

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060321\
 Data File : BF124119.D
 Acq On : 3 Jun 2021 13:41
 Operator : JU/CG
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

mohammad
 6/4/2021 2:48:18 PM

Quant Time: Jun 03 15:44:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 03 15:43:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	38823	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	142592	20.000	ng	# 0.00
39) Acenaphthene-d10	9.904	164	85302	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	164903	20.000	ng	# 0.00
76) Chrysene-d12	14.033	240	140240	20.000	ng	0.00
86) Perylene-d12	15.509	264	109891	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.486	112	24960	9.027	ng	0.00
7) Phenol-d6	6.492	99	34222	9.541	ng	-0.02
23) Nitrobenzene-d5	7.421	82	35123	10.613	ng	-0.01
42) 2,4,6-Tribromophenol	10.686	330	11747	11.465	ng	-0.01
45) 2-Fluorobiphenyl	9.221	172	68289	9.852	ng	0.00
79) Terphenyl-d14	12.968	244	88969	9.608	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.628	88	5826	4.798	ng	# 93
3) Pyridine	3.398	79	14421	4.601	ng	97
4) n-Nitrosodimethylamine	3.328	42	7859	3.861	ng	# 87
6) Aniline	6.522	93	20154	4.949	ng	96
8) 2-Chlorophenol	6.651	128	13797	4.794	ng	90
9) Benzaldehyde	6.410	77	8654	5.576	ng	99
10) Phenol	6.504	94	18550	5.117	ng	92
11) bis(2-Chloroethyl)ether	6.592	93	13925	4.942	ng	95
12) 1,3-Dichlorobenzene	6.804	146	15403	4.595	ng	91
13) 1,4-Dichlorobenzene	6.880	146	16015	4.669	ng	98
14) 1,2-Dichlorobenzene	7.033	146	15988	4.989	ng	98
15) Benzyl Alcohol	7.004	79	14516	4.685	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.133	45	21584	3.805	ng	83
17) 2-Methylphenol	7.116	107	11398	4.885	ng	# 89
18) Hexachloroethane	7.374	117	6494	5.109	ng	95
19) n-Nitroso-di-n-propyla...	7.263	70	12174	5.023	ng	100
20) 3+4-Methylphenols	7.269	107	15472	5.034	ng	# 66
22) Acetophenone	7.269	105	19664	4.856	ng	# 97
24) Nitrobenzene	7.439	77	18256	5.395	ng	92
25) Isophorone	7.680	82	34656	5.444	ng	# 94
26) 2-Nitrophenol	7.763	139	8039	6.315	ng	97
27) 2,4-Dimethylphenol	7.798	122	11615	4.765	ng	82
28) bis(2-Chloroethoxy)met...	7.892	93	15767	4.925	ng	96
29) 2,4-Dichlorophenol	8.004	162	13078	5.301	ng	87
30) 1,2,4-Trichlorobenzene	8.086	180	14963	5.394	ng	# 94
31) Naphthalene	8.168	128	43851	5.196	ng	95
33) 4-Chloroaniline	8.216	127	14445	4.497	ng	93
34) Hexachlorobutadiene	8.286	225	11108	6.117	ng	97
35) Caprolactam	8.545	113	3672	5.467	ng	# 66
36) 4-Chloro-3-methylphenol	8.698	107	14620	5.419	ng	91
37) 2-Methylnaphthalene	8.857	142	29202	5.105	ng	88
38) 1-Methylnaphthalene	8.957	142	28460	5.332	ng	89
40) 1,2,4,5-Tetrachloroben...	9.027	216	15010	4.879	ng	# 93
43) 2,4,6-Trichlorophenol	9.139	196	9571	4.795	ng	95
44) 2,4,5-Trichlorophenol	9.180	196	10566	5.095	ng	# 94
46) 1,1'-Biphenyl	9.321	154	35186	4.683	ng	92

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	9.345	162	29849	5.002	ng	99
48) 2-Nitroaniline	9.439	65	9441	4.685	ng	97
49) Acenaphthylene	9.762	152	42536	4.824	ng	94
50) Dimethylphthalate	9.615	163	37424	5.494	ng	98
51) 2,6-Dinitrotoluene	9.680	165	8124	5.898	ng	# 83
52) Acenaphthene	9.933	154	27291m	4.522	ng	
53) 3-Nitroaniline	9.851	138	8031	5.867	ng	97
54) 2,4-Dinitrophenol	9.957	184	2177	3.536	ng	89
55) Dibenzofuran	10.104	168	44024	5.079	ng	95
56) 4-Nitrophenol	10.015	139	4887	4.316	ng	96
57) 2,4-Dinitrotoluene	10.080	165	10323	6.110	ng	# 95
58) Fluorene	10.451	166	32991	5.050	ng	93
59) 2,3,4,6-Tetrachlorophenol	10.227	232	7523	4.411	ng	# 99
60) Diethylphthalate	10.315	149	35839	5.313	ng	97
61) 4-Chlorophenyl-phenyle...	10.439	204	16538	5.052	ng	97
62) 4-Nitroaniline	10.457	138	7951	5.840	ng	# 85
63) Azobenzene	10.598	77	40081	5.342	ng	95
65) 4,6-Dinitro-2-methylph...	10.492	198	6324	6.790	ng	# 55
66) n-Nitrosodiphenylamine	10.557	169	31742	5.038	ng	93
67) 4-Bromophenyl-phenylether	10.933	248	11699	5.398	ng	95
68) Hexachlorobenzene	10.998	284	12963	5.358	ng	95
69) Atrazine	11.080	200	8639	4.855	ng	92
71) Phenanthrene	11.409	178	52373	5.011	ng	99
72) Anthracene	11.462	178	53560	5.057	ng	96
73) Carbazole	11.615	167	48327	5.216	ng	# 95
74) Di-n-butylphthalate	11.945	149	59007	5.089	ng	# 97
75) Fluoranthene	12.598	202	56978	5.265	ng	96
78) Pyrene	12.833	202	59130m	4.749	ng	
80) Butylbenzylphthalate	13.450	149	25496	5.222	ng	95
81) Benzo(a)anthracene	14.021	228	51712	5.099	ng	95
82) 3,3'-Dichlorobenzidine	13.980	252	17138	5.439	ng	97
83) Chrysene	14.056	228	51290	5.258	ng	95
84) Bis(2-ethylhexyl)phtha...	14.009	149	31030	4.504	ng	# 95
85) Di-n-octyl phthalate	14.621	149	39485	3.791	ng	# 93
87) Indeno(1,2,3-cd)pyrene	16.991	276	26765	3.154	ng	95
88) Benzo(b)fluoranthene	15.074	252	42213	4.966	ng	99
89) Benzo(k)fluoranthene	15.103	252	40702m	5.254	ng	
90) Benzo(a)pyrene	15.444	252	35359	4.812	ng	99
91) Dibenzo(a,h)anthracene	17.003	278	25443	3.560	ng	# 83
92) Benzo(g,h,i)perylene	17.438	276	23637	3.298	ng	# 92

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

