

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102424\
 Data File : BM048215.D
 Acq On : 24 Oct 2024 13:51
 Operator : RC/JU
 Sample : P4368-07
 Misc : LOD-MDL-WATER-04 ng
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2024

Quant Time: Oct 24 14:39:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 18:21:59 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/25/2024
 Supervised By :mohammad ahmed 10/26/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	252931	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	1042873	20.000	ng	0.00
39) Acenaphthene-d10	14.368	164	665925	20.000	ng	0.00
64) Phenanthrene-d10	17.109	188	1330654	20.000	ng	0.00
76) Chrysene-d12	21.333	240	1195942	20.000	ng	0.00
86) Perylene-d12	24.285	264	1321753	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.304	112	1595882	106.060	ng	0.00
7) Phenol-d6	6.904	99	2117703	108.068	ng	0.00
23) Nitrobenzene-d5	8.880	82	1345570	72.851	ng	0.00
42) 2,4,6-Tribromophenol	15.857	330	860776	105.604	ng	0.00
45) 2-Fluorobiphenyl	12.986	172	3017197	69.525	ng	0.00
79) Terphenyl-d14	19.733	244	3759001	64.143	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.204	88	22807	3.703	ng	98
3) Pyridine	3.616	79	49419	3.020	ng	97
4) n-Nitrosodimethylamine	3.516	42	27739	3.993	ng	# 94
6) Aniline	7.051	93	72945	4.484	ng	98
8) 2-Chlorophenol	7.286	128	60435	3.697	ng	95
9) Benzaldehyde	6.863	77	53439	5.551	ng	99
10) Phenol	6.928	94	70398	3.571	ng	92
11) bis(2-Chloroethyl)ether	7.151	93	58617	3.733	ng	96
12) 1,3-Dichlorobenzene	7.610	146	66441	3.525	ng	98
13) 1,4-Dichlorobenzene	7.757	146	68224	3.563	ng	100
14) 1,2-Dichlorobenzene	8.069	146	65887	3.561	ng	99
15) Benzyl Alcohol	7.969	79	42045	3.363	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.239	45	86356	3.760	ng	99
17) 2-Methylphenol	8.175	107	42955	3.279	ng	98
18) Hexachloroethane	8.792	117	24356	3.471	ng	98
19) n-Nitroso-di-n-propyla...	8.528	70	43032	3.471	ng	99
20) 3+4-Methylphenols	8.504	107	56256	3.100	ng	96
22) Acetophenone	8.545	105	90189	3.544	ng	# 97
24) Nitrobenzene	8.922	77	65933	3.445	ng	98
25) Isophorone	9.451	82	111578	3.296	ng	99
26) 2-Nitrophenol	9.639	139	26170	2.962	ng	96
27) 2,4-Dimethylphenol	9.698	122	24024	2.241	ng	99
28) bis(2-Chloroethoxy)met...	9.933	93	70694	3.300	ng	96
29) 2,4-Dichlorophenol	10.180	162	46194	2.995	ng	100
30) 1,2,4-Trichlorobenzene	10.375	180	59627	3.376	ng	99
31) Naphthalene	10.563	128	187168	3.441	ng	99
32) Benzoic acid	9.786	122	25912m	2.171	ng	
33) 4-Chloroaniline	10.686	127	59852	3.393	ng	97
34) Hexachlorobutadiene	10.851	225	35358	3.230	ng	96
35) Caprolactam	11.463	113	13720	2.739	ng	93
36) 4-Chloro-3-methylphenol	11.816	107	47741	2.965	ng	99
37) 2-Methylnaphthalene	12.180	142	121033	3.330	ng	98
38) 1-Methylnaphthalene	12.398	142	114038	3.165	ng	97
40) 1,2,4,5-Tetrachloroben...	12.557	216	64610	3.395	ng	98
41) Hexachlorocyclopentadiene	12.533	237	13926	2.162	ng	85
43) 2,4,6-Trichlorophenol	12.804	196	38429	3.042	ng	96

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44) 2,4,5-Trichlorophenol	12.880	196	39887	2.832	ng	99
46) 1,1'-Biphenyl	13.198	154	171964	3.617	ng	99
47) 2-Chloronaphthalene	13.239	162	125504	3.348	ng	98
48) 2-Nitroaniline	13.457	65	28432	2.796	ng	99
49) Acenaphthylene	14.086	152	194628	3.559	ng	99
50) Dimethylphthalate	13.827	163	152403	3.352	ng	100
51) 2,6-Dinitrotoluene	13.945	165	28393	2.799	ng	93
52) Acenaphthene	14.427	154	115281	3.317	ng	99
53) 3-Nitroaniline	14.286	138	25354	2.650	ng	99
54) 2,4-Dinitrophenol	14.498	184	10913	1.841	ng	89
55) Dibenzofuran	14.768	168	191786	3.474	ng	100
56) 4-Nitrophenol	14.621	139	15921	1.865	ng	95
57) 2,4-Dinitrotoluene	14.739	165	33670	2.514	ng	# 86
58) Fluorene	15.415	166	147209	3.372	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.998	232	36526	3.128	ng	97
60) Diethylphthalate	15.192	149	146107	3.174	ng	99
61) 4-Chlorophenyl-phenyle...	15.409	204	74865	3.430	ng	94
62) 4-Nitroaniline	15.451	138	21047m	2.126	ng	
63) Azobenzene	15.704	77	132432	3.210	ng	98
65) 4,6-Dinitro-2-methylph...	15.509	198	17703	2.293	ng	97
66) n-Nitrosodiphenylamine	15.627	169	119758	3.278	ng	99
67) 4-Bromophenyl-phenylether	16.304	248	44296	3.283	ng	95
68) Hexachlorobenzene	16.421	284	54644	3.372	ng	96
69) Atrazine	16.580	200	40216	3.935	ng	96
70) Pentachlorophenol	16.768	266	28706	2.698	ng	99
71) Phenanthrene	17.151	178	232378	3.527	ng	99
72) Anthracene	17.245	178	209760	3.289	ng	98
73) Carbazole	17.521	167	194726	3.114	ng	98
74) Di-n-butylphthalate	18.074	149	230922	2.967	ng	99
75) Fluoranthene	19.168	202	250026	3.181	ng	99
77) Benzidine	19.356	184	51983m	3.766	ng	
78) Pyrene	19.527	202	264148	3.300	ng	98
80) Butylbenzylphthalate	20.427	149	98996	2.957	ng	95
81) Benzo(a)anthracene	21.315	228	262469	3.434	ng	99
82) 3,3'-Dichlorobenzidine	21.250	252	84418	3.225	ng	99
83) Chrysene	21.374	228	251672	3.431	ng	100
84) Bis(2-ethylhexyl)phtha...	21.239	149	153798	3.161	ng	99
85) Di-n-octyl phthalate	22.350	149	229033	2.718	ng	# 99
87) Indeno(1,2,3-cd)pyrene	27.609	276	288048	3.296	ng	99
88) Benzo(b)fluoranthene	23.356	252	248466	3.279	ng	99
89) Benzo(k)fluoranthene	23.415	252	251630	3.434	ng	98
90) Benzo(a)pyrene	24.144	252	219554	3.332	ng	99
91) Dibenzo(a,h)anthracene	27.656	278	237156	3.258	ng	96
92) Benzo(g,h,i)perylene	28.650	276	225355	3.024	ng	97

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

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