

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070721\
 Data File : BF124536.D
 Acq On : 7 Jul 2021 18:56
 Operator : JU/CG
 Sample : SSTDICC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

mohammad
 7/8/2021 12:39:36 PM

Quant Time: Jul 08 04:21:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 04:18:19 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.928	152	5267	20.000	ng	0.00
21) Naphthalene-d8	8.210	136	20305	20.000	ng	0.00
39) Acenaphthene-d10	9.969	164	11164	20.000	ng	# 0.00
64) Phenanthrene-d10	11.451	188	19729	20.000	ng	0.00
76) Chrysene-d12	14.098	240	15283	20.000	ng	0.00
86) Perylene-d12	15.586	264	11441	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	12775	38.145	ng	0.00
7) Phenol-d6	6.539	99	15546	35.551	ng	0.00
23) Nitrobenzene-d5	7.486	82	13425	27.414	ng	0.00
42) 2,4,6-Tribromophenol	10.751	330	3865	28.977	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	31812	34.906	ng	0.00
79) Terphenyl-d14	13.039	244	35400	34.693	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.775	88	2213	15.367	ng	# 100
3) Pyridine	3.522	79	5241	16.596	ng	91
4) n-Nitrosodimethylamine	3.475	42	4186	24.570	ng	98
6) Aniline	6.586	93	8193	16.771	ng	99
8) 2-Chlorophenol	6.704	128	7201	19.347	ng	97
9) Benzaldehyde	6.475	77	4631	22.750	ng	94
10) Phenol	6.551	94	7588	16.048	ng	97
11) bis(2-Chloroethyl)ether	6.657	93	6110	17.868	ng	97
12) 1,3-Dichlorobenzene	6.869	146	7675	16.978	ng	96
13) 1,4-Dichlorobenzene	6.945	146	7990	17.431	ng	98
14) 1,2-Dichlorobenzene	7.098	146	7255	16.806	ng	96
15) Benzyl Alcohol	7.057	79	5133	13.685	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.198	45	10418	23.749	ng	97
17) 2-Methylphenol	7.163	107	5332	17.903	ng	92
18) Hexachloroethane	7.439	117	2996	17.940	ng	95
19) n-Nitroso-di-n-propyla...	7.333	70	4415	15.013	ng	# 99
20) 3+4-Methylphenols	7.316	107	7048	18.129	ng	# 80
22) Acetophenone	7.333	105	9594	16.973	ng	96
24) Nitrobenzene	7.504	77	6335	12.949	ng	96
25) Isophorone	7.745	82	12002	14.255	ng	94
26) 2-Nitrophenol	7.822	139	3404	14.304	ng	98
27) 2,4-Dimethylphenol	7.851	122	5702	17.977	ng	98
28) bis(2-Chloroethoxy)met...	7.951	93	7319	17.150	ng	97
29) 2,4-Dichlorophenol	8.057	162	6087	16.166	ng	99
30) 1,2,4-Trichlorobenzene	8.145	180	6768	15.850	ng	97
31) Naphthalene	8.233	128	21440	17.947	ng	97
32) Benzoic acid	7.933	122	3491	21.239	ng	94
33) 4-Chloroaniline	8.275	127	8194	18.262	ng	98
34) Hexachlorobutadiene	8.351	225	3908	13.689	ng	97
35) Caprolactam	8.627	113	1346	13.642	ng	# 87
36) 4-Chloro-3-methylphenol	8.745	107	5962	15.677	ng	91
37) 2-Methylnaphthalene	8.922	142	14629	17.305	ng	97
38) 1-Methylnaphthalene	9.022	142	13931	17.717	ng	91
40) 1,2,4,5-Tetrachloroben...	9.086	216	6192	15.433	ng	96
41) Hexachlorocyclopentadiene	9.074	237	3797	121.966	ng	99
43) 2,4,6-Trichlorophenol	9.192	196	4213	16.733	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.227	196	4673	17.590	ng	98
46) 1,1'-Biphenyl	9.386	154	17189	18.115	ng	96
47) 2-Chloronaphthalene	9.410	162	13398	17.262	ng	97
48) 2-Nitroaniline	9.498	65	3221	12.615	ng	97
49) Acenaphthylene	9.827	152	20812	18.550	ng	97
50) Dimethylphthalate	9.680	163	16199	18.404	ng	# 96
51) 2,6-Dinitrotoluene	9.745	165	2926	14.391	ng	96
52) Acenaphthene	9.998	154	13845	19.435	ng	97
53) 3-Nitroaniline	9.910	138	3490	19.391	ng	94
54) 2,4-Dinitrophenol	10.010	184	1437	23.553	ng	# 69
55) Dibenzofuran	10.169	168	19969	17.587	ng	99
56) 4-Nitrophenol	10.051	139	2866	30.355	ng	79
57) 2,4-Dinitrotoluene	10.145	165	4085	15.157	ng	# 85
58) Fluorene	10.516	166	15917	18.244	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.280	232	3698	19.033	ng	# 93
60) Diethylphthalate	10.380	149	17000	19.410	ng	99
61) 4-Chlorophenyl-phenyle...	10.504	204	7218	16.695	ng	93
62) 4-Nitroaniline	10.521	138	3650	21.078	ng	87
63) Azobenzene	10.668	77	14817	16.995	ng	96
65) 4,6-Dinitro-2-methylph...	10.551	198	2193	15.507	ng	# 70
66) n-Nitrosodiphenylamine	10.621	169	13375	17.672	ng	93
67) 4-Bromophenyl-phenylether	10.998	248	4236	15.960	ng	97
68) Hexachlorobenzene	11.063	284	4149	14.388	ng	92
69) Atrazine	11.145	200	4078	20.885	ng	92
70) Pentachlorophenol	11.251	266	2689	44.658	ng	89
71) Phenanthrene	11.480	178	22633	18.198	ng	97
72) Anthracene	11.527	178	23234	18.750	ng	100
73) Carbazole	11.680	167	21312	19.734	ng	100
74) Di-n-butylphthalate	12.010	149	28700	22.317	ng	98
75) Fluoranthene	12.668	202	23091	18.450	ng	98
77) Benzidine	12.786	184	11602	37.879	ng	96
78) Pyrene	12.898	202	23926	17.151	ng	99
80) Butylbenzylphthalate	13.515	149	11597	20.935	ng	96
81) Benzo(a)anthracene	14.086	228	20251	18.379	ng	97
82) 3,3'-Dichlorobenzidine	14.045	252	5523	15.644	ng	99
83) Chrysene	14.121	228	19477	17.837	ng	98
84) Bis(2-ethylhexyl)phtha...	14.074	149	15839	21.341	ng	95
85) Di-n-octyl phthalate	14.692	149	23505	20.334	ng	99
87) Indeno(1,2,3-cd)pyrene	17.092	276	12385	13.727	ng	98
88) Benzo(b)fluoranthene	15.145	252	17815m	20.548	ng	
89) Benzo(k)fluoranthene	15.174	252	15238m	19.211	ng	
90) Benzo(a)pyrene	15.521	252	13979	18.145	ng	98
91) Dibenzo(a,h)anthracene	17.115	278	10546	13.570	ng	# 88
92) Benzo(g,h,i)perylene	17.556	276	10212	13.354	ng	94

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

