

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090624\
 Data File : BM047461.D
 Acq On : 06 Sep 2024 17:01
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDICC080

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 09/07/2024

Supervised By :mohammad ahmed 09/07/2024

Quant Time: Sep 07 01:16:52 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090624.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sat Sep 07 01:06:03 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.334	152	235353	20.000	ng	0.00
21) Naphthalene-d8	10.092	136	867292	20.000	ng	0.00
39) Acenaphthene-d10	14.004	164	514267	20.000	ng	0.00
64) Phenanthrene-d10	16.774	188	1008921	20.000	ng	0.00
76) Chrysene-d12	21.021	240	797244	20.000	ng	0.00
86) Perylene-d12	23.727	264	978099	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.999	112	2068691	151.436	ng	0.00
7) Phenol-d6	6.569	99	2670270	154.443	ng	0.00
23) Nitrobenzene-d5	8.493	82	2537136	151.977	ng	0.00
42) 2,4,6-Tribromophenol	15.522	330	869843	155.299	ng	0.00
45) 2-Fluorobiphenyl	12.604	172	4924577	148.204	ng	0.00
79) Terphenyl-d14	19.439	244	6197189	154.236	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.011	88	434373	75.280	ng	98
3) Pyridine	3.387	79	1133058m	78.227	ng	
4) n-Nitrosodimethylamine	3.311	42	652410	80.561	ng	99
6) Aniline	6.693	93	1100417	71.752	ng	99
8) 2-Chlorophenol	6.922	128	1064946	74.414	ng	98
9) Benzaldehyde	6.510	77	627847	61.057	ng	99
10) Phenol	6.593	94	1311845	78.225	ng	98
11) bis(2-Chloroethyl)ether	6.793	93	1124731	76.368	ng	98
12) 1,3-Dichlorobenzene	7.222	146	1278579	73.924	ng	100
13) 1,4-Dichlorobenzene	7.369	146	1285749	74.022	ng	99
14) 1,2-Dichlorobenzene	7.681	146	1184003	72.852	ng	99
15) Benzyl Alcohol	7.604	79	1011272	78.517	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.875	45	1945074	74.663	ng	100
17) 2-Methylphenol	7.816	107	871182	76.986	ng	99
18) Hexachloroethane	8.381	117	516624	74.686	ng	99
19) n-Nitroso-di-n-propyla...	8.151	70	801722	74.769	ng	99
20) 3+4-Methylphenols	8.146	107	1172559	77.770	ng	96
22) Acetophenone	8.163	105	1568017	75.889	ng	# 98
24) Nitrobenzene	8.534	77	1308919	76.431	ng	99
25) Isophorone	9.057	82	2236397	76.530	ng	99
26) 2-Nitrophenol	9.234	139	637252	81.726	ng	97
27) 2,4-Dimethylphenol	9.322	122	702905	78.980	ng	99
28) bis(2-Chloroethoxy)met...	9.540	93	1401481	78.118	ng	99
29) 2,4-Dichlorophenol	9.787	162	1026942	80.786	ng	100
30) 1,2,4-Trichlorobenzene	9.963	180	1131842	76.715	ng	100
31) Naphthalene	10.140	128	3300418	75.043	ng	99
32) Benzoic acid	9.575	122	802532	94.868	ng	100
33) 4-Chloroaniline	10.287	127	1170784	77.925	ng	99
34) Hexachlorobutadiene	10.416	225	721903	75.788	ng	98
35) Caprolactam	11.116	113	336230	82.076	ng	99
36) 4-Chloro-3-methylphenol	11.451	107	1083324	78.860	ng	99
37) 2-Methylnaphthalene	11.769	142	2204155	76.341	ng	99
38) 1-Methylnaphthalene	11.992	142	2155744	75.466	ng	100
40) 1,2,4,5-Tetrachloroben...	12.157	216	1170778	76.911	ng	100
41) Hexachlorocyclopentadiene	12.116	237	291392	80.915	ng	95
43) 2,4,6-Trichlorophenol	12.428	196	786126	80.830	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.522	196	866467	82.484	ng	98
46) 1,1'-Biphenyl	12.822	154	2772691	75.765	ng	99
47) 2-Chloronaphthalene	12.863	162	2139454	75.248	ng	99
48) 2-Nitroaniline	13.104	65	781902	80.810	ng	99
49) Acenaphthylene	13.722	152	3149248	75.631	ng	99
50) Dimethylphthalate	13.486	163	2625853	77.036	ng	99
51) 2,6-Dinitrotoluene	13.616	165	638830	80.170	ng	96
52) Acenaphthene	14.069	154	1986281	75.334	ng	99
53) 3-Nitroaniline	13.957	138	646745	82.364	ng	99
54) 2,4-Dinitrophenol	14.192	184	394037	80.938	ng	98
55) Dibenzofuran	14.410	168	3104662	75.178	ng	98
56) 4-Nitrophenol	14.333	139	405519	49.065	ng	98
57) 2,4-Dinitrotoluene	14.427	165	822183	79.181	ng	99
58) Fluorene	15.069	166	2440065	74.411	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.663	232	674332	79.433	ng	98
60) Diethylphthalate	14.869	149	2604937	75.505	ng	100
61) 4-Chlorophenyl-phenyle...	15.069	204	1205861	74.863	ng	99
62) 4-Nitroaniline	15.139	138	630589	82.416	ng	97
63) Azobenzene	15.363	77	2545377	75.757	ng	100
65) 4,6-Dinitro-2-methylph...	15.204	198	503758	91.261	ng	96
66) n-Nitrosodiphenylamine	15.298	169	2076539	75.218	ng	99
67) 4-Bromophenyl-phenylether	15.969	248	758606	76.509	ng	99
68) Hexachlorobenzene	16.080	284	872116	75.776	ng	100
69) Atrazine	16.269	200	275250	41.021	ng	98
70) Pentachlorophenol	16.445	266	566669	87.097	ng	97
71) Phenanthrene	16.816	178	3643340	74.964	ng	99
72) Anthracene	16.910	178	3652945	76.239	ng	99
73) Carbazole	17.198	167	3632129	76.118	ng	99
74) Di-n-butylphthalate	17.763	149	4377579	75.749	ng	100
75) Fluoranthene	18.857	202	4349931	75.494	ng	99
77) Benzidine	19.074	184	1366674	82.139	ng	100
78) Pyrene	19.221	202	4485208	78.759	ng	99
80) Butylbenzylphthalate	20.151	149	1944062	80.021	ng	98
81) Benzo(a)anthracene	21.003	228	3929537	77.518	ng	99
82) 3,3'-Dichlorobenzidine	20.945	252	1372859	76.472	ng	99
83) Chrysene	21.062	228	3768293	75.766	ng	99
84) Bis(2-ethylhexyl)phtha...	20.945	149	2616882	76.964	ng	100
85) Di-n-octyl phthalate	21.974	149	4704404	78.599	ng	100
87) Indeno(1,2,3-cd)pyrene	26.756	276	4826520	76.255	ng	100
88) Benzo(b)fluoranthene	22.880	252	4100421	76.696	ng	99
89) Benzo(k)fluoranthene	22.939	252	3926747	73.097	ng	99
90) Benzo(a)pyrene	23.603	252	3725868	76.418	ng	99
91) Dibenzo(a,h)anthracene	26.785	278	3908955	76.101	ng	99
92) Benzo(g,h,i)perylene	27.691	276	4206014	77.678	ng	99

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

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