

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111124\
 Data File : BP023012.D
 Acq On : 11 Nov 2024 13:36
 Operator : RC/JU
 Sample : P4368-09
 Misc : MDL-MDL-WATER 8 PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MDL-WATER-03-QT4-2024

Quant Time: Nov 11 14:38:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 23:36:11 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 11/12/2024
 Supervised By :mohammad ahmed 11/12/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	55105	20.000	ng	0.00
21) Naphthalene-d8	10.340	136	205503	20.000	ng	0.00
39) Acenaphthene-d10	14.210	164	173650	20.000	ng	0.00
64) Phenanthrene-d10	17.016	188	457647	20.000	ng	0.00
76) Chrysene-d12	21.457	240	547682	20.000	ng	0.00
86) Perylene-d12	24.704	264	679415	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.228	112	332914	114.653	ng	0.00
7) Phenol-d6	6.787	99	478282	118.112	ng	0.00
23) Nitrobenzene-d5	8.728	82	392909	90.327	ng	-0.01
42) 2,4,6-Tribromophenol	15.728	330	454466	135.021	ng	-0.02
45) 2-Fluorobiphenyl	12.816	172	1072531	88.252	ng	0.00
79) Terphenyl-d14	19.763	244	2644265	90.673	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.175	88	10523	9.077	ng	96
3) Pyridine	3.587	79	23014m	7.230	ng	
4) n-Nitrosodimethylamine	3.493	42	13226m	8.619	ng	
6) Aniline	6.934	93	34401	10.191	ng	99
8) 2-Chlorophenol	7.163	128	24443	8.031	ng	95
9) Benzaldehyde	6.763	77	25035	10.697	ng	95
10) Phenol	6.811	94	33719	8.408	ng	86
11) bis(2-Chloroethyl)ether	7.028	93	25235	8.329	ng	97
12) 1,3-Dichlorobenzene	7.481	146	29095	7.491	ng	94
13) 1,4-Dichlorobenzene	7.622	146	30160	7.604	ng	96
14) 1,2-Dichlorobenzene	7.934	146	29836	7.780	ng	94
15) Benzyl Alcohol	7.846	79	22061	7.193	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.099	45	29819	9.071	ng	98
17) 2-Methylphenol	8.046	107	21797	8.007	ng	92
18) Hexachloroethane	8.640	117	11905	8.113	ng	91
19) n-Nitroso-di-n-propyla...	8.381	70	23105	8.336	ng	# 92
20) 3+4-Methylphenols	8.375	107	29315	7.705	ng	98
22) Acetophenone	8.405	105	47131	8.616	ng	# 90
24) Nitrobenzene	8.775	77	37786	8.552	ng	96
25) Isophorone	9.293	82	62677	8.513	ng	100
26) 2-Nitrophenol	9.481	139	13365	7.366	ng	93
27) 2,4-Dimethylphenol	9.552	122	15762	7.608	ng	89
28) bis(2-Chloroethoxy)met...	9.763	93	35194	8.416	ng	# 98
29) 2,4-Dichlorophenol	10.034	162	27243	7.596	ng	97
30) 1,2,4-Trichlorobenzene	10.210	180	37312	8.138	ng	94
31) Naphthalene	10.387	128	84445	8.201	ng	99
32) Benzoic acid	9.669	122	16100m	6.471	ng	
33) 4-Chloroaniline	10.522	127	31953	8.718	ng	96
34) Hexachlorobutadiene	10.663	225	26131	7.771	ng	99
35) Caprolactam	11.304	113	8131m	7.573	ng	
36) 4-Chloro-3-methylphenol	11.651	107	31054	7.766	ng	97
37) 2-Methylnaphthalene	12.004	142	64068	8.308	ng	96
38) 1-Methylnaphthalene	12.222	142	60894	7.956	ng	97
40) 1,2,4,5-Tetrachloroben...	12.381	216	49127	8.357	ng	99
41) Hexachlorocyclopentadiene	12.346	237	12894	7.810	ng	98
43) 2,4,6-Trichlorophenol	12.640	196	29206	7.811	ng	98

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44) 2,4,5-Trichlorophenol	12.728	196	30316	7.123	ng	96
46) 1,1'-Biphenyl	13.022	154	99723	8.639	ng	99
47) 2-Chloronaphthalene	13.069	162	77960	8.138	ng	95
48) 2-Nitroaniline	13.298	65	20900	7.488	ng	95
49) Acenaphthylene	13.934	152	116893	8.737	ng	98
50) Dimethylphthalate	13.657	163	106210	8.354	ng	99
51) 2,6-Dinitrotoluene	13.793	165	21672	7.497	ng	# 86
52) Acenaphthene	14.275	154	67791	7.880	ng	99
53) 3-Nitroaniline	14.151	138	17201	7.323	ng	93
54) 2,4-Dinitrophenol	14.369	184	10441	5.820	ng	97
55) Dibenzofuran	14.622	168	129892	8.598	ng	99
56) 4-Nitrophenol	14.522	139	10256m	5.309	ng	
57) 2,4-Dinitrotoluene	14.616	165	31123	7.447	ng	95
58) Fluorene	15.287	166	103659	8.258	ng	97
59) 2,3,4,6-Tetrachlorophenol	14.869	232	33791	8.232	ng	99
60) Diethylphthalate	15.045	149	106042	8.386	ng	98
61) 4-Chlorophenyl-phenyle...	15.275	204	61301	8.286	ng	93
62) 4-Nitroaniline	15.334	138	16311	6.768	ng	92
63) Azobenzene	15.569	77	96770	8.900	ng	95
65) 4,6-Dinitro-2-methylph...	15.392	198	19136	6.873	ng	97
66) n-Nitrosodiphenylamine	15.492	169	91986	8.130	ng	98
67) 4-Bromophenyl-phenylether	16.181	248	43246	7.879	ng	92
68) Hexachlorobenzene	16.304	284	55364	7.875	ng	94
69) Atrazine	16.469	200	39966	10.480	ng	96
70) Pentachlorophenol	16.681	266	28965	6.451	ng	90
71) Phenanthrene	17.063	178	183665	8.293	ng	99
72) Anthracene	17.157	178	181219	8.526	ng	99
73) Carbazole	17.451	167	163133	8.036	ng	100
74) Di-n-butylphthalate	18.022	149	173388	7.664	ng	99
75) Fluoranthene	19.163	202	240240	7.819	ng	98
77) Benzidine	19.369	184	102554	9.477	ng	98
78) Pyrene	19.539	202	256167	7.902	ng	98
80) Butylbenzylphthalate	20.492	149	80399	7.274	ng	94
81) Benzo(a)anthracene	21.439	228	284624	8.351	ng	99
82) 3,3'-Dichlorobenzidine	21.369	252	107069	7.761	ng	98
83) Chrysene	21.510	228	276093	8.649	ng	99
84) Bis(2-ethylhexyl)phtha...	21.369	149	121287	7.646	ng	98
85) Di-n-octyl phthalate	22.604	149	203201	7.665	ng	99
87) Indeno(1,2,3-cd)pyrene	28.380	276	386990	8.445	ng	# 95
88) Benzo(b)fluoranthene	23.680	252	322196	8.072	ng	95
89) Benzo(k)fluoranthene	23.751	252	315734m	8.377	ng	
90) Benzo(a)pyrene	24.557	252	289834	8.486	ng	100
91) Dibenzo(a,h)anthracene	28.433	278	319553	8.504	ng	97
92) Benzo(g,h,i)perylene	29.515	276	295621	7.677	ng	99

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

