

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011923\
 Data File : BG056351.D
 Acq On : 19 Jan 2023 17:43
 Operator : CG/JU
 Sample : 01172-02MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-FMS

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 01/20/2023
 Supervised By : Jagrut Upadhyay 01/20/2023

Quant Time: Jan 20 01:35:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 19 00:31:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.307	152	39035	20.000	ng	0.00
21) Naphthalene-d8	11.133	136	142162	20.000	ng	0.00
39) Acenaphthene-d10	14.917	164	110745	20.000	ng	0.00
64) Phenanthrene-d10	17.655	188	302189	20.000	ng	0.00
76) Chrysene-d12	21.950	240	304703	20.000	ng	# 0.00
86) Perylene-d12	25.393	264	337116	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.828	112	289266	132.313	ng	0.00
7) Phenol-d6	7.449	99	368575	109.779	ng	0.00
23) Nitrobenzene-d5	9.482	82	226889	81.637	ng	0.00
42) 2,4,6-Tribromophenol	16.404	330	216412	154.297	ng	0.00
45) 2-Fluorobiphenyl	13.548	172	723146	90.170	ng	0.00
79) Terphenyl-d14	20.228	244	1654050	97.225	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.648	88	37965	38.019	ng	# 76
3) Pyridine	4.071	79	92883	30.539	ng	# 93
4) n-Nitrosodimethylamine	3.965	42	23671	38.260	ng	# 94
6) Aniline	7.620	93	67408	15.310	ng	99
8) 2-Chlorophenol	7.866	128	107768	49.537	ng	95
9) Benzaldehyde	7.432	77	46525	23.306	ng	96
10) Phenol	7.479	94	126655	37.198	ng	96
11) bis(2-Chloroethyl)ether	7.708	93	98118	42.523	ng	94
12) 1,3-Dichlorobenzene	8.195	146	127109m	48.007	ng	
13) 1,4-Dichlorobenzene	8.342	146	124490	46.821	ng	98
14) 1,2-Dichlorobenzene	8.666	146	120913	47.131	ng	96
15) Benzyl Alcohol	8.536	79	98615	40.358	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.818	45	19851	48.684	ng	# 72
17) 2-Methylphenol	8.742	107	100764	40.331	ng	95
18) Hexachloroethane	9.400	117	36868	45.645	ng	99
19) n-Nitroso-di-n-propyla...	9.106	70	72030	35.281	ng	98
20) 3+4-Methylphenols	9.071	107	136655	38.904	ng	97
22) Acetophenone	9.130	105	169939	42.277	ng	# 99
24) Nitrobenzene	9.523	77	109848	39.883	ng	93
25) Isophorone	10.040	82	220112	37.642	ng	98
26) 2-Nitrophenol	10.240	139	66907	47.479	ng	98
27) 2,4-Dimethylphenol	10.287	122	114875	44.694	ng	98
28) bis(2-Chloroethoxy)met...	10.516	93	124848	41.079	ng	98
29) 2,4-Dichlorophenol	10.781	162	133407	46.948	ng	99
30) 1,2,4-Trichlorobenzene	10.992	180	142229	44.107	ng	97
31) Naphthalene	11.186	128	325726	43.536	ng	99
32) Benzoic acid	10.416	122	55081m	28.883	ng	
33) 4-Chloroaniline	11.298	127	18004	5.040	ng	# 89
34) Hexachlorobutadiene	11.462	225	99037	42.703	ng	97
35) Caprolactam	12.056	113	33995m	35.335	ng	
36) 4-Chloro-3-methylphenol	12.391	107	122959	42.644	ng	99
37) 2-Methylnaphthalene	12.773	142	255914	40.510	ng	98
38) 1-Methylnaphthalene	12.984	142	240362	40.973	ng	97
40) 1,2,4,5-Tetrachloroben...	13.131	216	187476	45.167	ng	99
41) Hexachlorocyclopentadiene	13.107	237	223129	94.290	ng	96
43) 2,4,6-Trichlorophenol	13.366	196	133676	48.705	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.436	196	144080	44.377	ng	97
46) 1,1'-Biphenyl	13.754	154	378874	47.770	ng	99
47) 2-Chloronaphthalene	13.807	162	298820	47.733	ng	98
48) 2-Nitroaniline	14.000	65	63404	47.536	ng	93
49) Acenaphthylene	14.647	152	472056	46.337	ng	99
50) Dimethylphthalate	14.359	163	414085	46.318	ng	98
51) 2,6-Dinitrotoluene	14.482	165	89437	46.950	ng	94
52) Acenaphthene	14.982	154	345957m	52.214	ng	
53) 3-Nitroaniline	14.817	138	31273	17.250	ng	99
54) 2,4-Dinitrophenol	15.023	184	117232	86.144	ng	96
55) Dibenzofuran	15.317	168	497981	45.418	ng	98
56) 4-Nitrophenol	15.117	139	131767	94.502	ng	92
57) 2,4-Dinitrotoluene	15.270	165	131965	49.534	ng	90
58) Fluorene	15.957	166	402875	45.683	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.534	232	144726m	47.475	ng	
60) Diethylphthalate	15.710	149	380197	45.348	ng	99
61) 4-Chlorophenyl-phenylether	15.939	204	246105	44.718	ng	99
62) 4-Nitroaniline	15.975	138	87929	47.348	ng	91
63) Azobenzene	16.233	77	274233	42.387	ng	97
65) 4,6-Dinitro-2-methylph...	16.033	198	88295	43.368	ng	96
66) n-Nitrosodiphenylamine	16.157	169	377320	45.003	ng	96
67) 4-Bromophenyl-phenylether	16.832	248	171999	44.697	ng	98
68) Hexachlorobenzene	16.962	284	175923	49.001	ng	99
69) Atrazine	17.091	200	173975	50.068	ng	99
70) Pentachlorophenol	17.308	266	232499	84.304	ng	97
71) Phenanthrene	17.696	178	733276	46.824	ng	99
72) Anthracene	17.790	178	746703	46.222	ng	99
73) Carbazole	18.055	167	678001	49.929	ng	98
74) Di-n-butylphthalate	18.583	149	677962	50.084	ng	100
75) Fluoranthene	19.688	202	961788	47.083	ng	99
77) Benzidine	19.852	184	96313	9.020	ng	98
78) Pyrene	20.052	202	992860	46.229	ng	99
80) Butylbenzylphthalate	20.910	149	295540	47.966	ng	95
81) Benzo(a)anthracene	21.932	228	988785	46.202	ng	99
82) 3,3'-Dichlorobenzidine	21.827	252	135693	17.613	ng	97
83) Chrysene	22.003	228	922819	46.033	ng	99
84) Bis(2-ethylhexyl)phtha...	21.791	149	427839	48.197	ng	99
85) Di-n-octyl phthalate	23.072	149	717846	47.902	ng	100
87) Indeno(1,2,3-cd)pyrene	29.359	276	1165313	47.530	ng	# 97
88) Benzo(b)fluoranthene	24.294	252	1040696	48.587	ng	99
89) Benzo(k)fluoranthene	24.365	252	1005673	47.150	ng	99
90) Benzo(a)pyrene	25.228	252	910018	43.560	ng	99
91) Dibenzo(a,h)anthracene	29.418	278	978709	47.614	ng	99
92) Benzo(g,h,i)perylene	30.605	276	941451	47.230	ng	99

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

