

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011024\
 Data File : BF137011.D
 Acq On : 10 Jan 2024 10:24
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

01/11/2024
 Supervised By :mohammad
 ahmed

Quant Time: Jan 10 10:44:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	57642	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	228224	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	121243	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	240878	20.000	ng	0.00
76) Chrysene-d12	13.980	240	139609	20.000	ng	# 0.01
86) Perylene-d12	15.468	264	112349	20.000	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	308723	83.567	ng	0.00
7) Phenol-d6	6.440	99	383364	80.085	ng	0.00
23) Nitrobenzene-d5	7.351	82	354276	79.434	ng	0.00
42) 2,4,6-Tribromophenol	10.622	330	114473	80.168	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	728555	80.821	ng	0.00
79) Terphenyl-d14	12.916	244	834779	83.560	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.540	88	62593	40.664	ng	95
3) Pyridine	3.304	79	148363	39.504	ng	93
4) n-Nitrosodimethylamine	3.281	42	72740	36.269	ng	92
6) Aniline	6.457	93	188498	39.385	ng	# 50
8) 2-Chlorophenol	6.581	128	164424	40.671	ng	93
9) Benzaldehyde	6.345	77	103847	38.140	ng	98
10) Phenol	6.451	94	203325	39.789	ng	86
11) bis(2-Chloroethyl)ether	6.528	93	152318	39.195	ng	97
12) 1,3-Dichlorobenzene	6.728	146	178553	40.486	ng	97
13) 1,4-Dichlorobenzene	6.804	146	182615	40.524	ng	98
14) 1,2-Dichlorobenzene	6.957	146	173021	41.262	ng	98
15) Benzyl Alcohol	6.934	79	129541	40.112	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.063	45	238744	37.576	ng	77
17) 2-Methylphenol	7.051	107	131230	39.182	ng	94
18) Hexachloroethane	7.292	117	64185	39.219	ng	91
19) n-Nitroso-di-n-propyla...	7.204	70	115895	39.939	ng	94
20) 3+4-Methylphenols	7.204	107	172178	41.150	ng	# 70
22) Acetophenone	7.198	105	233009	40.733	ng	# 89
24) Nitrobenzene	7.375	77	177891	39.817	ng	93
25) Isophorone	7.610	82	320637	39.509	ng	98
26) 2-Nitrophenol	7.687	139	87954	41.147	ng	99
27) 2,4-Dimethylphenol	7.728	122	106261	40.832	ng	96
28) bis(2-Chloroethoxy)met...	7.816	93	193434	39.212	ng	99
29) 2,4-Dichlorophenol	7.934	162	137638	41.082	ng	96
30) 1,2,4-Trichlorobenzene	8.004	180	154427	41.369	ng	99
31) Naphthalene	8.087	128	504723	40.255	ng	99
32) Benzoic acid	7.869	122	95883m	38.079	ng	
33) 4-Chloroaniline	8.139	127	185826	40.142	ng	97
34) Hexachlorobutadiene	8.198	225	91033	40.138	ng	99
35) Caprolactam	8.528	113	46879m	42.415	ng	
36) 4-Chloro-3-methylphenol	8.634	107	151135	40.070	ng	97
37) 2-Methylnaphthalene	8.781	142	327846	40.630	ng	98
38) 1-Methylnaphthalene	8.875	142	320591	40.497	ng	100
40) 1,2,4,5-Tetrachloroben...	8.945	216	152064	40.498	ng	99
41) Hexachlorocyclopentadiene	8.928	237	38383	34.996	ng	98
43) 2,4,6-Trichlorophenol	9.063	196	97553	40.511	ng	100

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44) 2,4,5-Trichlorophenol	9.116	196	103419	38.540	ng	99
46) 1,1'-Biphenyl	9.245	154	411839	40.479	ng	99
47) 2-Chloronaphthalene	9.269	162	313773	40.568	ng	99
48) 2-Nitroaniline	9.375	65	96036	39.651	ng	95
49) Acenaphthylene	9.686	152	477017	40.349	ng	99
50) Dimethylphthalate	9.545	163	356627	40.270	ng	99
51) 2,6-Dinitrotoluene	9.616	165	83533	41.562	ng	89
52) Acenaphthene	9.857	154	292538	39.919	ng	99
53) 3-Nitroaniline	9.792	138	86380	40.572	ng	92
54) 2,4-Dinitrophenol	9.904	184	41309	39.088	ng	# 83
55) Dibenzofuran	10.033	168	438944	39.513	ng	95
56) 4-Nitrophenol	9.975	139	57484	39.996	ng	92
57) 2,4-Dinitrotoluene	10.028	165	103635	39.719	ng	93
58) Fluorene	10.375	166	366872	41.622	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.157	232	78689	38.094	ng	99
60) Diethylphthalate	10.251	149	369823	40.248	ng	99
61) 4-Chlorophenyl-phenyle...	10.363	204	176410	41.175	ng	99
62) 4-Nitroaniline	10.410	138	83865	40.861	ng	92
63) Azobenzene	10.528	77	333202	38.825	ng	98
65) 4,6-Dinitro-2-methylph...	10.445	198	55028	39.862	ng	83
66) n-Nitrosodiphenylamine	10.492	169	314891	40.243	ng	97
67) 4-Bromophenyl-phenylether	10.857	248	106976	39.989	ng	98
68) Hexachlorobenzene	10.928	284	120567	40.443	ng	99
69) Atrazine	11.022	200	79101	39.524	ng	99
70) Pentachlorophenol	11.128	266	45540	32.773	ng	96
71) Phenanthrene	11.345	178	503022	40.699	ng	99
72) Anthracene	11.398	178	511061	40.829	ng	99
73) Carbazole	11.557	167	479555	40.636	ng	99
74) Di-n-butylphthalate	11.880	149	591807	41.097	ng	99
75) Fluoranthene	12.539	202	549968	41.600	ng	97
77) Benzidine	12.663	184	172294	41.855	ng	99
78) Pyrene	12.769	202	566726	41.352	ng	99
80) Butylbenzylphthalate	13.386	149	218070	43.729	ng	95
81) Benzo(a)anthracene	13.968	228	407379	42.023	ng	99
82) 3,3'-Dichlorobenzidine	13.933	252	112104	40.719	ng	99
83) Chrysene	14.004	228	384699	40.580	ng	98
84) Bis(2-ethylhexyl)phtha...	13.951	149	294895	47.000	ng	97
85) Di-n-octyl phthalate	14.568	149	413069	42.560	ng	99
87) Indeno(1,2,3-cd)pyrene	16.998	276	332750	41.021	ng	99
88) Benzo(b)fluoranthene	15.027	252	316830	42.222	ng	99
89) Benzo(k)fluoranthene	15.057	252	280483	39.558	ng	99
90) Benzo(a)pyrene	15.404	252	252984	39.940	ng	97
91) Dibenzo(a,h)anthracene	17.009	278	272628	40.967	ng	99
92) Benzo(g,h,i)perylene	17.462	276	283463	41.588	ng	98

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

