pyrene	26.439	276	411986	3.088	ng	# 88
88) Benzo(b)fluoranthene	23.122	252	397246	3.469	ng	# 95
89) Benzo(k)fluoranthene	23.169	252	385388	3.340	ng	# 96
90) Benzo(a)pyrene	23.763	252	318825	3.370	ng	# 95
91) Dibenzo(a,h)anthracene	26.457	278	367091	3.314	ng	# 93
92) Benzo(g,h,i)perylene	27.227	276	357118	3.307	ng	# 90

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

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