

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011122\
 Data File : BF126585.D
 Acq On : 11 Jan 2022 11:58
 Operator : JU/CG
 Sample : PB141938BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB141938BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/12/2022
 Supervised By : Jagrut Upadhyay 01/12/2022

Quant Time: Jan 11 14:07:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 10 14:50:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	87406	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	378184	20.000	ng	# 0.00
39) Acenaphthene-d10	9.816	164	184300	20.000	ng	0.00
64) Phenanthrene-d10	11.304	188	304723	20.000	ng	# 0.00
76) Chrysene-d12	13.951	240	216044	20.000	ng	0.00
86) Perylene-d12	15.410	264	187588	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.398	112	683915	112.366	ng	0.02
7) Phenol-d6	6.416	99	882572	111.954	ng	0.00
23) Nitrobenzene-d5	7.339	82	582180	79.043	ng	0.00
42) 2,4,6-Tribromophenol	10.610	330	218557	153.463	ng	0.00
45) 2-Fluorobiphenyl	9.133	172	875127	83.559	ng	0.00
79) Terphenyl-d14	12.892	244	893161	80.324	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.552	88	100037	32.775	ng	99
3) Pyridine	3.275	79	222436	27.017	ng	96
4) n-Nitrosodimethylamine	3.222	42	128293	37.901	ng	95
6) Aniline	6.434	93	226865	23.143	ng	# 86
8) 2-Chlorophenol	6.557	128	289229	47.650	ng	93
9) Benzaldehyde	6.316	77	179405	36.762	ng	95
10) Phenol	6.428	94	343265	39.018	ng	85
11) bis(2-Chloroethyl)ether	6.510	93	291758	43.432	ng	96
12) 1,3-Dichlorobenzene	6.710	146	273366	45.068	ng	99
13) 1,4-Dichlorobenzene	6.787	146	275647	45.527	ng	98
14) 1,2-Dichlorobenzene	6.939	146	269342	45.657	ng	97
15) Benzyl Alcohol	6.916	79	249814	45.291	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.051	45	456720	43.291	ng	100
17) 2-Methylphenol	7.034	107	231296	43.320	ng	98
18) Hexachloroethane	7.281	117	108550	45.000	ng	95
19) n-Nitroso-di-n-propyla...	7.192	70	190805	42.415	ng	96
20) 3+4-Methylphenols	7.187	107	267007	42.004	ng	94
22) Acetophenone	7.181	105	382984	43.805	ng	# 97
24) Nitrobenzene	7.357	77	303048	41.117	ng	95
25) Isophorone	7.598	82	549856	41.703	ng	98
26) 2-Nitrophenol	7.675	139	150004	45.910	ng	95
27) 2,4-Dimethylphenol	7.716	122	251201	49.169	ng	98
28) bis(2-Chloroethoxy)met...	7.810	93	354842	41.197	ng	100
29) 2,4-Dichlorophenol	7.916	162	219224	46.994	ng	98
30) 1,2,4-Trichlorobenzene	7.998	180	222009	45.663	ng	100
31) Naphthalene	8.081	128	806671	45.191	ng	99
32) Benzoic acid	7.863	122	161752	44.285	ng	95
33) 4-Chloroaniline	8.128	127	121136	16.199	ng	95
34) Hexachlorobutadiene	8.198	225	113683	48.108	ng	98
35) Caprolactam	8.510	113	82587m	69.591	ng	
36) 4-Chloro-3-methylphenol	8.622	107	255474	44.874	ng	98
37) 2-Methylnaphthalene	8.769	142	521028	49.480	ng	99
38) 1-Methylnaphthalene	8.869	142	482767	47.420	ng	98
40) 1,2,4,5-Tetrachloroben...	8.939	216	187506	46.820	ng	97
41) Hexachlorocyclopentadiene	8.928	237	184498	113.152	ng	98
43) 2,4,6-Trichlorophenol	9.051	196	134847	45.530	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.092	196	142082	44.568	ng	99
46) 1,1'-Biphenyl	9.233	154	622142	46.025	ng	100
47) 2-Chloronaphthalene	9.263	162	442286	43.199	ng	99
48) 2-Nitroaniline	9.357	65	161838	39.144	ng	89
49) Acenaphthylene	9.675	152	701506	45.176	ng	100
50) Dimethylphthalate	9.545	163	534025	46.415	ng	100
51) 2,6-Dinitrotoluene	9.604	165	118166	47.557	ng	90
52) Acenaphthene	9.851	154	499914m	49.107	ng	
53) 3-Nitroaniline	9.769	138	64241	18.857	ng	87
54) 2,4-Dinitrophenol	9.880	184	115346	89.816	ng	91
55) Dibenzofuran	10.022	168	607617	44.151	ng	97
56) 4-Nitrophenol	9.945	139	204863	84.552	ng	95
57) 2,4-Dinitrotoluene	10.010	165	155094	48.380	ng	96
58) Fluorene	10.363	166	455014	46.762	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.145	232	113271	49.503	ng	91
60) Diethylphthalate	10.245	149	505520	45.590	ng	98
61) 4-Chlorophenyl-phenyle...	10.357	204	199266	46.927	ng	95
62) 4-Nitroaniline	10.386	138	130877	40.790	ng	93
63) Azobenzene	10.516	77	571466	40.011	ng	97
65) 4,6-Dinitro-2-methylph...	10.422	198	77648	47.054	ng	91
66) n-Nitrosodiphenylamine	10.480	169	422209	44.659	ng	98
67) 4-Bromophenyl-phenylether	10.845	248	128425	47.482	ng	96
68) Hexachlorobenzene	10.916	284	159165	50.591	ng	# 91
69) Atrazine	11.010	200	129079	79.290	ng	96
70) Pentachlorophenol	11.110	266	150809	98.625	ng	98
71) Phenanthrene	11.327	178	706960	45.379	ng	99
72) Anthracene	11.380	178	723768	46.430	ng	99
73) Carbazole	11.533	167	668998	43.827	ng	99
74) Di-n-butylphthalate	11.869	149	787744	45.329	ng	99
75) Fluoranthene	12.516	202	706714	49.703	ng	98
77) Benzidine	12.639	184	231068	65.979	ng	98
78) Pyrene	12.745	202	715979	39.215	ng	99
80) Butylbenzylphthalate	13.368	149	329397	40.715	ng	95
81) Benzo(a)anthracene	13.939	228	519267	41.794	ng	99
82) 3,3'-Dichlorobenzidine	13.904	252	143645	24.782	ng	# 97
83) Chrysene	13.980	228	562986	44.175	ng	98
84) Bis(2-ethylhexyl)phtha...	13.933	149	392385	46.799	ng	98
85) Di-n-octyl phthalate	14.563	149	747533	51.409	ng	97
87) Indeno(1,2,3-cd)pyrene	16.839	276	478377	38.105	ng	93
88) Benzo(b)fluoranthene	14.998	252	579672	51.873	ng	# 94
89) Benzo(k)fluoranthene	15.027	252	504355	46.591	ng	# 95
90) Benzo(a)pyrene	15.351	252	503128	53.696	ng	# 96
91) Dibenzo(a,h)anthracene	16.856	278	403608	39.842	ng	96
92) Benzo(g,h,i)perylene	17.268	276	404771	37.524	ng	# 93

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

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