

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102623\
 Data File : BF135967.D
 Acq On : 26 Oct 2023 17:10
 Operator : CG\JU
 Sample : 05039-04MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MMFB-4MS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/27/2023
 Supervised By :mohammad ahmed 10/27/2023

Quant Time: Oct 26 17:43:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 26 03:58:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	176735	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	671219	20.000	ng	0.00
39) Acenaphthene-d10	9.863	164	308909	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	463494	20.000	ng	0.00
76) Chrysene-d12	14.010	240	327569	20.000	ng	0.00
86) Perylene-d12	15.498	264	297892	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	1089024	95.150	ng	0.00
7) Phenol-d6	6.463	99	1353755	95.007	ng	0.00
23) Nitrobenzene-d5	7.392	82	773929	75.672	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	284387	105.178	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	1415213	73.358	ng	0.00
79) Terphenyl-d14	12.951	244	1246877	63.597	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.640	88	185523	34.895	ng	98
3) Pyridine	3.393	79	462432	31.512	ng	98
4) n-Nitrosodimethylamine	3.363	42	243359	38.033	ng	98
6) Aniline	6.487	93	608373	31.709	ng	100
8) 2-Chlorophenol	6.610	128	527646	44.533	ng	99
9) Benzaldehyde	6.375	77	152581	19.795	ng	99
10) Phenol	6.475	94	641613m	43.072	ng	
11) bis(2-Chloroethyl)ether	6.569	93	494988	42.690	ng	99
12) 1,3-Dichlorobenzene	6.769	146	529666	42.086	ng	96
13) 1,4-Dichlorobenzene	6.845	146	540989	42.801	ng	98
14) 1,2-Dichlorobenzene	6.992	146	511209	43.554	ng	98
15) Benzyl Alcohol	6.969	79	426603	42.352	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.104	45	712000	41.697	ng	98
17) 2-Methylphenol	7.081	107	409765	39.239	ng	98
18) Hexachloroethane	7.339	117	192517	44.793	ng	93
19) n-Nitroso-di-n-propyla...	7.245	70	338123	41.729	ng	96
20) 3+4-Methylphenols	7.234	107	503340	39.936	ng	# 85
22) Acetophenone	7.239	105	673572	44.431	ng	# 92
24) Nitrobenzene	7.410	77	500975	45.227	ng	99
25) Isophorone	7.645	82	951665	44.855	ng	100
26) 2-Nitrophenol	7.722	139	231533	55.412	ng	98
27) 2,4-Dimethylphenol	7.763	122	400540	41.422	ng	99
28) bis(2-Chloroethoxy)met...	7.863	93	589166	45.221	ng	98
29) 2,4-Dichlorophenol	7.963	162	415065	47.934	ng	98
30) 1,2,4-Trichlorobenzene	8.045	180	444955	45.385	ng	99
31) Naphthalene	8.128	128	1463019	45.083	ng	99
32) Benzoic acid	7.875	122	288775m	56.236	ng	
33) 4-Chloroaniline	8.175	127	398314	27.770	ng	100
34) Hexachlorobutadiene	8.245	225	255038	46.152	ng	97
35) Caprolactam	8.551	113	139572m	48.437	ng	
36) 4-Chloro-3-methylphenol	8.657	107	420888	45.739	ng	99
37) 2-Methylnaphthalene	8.822	142	914304	42.583	ng	100
38) 1-Methylnaphthalene	8.922	142	872788	43.503	ng	100
40) 1,2,4,5-Tetrachloroben...	8.986	216	417985	47.908	ng	100
41) Hexachlorocyclopentadiene	8.975	237	220068	51.154	ng	99
43) 2,4,6-Trichlorophenol	9.098	196	271175	53.103	ng	98

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44) 2,4,5-Trichlorophenol	9.139	196	288729	47.621	ng	96
46) 1,1'-Biphenyl	9.286	154	1124134	48.026	ng	99
47) 2-Chloronaphthalene	9.310	162	825400	47.674	ng	98
48) 2-Nitroaniline	9.404	65	241389	50.458	ng	99
49) Acenaphthylene	9.727	152	1274724	46.670	ng	99
50) Dimethylphthalate	9.592	163	955631	47.704	ng	99
51) 2,6-Dinitrotoluene	9.651	165	200102	48.163	ng	90
52) Acenaphthene	9.898	154	812122	46.395	ng	99
53) 3-Nitroaniline	9.816	138	200319	41.041	ng	96
54) 2,4-Dinitrophenol	9.922	184	99094	70.397	ng	94
55) Dibenzofuran	10.075	168	1094071	43.265	ng	97
56) 4-Nitrophenol	9.975	139	311731	77.359	ng	92
57) 2,4-Dinitrotoluene	10.051	165	242841	47.331	ng	92
58) Fluorene	10.416	166	798343	42.198	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.186	232	215115	47.150	ng	99
60) Diethylphthalate	10.292	149	886956	46.431	ng	99
61) 4-Chlorophenyl-phenyle...	10.410	204	391594	42.909	ng	97
62) 4-Nitroaniline	10.433	138	187121	38.994	ng	96
63) Azobenzene	10.569	77	794897	41.004	ng	100
65) 4,6-Dinitro-2-methylph...	10.457	198	72836	48.451	ng	98
66) n-Nitrosodiphenylamine	10.527	169	723404	51.664	ng	100
67) 4-Bromophenyl-phenylether	10.904	248	240856	50.742	ng	93
68) Hexachlorobenzene	10.963	284	250793	49.460	ng	94
69) Atrazine	11.063	200	232270	60.850	ng	99
70) Pentachlorophenol	11.157	266	269398	83.596	ng	99
71) Phenanthrene	11.380	178	1157787	49.374	ng	100
72) Anthracene	11.433	178	1108878	46.183	ng	100
73) Carbazole	11.586	167	950091	44.359	ng	99
74) Di-n-butylphthalate	11.927	149	1125490	49.382	ng	100
75) Fluoranthene	12.574	202	1159712	47.009	ng	98
77) Benzidine	12.704	184	635455	100.606	ng	99
78) Pyrene	12.804	202	1250246	45.457	ng	100
80) Butylbenzylphthalate	13.439	149	443413	43.084	ng	99
81) Benzo(a)anthracene	13.998	228	1089363	48.910	ng	99
82) 3,3'-Dichlorobenzidine	13.968	252	352793	52.063	ng	97
83) Chrysene	14.039	228	1024774	47.240	ng	99
84) Bis(2-ethylhexyl)phtha...	14.004	149	568318	50.249	ng	99
85) Di-n-octyl phthalate	14.627	149	1059858	59.488	ng	100
87) Indeno(1,2,3-cd)pyrene	17.003	276	733644	43.433	ng	99
88) Benzo(b)fluoranthene	15.062	252	923023	52.065	ng	100
89) Benzo(k)fluoranthene	15.092	252	825326	47.401	ng	99
90) Benzo(a)pyrene	15.433	252	760281	46.532	ng	99
91) Dibenzo(a,h)anthracene	17.033	278	599576	41.498	ng	97
92) Benzo(g,h,i)perylene	17.462	276	581968	36.411	ng	99

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

