

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061621\
 Data File : BG048992.D
 Acq On : 16 Jun 2021 9:59
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/17/2021 12:15:29 PM

Quant Time: Jun 16 12:05:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 16 12:04:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.114	152	33522	20.000	ng	0.00
21) Naphthalene-d8	10.940	136	137589	20.000	ng	# 0.00
39) Acenaphthene-d10	14.759	164	104964	20.000	ng	0.00
64) Phenanthrene-d10	17.508	188	253118	20.000	ng	# 0.00
76) Chrysene-d12	21.797	240	252620	20.000	ng	0.00
86) Perylene-d12	25.129	264	260614	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.663	112	159625	74.126	ng	0.00
7) Phenol-d6	7.267	99	232587	72.062	ng	0.00
23) Nitrobenzene-d5	9.283	82	234217	73.403	ng	0.00
42) 2,4,6-Tribromophenol	16.245	330	116332	74.806	ng	0.00
45) 2-Fluorobiphenyl	13.384	172	503270	67.733	ng	0.00
79) Terphenyl-d14	20.111	244	994354	72.087	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.537	88	38129	36.269	ng	# 76
3) Pyridine	3.942	79	108285	37.969	ng	92
4) n-Nitrosodimethylamine	3.854	42	58968	43.177	ng	# 77
6) Aniline	7.432	93	143465	35.296	ng	# 92
8) 2-Chlorophenol	7.679	128	86881	36.604	ng	84
9) Benzaldehyde	7.244	77	56035	33.552	ng	88
10) Phenol	7.291	94	117945	35.373	ng	96
11) bis(2-Chloroethyl)ether	7.526	93	85747	34.659	ng	81
12) 1,3-Dichlorobenzene	8.002	146	94291	35.562	ng	96
13) 1,4-Dichlorobenzene	8.149	146	95736	36.216	ng	94
14) 1,2-Dichlorobenzene	8.472	146	90981m	35.777	ng	
15) Benzyl Alcohol	8.349	79	104597	35.386	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.642	45	160757	34.342	ng	83
17) 2-Methylphenol	8.554	107	78302	35.809	ng	92
18) Hexachloroethane	9.206	117	38609	36.407	ng	98
19) n-Nitroso-di-n-propyla...	8.924	70	85938	34.064	ng	97
20) 3+4-Methylphenols	8.883	107	107472	35.948	ng	99
22) Acetophenone	8.936	105	147408	35.215	ng	# 93
24) Nitrobenzene	9.324	77	131041	36.458	ng	99
25) Isophorone	9.847	82	232819	33.492	ng	99
26) 2-Nitrophenol	10.041	139	49546	40.638	ng	97
27) 2,4-Dimethylphenol	10.094	122	75649	34.222	ng	91
28) bis(2-Chloroethoxy)met...	10.334	93	107280	33.328	ng	95
29) 2,4-Dichlorophenol	10.581	162	87620	35.331	ng	94
30) 1,2,4-Trichlorobenzene	10.799	180	101537	35.711	ng	94
31) Naphthalene	10.987	128	278979	34.035	ng	99
32) Benzoic acid	10.235	122	64242	44.493	ng	83
33) 4-Chloroaniline	11.092	127	120721	34.687	ng	96
34) Hexachlorobutadiene	11.275	225	74039	36.536	ng	95
35) Caprolactam	11.862	113	31188	43.505	ng	# 38
36) 4-Chloro-3-methylphenol	12.209	107	108270	35.718	ng	# 90
37) 2-Methylnaphthalene	12.591	142	209396	33.821	ng	98
38) 1-Methylnaphthalene	12.808	142	198143	34.089	ng	95
40) 1,2,4,5-Tetrachloroben...	12.955	216	112556	33.770	ng	98
41) Hexachlorocyclopentadiene	12.937	237	74994	35.001	ng	88
43) 2,4,6-Trichlorophenol	13.190	196	86015	38.570	ng	96

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44) 2,4,5-Trichlorophenol	13.272	196	91647	39.743	ng	95
46) 1,1'-Biphenyl	13.589	154	260423	33.653	ng	97
47) 2-Chloronaphthalene	13.636	162	214831	34.764	ng	99
48) 2-Nitroaniline	13.830	65	88000	40.502	ng	95
49) Acenaphthylene	14.483	152	343879	33.394	ng	96
50) Dimethylphthalate	14.206	163	290566	35.538	ng	98
51) 2,6-Dinitrotoluene	14.324	165	65305	39.221	ng	99
52) Acenaphthene	14.823	154	229664	34.602	ng	94
53) 3-Nitroaniline	14.653	138	64500	38.262	ng	86
54) 2,4-Dinitrophenol	14.853	184	35115	45.528	ng	92
55) Dibenzofuran	15.152	168	348227	33.752	ng	97
56) 4-Nitrophenol	14.970	139	55896	40.671	ng	89
57) 2,4-Dinitrotoluene	15.111	165	95627	46.756	ng	# 95
58) Fluorene	15.805	166	285364	33.824	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.381	232	84481	36.193	ng	97
60) Diethylphthalate	15.570	149	316531	35.745	ng	98
61) 4-Chlorophenyl-phenylether	15.793	204	156879	35.562	ng	96
62) 4-Nitroaniline	15.816	138	69950	41.030	ng	84
63) Azobenzene	16.087	77	334929	33.056	ng	87
65) 4,6-Dinitro-2-methylph...	15.875	198	57404	47.825	ng	95
66) n-Nitrosodiphenylamine	16.010	169	255087	32.338	ng	97
67) 4-Bromophenyl-phenylether	16.692	248	103120	32.628	ng	95
68) Hexachlorobenzene	16.809	284	118511	34.597	ng	97
69) Atrazine	16.956	200	94450	37.303	ng	97
70) Pentachlorophenol	17.150	266	72592	35.262	ng	97
71) Phenanthrene	17.550	178	482695	33.892	ng	98
72) Anthracene	17.644	178	470858	32.901	ng	98
73) Carbazole	17.908	167	468833	34.176	ng	97
74) Di-n-butylphthalate	18.466	149	561251	36.477	ng	99
75) Fluoranthene	19.553	202	611233	35.589	ng	96
77) Benzidine	19.729	184	239375	47.038	ng	96
78) Pyrene	19.917	202	621787	33.911	ng	97
80) Butylbenzylphthalate	20.799	149	251941	36.419	ng	98
81) Benzo(a)anthracene	21.780	228	630816	34.778	ng	98
82) 3,3'-Dichlorobenzidine	21.692	252	227953	39.300	ng	100
83) Chrysene	21.850	228	610551	35.108	ng	97
84) Bis(2-ethylhexyl)phtha...	21.686	149	355435	36.045	ng	98
85) Di-n-octyl phthalate	22.937	149	605356	36.196	ng	96
87) Indeno(1,2,3-cd)pyrene	28.954	276	741680	37.815	ng	# 96
88) Benzo(b)fluoranthene	24.071	252	670693	36.101	ng	# 95
89) Benzo(k)fluoranthene	24.142	252	634333	35.987	ng	# 99
90) Benzo(a)pyrene	24.976	252	621510	36.648	ng	# 99
91) Dibenzo(a,h)anthracene	29.018	278	628148	38.320	ng	# 97
92) Benzo(g,h,i)perylene	30.141	276	608764	36.829	ng	# 95

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

