

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102822\
 Data File : BM037256.D
 Acq On : 28 Oct 2022 15:19
 Operator : CG/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/31/2022
 Supervised By : Jagrut Upadhyay 11/01/2022

Quant Time: Oct 29 02:56:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 02:48:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	393373	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1519292	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	848205	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	1592292	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1556247	20.000	ng	0.00
86) Perylene-d12	23.786	264	1470863	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1795572	81.008	ng	0.00
7) Phenol-d6	7.046	99	2403356	77.192	ng	0.00
23) Nitrobenzene-d5	9.045	82	2346547	82.612	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	789413	118.665	ng	0.00
45) 2-Fluorobiphenyl	13.139	172	5019392	91.246	ng	0.00
79) Terphenyl-d14	19.868	244	6596639	84.454	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.328	88	362751	37.982	ng	100
3) Pyridine	3.734	79	1095889m	37.174	ng	
4) n-Nitrosodimethylamine	3.646	42	475784	37.184	ng	100
6) Aniline	7.210	93	1463080	37.010	ng	100
8) 2-Chlorophenol	7.440	128	1069041	40.420	ng	100
9) Benzaldehyde	7.022	77	603788	30.277	ng	100
10) Phenol	7.075	94	1332749	37.373	ng	100
11) bis(2-Chloroethyl)ether	7.304	93	1049463	36.028	ng	100
12) 1,3-Dichlorobenzene	7.763	146	1156828	39.923	ng	98
13) 1,4-Dichlorobenzene	7.910	146	1186060	39.939	ng	100
14) 1,2-Dichlorobenzene	8.228	146	1140870	40.107	ng	100
15) Benzyl Alcohol	8.122	79	849521	35.020	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.410	45	1476162	34.697	ng	100
17) 2-Methylphenol	8.328	107	869917	37.361	ng	100
18) Hexachloroethane	8.951	117	418317	38.185	ng	100
19) n-Nitroso-di-n-propyla...	8.693	70	797940	33.740	ng	100
20) 3+4-Methylphenols	8.657	107	1183396	36.972	ng	100
22) Acetophenone	8.704	105	1532757	40.194	ng	100
24) Nitrobenzene	9.087	77	1194034	39.749	ng	100
25) Isophorone	9.616	82	1938192	36.910	ng	100
26) 2-Nitrophenol	9.798	139	545642	54.226	ng	100
27) 2,4-Dimethylphenol	9.857	122	789934	40.654	ng	100
28) bis(2-Chloroethoxy)met...	10.098	93	1191462	37.509	ng	100
29) 2,4-Dichlorophenol	10.328	162	910659	44.387	ng	100
30) 1,2,4-Trichlorobenzene	10.539	180	1018406	46.100	ng	100
31) Naphthalene	10.728	128	3302755	40.239	ng	100
32) Benzoic acid	10.010	122	619371	41.763	ng	100
33) 4-Chloroaniline	10.845	127	1324186	40.134	ng	100
34) Hexachlorobutadiene	11.004	225	564702	46.740	ng	100
35) Caprolactam	11.651	113	274752	38.807	ng	100
36) 4-Chloro-3-methylphenol	11.975	107	925459	38.369	ng	100
37) 2-Methylnaphthalene	12.339	142	2227700	40.404	ng	100
38) 1-Methylnaphthalene	12.557	142	2174451	40.297	ng	100
40) 1,2,4,5-Tetrachloroben...	12.704	216	1004371	48.235	ng	100
41) Hexachlorocyclopentadiene	12.675	237	479807	47.042	ng	100
43) 2,4,6-Trichlorophenol	12.945	196	701664	49.151	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.022	196	768340	48.354	ng	100
46) 1,1'-Biphenyl	13.345	154	2687301	42.974	ng	100
47) 2-Chloronaphthalene	13.392	162	2149990	44.772	ng	100
48) 2-Nitroaniline	13.604	65	634238	41.492	ng	100
49) Acenaphthylene	14.233	152	3329671	42.816	ng	100
50) Dimethylphthalate	13.975	163	2478082	41.429	ng	100
51) 2,6-Dinitrotoluene	14.098	165	537083	44.938	ng	100
52) Acenaphthene	14.575	154	2025693	41.916	ng	100
53) 3-Nitroaniline	14.427	138	612409	42.682	ng	100
54) 2,4-Dinitrophenol	14.633	184	295907	49.335	ng	100
55) Dibenzofuran	14.910	168	3111468	42.525	ng	100
56) 4-Nitrophenol	14.739	139	483691	42.168	ng	100
57) 2,4-Dinitrotoluene	14.886	165	714594	46.009	ng	100
58) Fluorene	15.557	166	2615954	43.639	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.133	232	631889	48.082	ng	100
60) Diethylphthalate	15.333	149	2396729	38.633	ng	100
61) 4-Chlorophenyl-phenyle...	15.551	204	1189552	45.408	ng	100
62) 4-Nitroaniline	15.592	138	634499	43.124	ng	100
63) Azobenzene	15.845	77	2442762	35.405	ng	100
65) 4,6-Dinitro-2-methylph...	15.645	198	402409	50.828	ng	100
66) n-Nitrosodiphenylamine	15.769	169	2113398	42.959	ng	100
67) 4-Bromophenyl-phenylether	16.445	248	691048	47.063	ng	100
68) Hexachlorobenzene	16.551	284	829780	51.032	ng	100
69) Atrazine	16.721	200	533379	39.357	ng	100
70) Pentachlorophenol	16.904	266	490745	49.177	ng	100
71) Phenanthrene	17.292	178	3831683	42.982	ng	100
72) Anthracene	17.386	178	3691102	43.422	ng	100
73) Carbazole	17.657	167	3426123	43.182	ng	100
74) Di-n-butylphthalate	18.215	149	3771913	37.964	ng	100
75) Fluoranthene	19.304	202	4197560	45.765	ng	100
77) Benzidine	19.498	184	1036167	31.734	ng	100
78) Pyrene	19.668	202	4406653	37.601	ng	100
80) Butylbenzylphthalate	20.562	149	1642436	33.997	ng	100
81) Benzo(a)anthracene	21.409	228	4282781	42.370	ng	100
82) 3,3'-Dichlorobenzidine	21.345	252	1318698	48.060	ng	100
83) Chrysene	21.462	228	4125500	42.374	ng	100
84) Bis(2-ethylhexyl)phtha...	21.333	149	2290252	32.294	ng	100
85) Di-n-octyl phthalate	22.245	149	3690572	32.316	ng	100
87) Indeno(1,2,3-cd)pyrene	26.233	276	4692499	48.309	ng	100
88) Benzo(b)fluoranthene	23.062	252	3908804	41.936	ng	100
89) Benzo(k)fluoranthene	23.115	252	4154740	43.322	ng	99
90) Benzo(a)pyrene	23.680	252	3269757	43.619	ng	100
91) Dibenzo(a,h)anthracene	26.244	278	3896563	48.154	ng	100
92) Benzo(g,h,i)perylene	26.985	276	3762455	47.891	ng	100

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

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