

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110824\
 Data File : BP022933.D
 Acq On : 08 Nov 2024 14:09
 Operator : RC/JU
 Sample : P4368-07
 Misc : LOD-MDL-WATER-4 PPM
 ALS Vial : 10 Sample Multiplier: 1

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

Quant Time: Nov 08 14:39:12 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 :Jagrut Upadhyay 11/08/2024
 Supervised By :mohammad ahmed 11/11/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	55139	20.000	ng	0.00
21) Naphthalene-d8	10.346	136	201264	20.000	ng	0.00
39) Acenaphthene-d10	14.210	164	165538	20.000	ng	0.00
64) Phenanthrene-d10	17.022	188	412067	20.000	ng	0.00
76) Chrysene-d12	21.469	240	496962	20.000	ng	0.00
86) Perylene-d12	24.698	264	588616	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.228	112	320467	110.298	ng	0.00
7) Phenol-d6	6.793	99	430785	106.317	ng	0.00
23) Nitrobenzene-d5	8.740	82	347512	81.574	ng	0.00
42) 2,4,6-Tribromophenol	15.751	330	369560	115.176	ng	0.00
45) 2-Fluorobiphenyl	12.816	172	924845	79.829	ng	0.00
79) Terphenyl-d14	19.763	244	2118226	80.048	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.181	88	5285	4.556	ng	96
3) Pyridine	3.599	79	9561	3.002	ng	98
4) n-Nitrosodimethylamine	3.499	42	5879	3.829	ng	# 90
6) Aniline	6.940	93	15479	4.583	ng	93
8) 2-Chlorophenol	7.169	128	11705	3.843	ng	92
9) Benzaldehyde	6.764	77	12848	5.487	ng	91
10) Phenol	6.816	94	14394	3.587	ng	83
11) bis(2-Chloroethyl)ether	7.034	93	10752	3.546	ng	90
12) 1,3-Dichlorobenzene	7.487	146	14446	3.717	ng	91
13) 1,4-Dichlorobenzene	7.628	146	14448	3.640	ng	93
14) 1,2-Dichlorobenzene	7.934	146	14273	3.720	ng	95
15) Benzyl Alcohol	7.852	79	8586	2.798	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.105	45	12704	3.862	ng	97
17) 2-Methylphenol	8.046	107	8724	3.203	ng	97
18) Hexachloroethane	8.640	117	5261	3.583	ng	84
19) n-Nitroso-di-n-propyla...	8.393	70	9526	3.435	ng	98
20) 3+4-Methylphenols	8.387	107	12896	3.387	ng	91
22) Acetophenone	8.411	105	20632	3.851	ng	# 92
24) Nitrobenzene	8.775	77	16366	3.782	ng	96
25) Isophorone	9.299	82	24937	3.459	ng	# 95
26) 2-Nitrophenol	9.493	139	5810	3.269	ng	98
27) 2,4-Dimethylphenol	9.558	122	7076	3.487	ng	91
28) bis(2-Chloroethoxy)met...	9.775	93	14261	3.482	ng	95
29) 2,4-Dichlorophenol	10.046	162	11061	3.149	ng	94
30) 1,2,4-Trichlorobenzene	10.216	180	16520	3.679	ng	99
31) Naphthalene	10.399	128	36702	3.639	ng	98
32) Benzoic acid	9.675	122	6851m	2.812	ng	
33) 4-Chloroaniline	10.540	127	12365	3.445	ng	# 94
34) Hexachlorobutadiene	10.669	225	11393	3.459	ng	99
35) Caprolactam	11.322	113	3539m	3.366	ng	
36) 4-Chloro-3-methylphenol	11.669	107	12755	3.257	ng	94
37) 2-Methylnaphthalene	12.010	142	27326	3.618	ng	98
38) 1-Methylnaphthalene	12.234	142	25807	3.443	ng	97
40) 1,2,4,5-Tetrachloroben...	12.387	216	20856	3.721	ng	98
41) Hexachlorocyclopentadiene	12.352	237	4326	2.749	ng	97
43) 2,4,6-Trichlorophenol	12.646	196	11235	3.152	ng	96

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44) 2,4,5-Trichlorophenol	12.734	196	11317	2.789	ng	98
46) 1,1'-Biphenyl	13.034	154	43601	3.962	ng	98
47) 2-Chloronaphthalene	13.075	162	32897	3.602	ng	97
48) 2-Nitroaniline	13.316	65	7634	2.869	ng	92
49) Acenaphthylene	13.940	152	46724	3.664	ng	98
50) Dimethylphthalate	13.675	163	43731	3.608	ng	98
51) 2,6-Dinitrotoluene	13.798	165	8757	3.178	ng	100
52) Acenaphthene	14.281	154	28706	3.500	ng	95
53) 3-Nitroaniline	14.146	138	6869	3.068	ng	# 99
54) 2,4-Dinitrophenol	14.398	184	3654	2.137	ng	# 80
55) Dibenzofuran	14.628	168	52940	3.676	ng	99
56) 4-Nitrophenol	14.546	139	3002m	1.630	ng	
57) 2,4-Dinitrotoluene	14.634	165	11395	2.860	ng	# 75
58) Fluorene	15.275	166	41794	3.493	ng	97
59) 2,3,4,6-Tetrachlorophenol	14.863	232	12640	3.230	ng	99
60) Diethylphthalate	15.057	149	41619	3.453	ng	99
61) 4-Chlorophenyl-phenyle...	15.281	204	24823	3.520	ng	# 89
62) 4-Nitroaniline	15.351	138	6652m	2.895	ng	
63) Azobenzene	15.575	77	35326	3.408	ng	95
65) 4,6-Dinitro-2-methylph...	15.416	198	6278	2.504	ng	95
66) n-Nitrosodiphenylamine	15.504	169	36002	3.534	ng	95
67) 4-Bromophenyl-phenylether	16.198	248	16853	3.410	ng	# 91
68) Hexachlorobenzene	16.316	284	22165	3.501	ng	96
69) Atrazine	16.492	200	15687	4.568	ng	95
70) Pentachlorophenol	16.687	266	10504	2.598	ng	88
71) Phenanthrene	17.063	178	74939	3.758	ng	98
72) Anthracene	17.175	178	71535	3.738	ng	97
73) Carbazole	17.463	167	60832	3.328	ng	99
74) Di-n-butylphthalate	18.034	149	64279	3.155	ng	98
75) Fluoranthene	19.169	202	91392	3.304	ng	99
77) Benzidine	19.381	184	39964	4.070	ng	98
78) Pyrene	19.551	202	96924	3.295	ng	99
80) Butylbenzylphthalate	20.498	149	30525	3.043	ng	99
81) Benzo(a)anthracene	21.451	228	112328	3.632	ng	99
82) 3,3'-Dichlorobenzidine	21.380	252	40480	3.234	ng	96
83) Chrysene	21.510	228	105014	3.625	ng	99
84) Bis(2-ethylhexyl)phtha...	21.380	149	43655	3.033	ng	99
85) Di-n-octyl phthalate	22.610	149	72431	3.011	ng	99
87) Indeno(1,2,3-cd)pyrene	28.380	276	143290	3.609	ng	# 96
88) Benzo(b)fluoranthene	23.674	252	121325m	3.508	ng	
89) Benzo(k)fluoranthene	23.751	252	120212	3.681	ng	98
90) Benzo(a)pyrene	24.539	252	107772	3.642	ng	98
91) Dibenzo(a,h)anthracene	28.439	278	118610	3.643	ng	96
92) Benzo(g,h,i)perylene	29.504	276	111269	3.335	ng	95

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

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