

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092524\
 Data File : BF139536.D
 Acq On : 25 Sep 2024 18:50
 Operator : RC/JU
 Sample : P4148-02MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STOCK-PILEMS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 09/26/2024
 Supervised By :mohammad ahmed 09/27/2024

Quant Time: Sep 26 05:03:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 14 02:58:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	82065	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	315781	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	174649	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	314578	20.000	ng	0.00
76) Chrysene-d12	14.039	240	146193	20.000	ng	0.00
86) Perylene-d12	15.510	264	159732	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	643467	127.913	ng	0.02
7) Phenol-d6	6.516	99	797764	120.821	ng	0.01
23) Nitrobenzene-d5	7.445	82	623847	92.874	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	242290	130.309	ng	0.01
45) 2-Fluorobiphenyl	9.234	172	1133101	90.994	ng	0.00
79) Terphenyl-d14	12.986	244	977967	109.802	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.834	88	87954	43.579	ng	# 77
3) Pyridine	3.546	79	222763	50.061	ng	98
4) n-Nitrosodimethylamine	3.481	42	163254	42.328	ng	# 82
6) Aniline	6.545	93	248921	49.954	ng	# 78
8) 2-Chlorophenol	6.669	128	259772	50.166	ng	98
9) Benzaldehyde	6.428	77	16977	4.450	ng	97
10) Phenol	6.534	94	305237	45.389	ng	97
11) bis(2-Chloroethyl)ether	6.616	93	274416	53.688	ng	99
12) 1,3-Dichlorobenzene	6.822	146	260664	44.262	ng	98
13) 1,4-Dichlorobenzene	6.898	146	269017	45.204	ng	99
14) 1,2-Dichlorobenzene	7.045	146	259349	46.120	ng	98
15) Benzyl Alcohol	7.022	79	262693	49.867	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.151	45	465978	59.155	ng	85
17) 2-Methylphenol	7.140	107	212007	51.531	ng	97
18) Hexachloroethane	7.387	117	100566	44.366	ng	98
19) n-Nitroso-di-n-propyla...	7.292	70	207628	53.064	ng	97
20) 3+4-Methylphenols	7.287	107	267127	47.590	ng	# 84
22) Acetophenone	7.287	105	374058	45.310	ng	98
24) Nitrobenzene	7.463	77	312093	45.144	ng	97
25) Isophorone	7.698	82	560611	52.772	ng	99
26) 2-Nitrophenol	7.775	139	142475	52.898	ng	92
27) 2,4-Dimethylphenol	7.816	122	220550m	62.077	ng	
28) bis(2-Chloroethoxy)met...	7.910	93	323567	53.443	ng	99
29) 2,4-Dichlorophenol	8.022	162	226877	50.267	ng	99
30) 1,2,4-Trichlorobenzene	8.098	180	234822	44.652	ng	98
31) Naphthalene	8.181	128	785043	47.620	ng	100
32) Benzoic acid	7.939	122	143410	43.133	ng	92
33) 4-Chloroaniline	8.234	127	121025m	23.997	ng	
34) Hexachlorobutadiene	8.292	225	150258	42.055	ng	99
35) Caprolactam	8.610	113	64901m	46.745	ng	
36) 4-Chloro-3-methylphenol	8.722	107	257243	49.870	ng	95
37) 2-Methylnaphthalene	8.875	142	518590	49.560	ng	100
38) 1-Methylnaphthalene	8.969	142	487145	47.419	ng	99
40) 1,2,4,5-Tetrachloroben...	9.039	216	247543	47.825	ng	99
41) Hexachlorocyclopentadiene	9.022	237	275699	152.551	ng	100
43) 2,4,6-Trichlorophenol	9.151	196	173208	51.435	ng	100

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44) 2,4,5-Trichlorophenol	9.198	196	182291	51.318	ng	98
46) 1,1'-Biphenyl	9.339	154	662323	49.294	ng	99
47) 2-Chloronaphthalene	9.363	162	489449	48.720	ng	99
48) 2-Nitroaniline	9.463	65	193529	51.792	ng	99
49) Acenaphthylene	9.781	152	814710	53.874	ng	100
50) Dimethylphthalate	9.639	163	604436	52.729	ng	100
51) 2,6-Dinitrotoluene	9.704	165	132512	52.182	ng	90
52) Acenaphthene	9.951	154	566904	53.712	ng	100
53) 3-Nitroaniline	9.875	138	107264	42.266	ng	89
54) 2,4-Dinitrophenol	9.981	184	115504	81.628	ng	# 85
55) Dibenzofuran	10.122	168	740691	50.350	ng	96
56) 4-Nitrophenol	10.039	139	179559	85.507	ng	# 81
57) 2,4-Dinitrotoluene	10.104	165	178895	52.180	ng	98
58) Fluorene	10.463	166	584809	48.723	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	154630	52.623	ng	96
60) Diethylphthalate	10.339	149	613661	51.876	ng	99
61) 4-Chlorophenyl-phenyle...	10.457	204	285519	48.947	ng	95
62) 4-Nitroaniline	10.486	138	128242	46.301	ng	89
63) Azobenzene	10.616	77	608449	52.296	ng	94
65) 4,6-Dinitro-2-methylph...	10.516	198	81166	48.886	ng	91
66) n-Nitrosodiphenylamine	10.575	169	508728	55.313	ng	100
67) 4-Bromophenyl-phenylether	10.945	248	168615	54.089	ng	99
68) Hexachlorobenzene	11.010	284	189052	51.673	ng	99
69) Atrazine	11.104	200	160193	73.162	ng	99
70) Pentachlorophenol	11.210	266	198939	90.619	ng	99
71) Phenanthrene	11.428	178	803227	52.924	ng	99
72) Anthracene	11.480	178	826145	55.674	ng	99
73) Carbazole	11.633	167	710938	50.058	ng	99
74) Di-n-butylphthalate	11.963	149	900920	56.945	ng	100
75) Fluoranthene	12.616	202	762307	47.033	ng	98
77) Benzidine	12.739	184	363941	173.759	ng	99
78) Pyrene	12.845	202	742213	59.128	ng	100
80) Butylbenzylphthalate	13.457	149	258860	60.585	ng	98
81) Benzo(a)anthracene	14.027	228	518385	53.531	ng	100
82) 3,3'-Dichlorobenzidine	13.992	252	143352	50.549	ng	99
83) Chrysene	14.069	228	490760	55.085	ng	100
84) Bis(2-ethylhexyl)phtha...	14.016	149	315187	57.988	ng	98
85) Di-n-octyl phthalate	14.627	149	499761	62.433	ng	99
87) Indeno(1,2,3-cd)pyrene	16.992	276	479035	50.435	ng	96
88) Benzo(b)fluoranthene	15.080	252	554047	56.241	ng	98
89) Benzo(k)fluoranthene	15.110	252	427789	47.097	ng	# 98
90) Benzo(a)pyrene	15.451	252	477853	58.488	ng	99
91) Dibenzo(a,h)anthracene	17.009	278	390611	49.769	ng	98
92) Benzo(g,h,i)perylene	17.439	276	335926	43.228	ng	# 96

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

