

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020922\
 Data File : BG052343.D
 Acq On : 9 Feb 2022 20:16
 Operator : CG/JU
 Sample : N1432-07MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-3MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/10/2022
 Supervised By : Jagrut Upadhyay 02/11/2022

Quant Time: Feb 10 00:08:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 08 01:21:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.225	152	24799	20.000	ng	0.00
21) Naphthalene-d8	11.068	136	110461	20.000	ng	# 0.00
39) Acenaphthene-d10	14.874	164	84394	20.000	ng	0.00
64) Phenanthrene-d10	17.624	188	210453	20.000	ng	0.00
76) Chrysene-d12	21.941	240	199794	20.000	ng	0.00
86) Perylene-d12	25.413	264	205605	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.757	112	123558	86.836	ng	0.00
7) Phenol-d6	7.373	99	190377	86.715	ng	0.00
23) Nitrobenzene-d5	9.399	82	125921	49.544	ng	0.00
42) 2,4,6-Tribromophenol	16.361	330	100200	82.642	ng	0.00
45) 2-Fluorobiphenyl	13.494	172	304264	50.097	ng	0.00
79) Terphenyl-d14	20.214	244	583006	51.529	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.578	88	23809	36.339	ng	96
3) Pyridine	3.989	79	58962	31.029	ng	94
4) n-Nitrosodimethylamine	3.895	42	39538	40.296	ng	91
6) Aniline	7.537	93	81782	32.063	ng	97
8) 2-Chlorophenol	7.790	128	81929	52.341	ng	92
9) Benzaldehyde	7.343	77	56029	47.567	ng	99
10) Phenol	7.402	94	114586	52.529	ng	99
11) bis(2-Chloroethyl)ether	7.625	93	72810	43.668	ng	92
12) 1,3-Dichlorobenzene	8.113	146	81832	45.884	ng	97
13) 1,4-Dichlorobenzene	8.260	146	84975	46.738	ng	94
14) 1,2-Dichlorobenzene	8.589	146	83333	46.475	ng	97
15) Benzyl Alcohol	8.460	79	87631	48.094	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.747	45	100178	41.607	ng	97
17) 2-Methylphenol	8.671	107	75445	52.413	ng	95
18) Hexachloroethane	9.329	117	28834	45.610	ng	94
19) n-Nitroso-di-n-propyla...	9.035	70	70970	42.879	ng	94
20) 3+4-Methylphenols	9.000	107	104714	50.848	ng	96
22) Acetophenone	9.053	105	122630	42.005	ng	# 95
24) Nitrobenzene	9.441	77	105357	43.700	ng	96
25) Isophorone	9.969	82	194030	44.867	ng	98
26) 2-Nitrophenol	10.163	139	47547	46.529	ng	93
27) 2,4-Dimethylphenol	10.222	122	86455	57.143	ng	92
28) bis(2-Chloroethoxy)met...	10.445	93	103950	42.355	ng	100
29) 2,4-Dichlorophenol	10.709	162	89601	49.408	ng	97
30) 1,2,4-Trichlorobenzene	10.927	180	92432	43.649	ng	96
31) Naphthalene	11.121	128	257693	44.500	ng	96
32) Benzoic acid	10.363	122	56450	56.496	ng	96
33) 4-Chloroaniline	11.215	127	44682	18.481	ng	97
34) Hexachlorobutadiene	11.403	225	58986	41.245	ng	98
35) Caprolactam	11.984	113	33773m	51.102	ng	
36) 4-Chloro-3-methylphenol	12.337	107	104584	50.691	ng	96
37) 2-Methylnaphthalene	12.713	142	198424	46.133	ng	98
38) 1-Methylnaphthalene	12.930	142	189929	45.571	ng	99
40) 1,2,4,5-Tetrachloroben...	13.083	216	119605	43.088	ng	98
41) Hexachlorocyclopentadiene	13.059	237	117011	87.018	ng	97
43) 2,4,6-Trichlorophenol	13.312	196	85890	47.311	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.388	196	95117	48.347	ng	94
46) 1,1'-Biphenyl	13.705	154	279217	45.119	ng	99
47) 2-Chloronaphthalene	13.758	162	214574	45.005	ng	96
48) 2-Nitroaniline	13.952	65	78303	46.130	ng	88
49) Acenaphthylene	14.598	152	366549	47.228	ng	99
50) Dimethylphthalate	14.316	163	284767	44.047	ng	100
51) 2,6-Dinitrotoluene	14.440	165	66321	47.538	ng	# 86
52) Acenaphthene	14.939	154	264367m	52.011	ng	
53) 3-Nitroaniline	14.769	138	37260	25.695	ng	99
54) 2,4-Dinitrophenol	14.980	184	83536	95.667	ng	96
55) Dibenzofuran	15.274	168	349218	43.688	ng	99
56) 4-Nitrophenol	15.080	139	112197	103.027	ng	86
57) 2,4-Dinitrotoluene	15.227	165	95538	48.114	ng	# 88
58) Fluorene	15.920	166	296574	46.476	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.497	232	92362	49.143	ng	# 98
60) Diethylphthalate	15.673	149	296812	45.448	ng	97
61) 4-Chlorophenyl-phenyle...	15.902	204	161926	44.546	ng	97
62) 4-Nitroaniline	15.932	138	67699	44.346	ng	89
63) Azobenzene	16.196	77	291971	43.671	ng	96
65) 4,6-Dinitro-2-methylph...	15.996	198	57460	45.387	ng	92
66) n-Nitrosodiphenylamine	16.114	169	264203	45.564	ng	97
67) 4-Bromophenyl-phenylether	16.801	248	111081	43.402	ng	96
68) Hexachlorobenzene	16.936	284	125919	45.362	ng	91
69) Atrazine	17.060	200	113179	48.279	ng	95
70) Pentachlorophenol	17.277	266	143618	105.756	ng	98
71) Phenanthrene	17.671	178	488221	45.001	ng	99
72) Anthracene	17.759	178	489592	46.052	ng	99
73) Carbazole	18.023	167	453204	43.708	ng	99
74) Di-n-butylphthalate	18.564	149	515193	45.603	ng	99
75) Fluoranthene	19.668	202	608001	44.012	ng	99
77) Benzidine	19.838	184	221201	51.814	ng	99
78) Pyrene	20.032	202	615870	46.067	ng	98
80) Butylbenzylphthalate	20.902	149	223407	48.108	ng	92
81) Benzo(a)anthracene	21.924	228	593431	44.885	ng	100
82) 3,3'-Dichlorobenzidine	21.824	252	120661	26.142	ng	99
83) Chrysene	21.994	228	550759	43.961	ng	100
84) Bis(2-ethylhexyl)phtha...	21.800	149	306179	47.549	ng	98
85) Di-n-octyl phthalate	23.093	149	511276	47.358	ng	96
87) Indeno(1,2,3-cd)pyrene	29.425	276	662805	44.053	ng	# 97
88) Benzo(b)fluoranthene	24.309	252	613159m	47.857	ng	
89) Benzo(k)fluoranthene	24.379	252	572044	45.196	ng	97
90) Benzo(a)pyrene	25.254	252	583861	53.946	ng	99
91) Dibenzo(a,h)anthracene	29.496	278	569913	45.853	ng	97
92) Benzo(g,h,i)perylene	30.688	276	563691	46.213	ng	97

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

