

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021523\
 Data File : BG056644.D
 Acq On : 15 Feb 2023 22:41
 Operator : CG/JU
 Sample : PB150774BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB150774BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/16/2023
 Supervised By : Jagrut Upadhyay 02/16/2023

Quant Time: Feb 16 01:55:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG021523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 16 01:33:33 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.437	152	13275	20.000	ng	0.00
21) Naphthalene-d8	11.275	136	57651	20.000	ng	0.00
39) Acenaphthene-d10	15.041	164	47098	20.000	ng	0.00
64) Phenanthrene-d10	17.773	188	120531	20.000	ng	0.00
76) Chrysene-d12	22.092	240	116270	20.000	ng	0.00
86) Perylene-d12	25.635	264	122491	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.940	112	121300	148.673	ng	0.00
7) Phenol-d6	7.556	99	174919	145.139	ng	0.00
23) Nitrobenzene-d5	9.612	82	100543	89.295	ng	0.00
42) 2,4,6-Tribromophenol	16.516	330	91439	141.803	ng	0.00
45) 2-Fluorobiphenyl	13.672	172	282285	80.257	ng	0.00
79) Terphenyl-d14	20.347	244	617496	86.892	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.749	88	12104	33.839	ng	# 92
3) Pyridine	4.172	79	34733	30.698	ng	97
4) n-Nitrosodimethylamine	4.078	42	21636	41.025	ng	# 84
6) Aniline	7.744	93	56881	33.960	ng	99
8) 2-Chlorophenol	7.991	128	39173	48.527	ng	96
9) Benzaldehyde	7.556	77	17347	27.217	ng	98
10) Phenol	7.585	94	60413	49.004	ng	99
11) bis(2-Chloroethyl)ether	7.832	93	39229	42.410	ng	99
12) 1,3-Dichlorobenzene	8.326	146	40991	42.620	ng	95
13) 1,4-Dichlorobenzene	8.472	146	41769	43.123	ng	96
14) 1,2-Dichlorobenzene	8.796	146	41192	44.244	ng	93
15) Benzyl Alcohol	8.660	79	49817	46.164	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.960	45	54020	43.499	ng	98
17) 2-Methylphenol	8.860	107	38509	44.449	ng	98
18) Hexachloroethane	9.542	117	14562	47.460	ng	95
19) n-Nitroso-di-n-propyla...	9.236	70	39593	43.883	ng	98
20) 3+4-Methylphenols	9.183	107	53192	44.695	ng	93
22) Acetophenone	9.260	105	70032	39.757	ng	# 97
24) Nitrobenzene	9.653	77	52736	40.647	ng	95
25) Isophorone	10.176	82	108592	39.628	ng	98
26) 2-Nitrophenol	10.370	139	20794	44.851	ng	92
27) 2,4-Dimethylphenol	10.411	122	46673	44.054	ng	98
28) bis(2-Chloroethoxy)met...	10.652	93	57023	40.832	ng	97
29) 2,4-Dichlorophenol	10.905	162	48076	45.785	ng	100
30) 1,2,4-Trichlorobenzene	11.134	180	51848	41.451	ng	98
31) Naphthalene	11.328	128	127500	39.904	ng	100
32) Benzoic acid	10.523	122	29882	43.761	ng	97
33) 4-Chloroaniline	11.416	127	36137	24.755	ng	97
34) Hexachlorobutadiene	11.604	225	36152	40.444	ng	96
35) Caprolactam	12.174	113	15883m	47.029	ng	
36) 4-Chloro-3-methylphenol	12.497	107	53552	46.247	ng	99
37) 2-Methylnaphthalene	12.897	142	97916	38.812	ng	99
38) 1-Methylnaphthalene	13.114	142	92254	39.189	ng	98
40) 1,2,4,5-Tetrachloroben...	13.255	216	69655	39.784	ng	99
41) Hexachlorocyclopentadiene	13.237	237	85314	89.839	ng	99
43) 2,4,6-Trichlorophenol	13.478	196	48332	45.498	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.549	196	53201	41.786	ng	99
46) 1,1'-Biphenyl	13.884	154	144451	40.488	ng	99
47) 2-Chloronaphthalene	13.931	162	113235	40.107	ng	96
48) 2-Nitroaniline	14.119	65	33203	41.205	ng	91
49) Acenaphthylene	14.771	152	182444	39.770	ng	100
50) Dimethylphthalate	14.477	163	161339	41.439	ng	99
51) 2,6-Dinitrotoluene	14.601	165	30200	40.282	ng	90
52) Acenaphthene	15.106	154	125613	44.391	ng	100
53) 3-Nitroaniline	14.930	138	24127	31.924	ng	98
54) 2,4-Dinitrophenol	15.129	184	40556	90.164	ng	# 61
55) Dibenzofuran	15.435	168	190209	39.788	ng	99
56) 4-Nitrophenol	15.206	139	53624	92.608	ng	85
57) 2,4-Dinitrotoluene	15.382	165	42816	45.091	ng	93
58) Fluorene	16.081	166	153284	40.553	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.646	232	52025m	46.624	ng	
60) Diethylphthalate	15.829	149	162332	41.333	ng	98
61) 4-Chlorophenyl-phenyle...	16.064	204	92069	39.871	ng	99
62) 4-Nitroaniline	16.087	138	32720	41.541	ng	92
63) Azobenzene	16.357	77	158079	40.217	ng	96
65) 4,6-Dinitro-2-methylph...	16.134	198	26204	42.100	ng	98
66) n-Nitrosodiphenylamine	16.275	169	141370	40.317	ng	99
67) 4-Bromophenyl-phenylether	16.957	248	63365	40.947	ng	95
68) Hexachlorobenzene	17.080	284	71029	45.051	ng	98
69) Atrazine	17.203	200	61451	45.640	ng	98
70) Pentachlorophenol	17.415	266	86263	82.618	ng	93
71) Phenanthrene	17.820	178	275096	41.951	ng	99
72) Anthracene	17.908	178	275875	41.115	ng	100
73) Carbazole	18.167	167	241519	41.755	ng	98
74) Di-n-butylphthalate	18.702	149	272508	42.099	ng	99
75) Fluoranthene	19.800	202	349077	40.344	ng	98
77) Benzidine	19.965	184	124630	44.816	ng	97
78) Pyrene	20.159	202	353195	40.223	ng	99
80) Butylbenzylphthalate	21.028	149	113168	42.918	ng	98
81) Benzo(a)anthracene	22.074	228	352573	40.776	ng	99
82) 3,3'-Dichlorobenzidine	21.968	252	98326	35.049	ng	98
83) Chrysene	22.145	228	322645	39.915	ng	99
84) Bis(2-ethylhexyl)phtha...	21.945	149	167165	42.307	ng	99
85) Di-n-octyl phthalate	23.273	149	284970	43.392	ng	99
87) Indeno(1,2,3-cd)pyrene	29.724	276	427404	44.553	ng	# 99
88) Benzo(b)fluoranthene	24.507	252	371379	46.408	ng	98
89) Benzo(k)fluoranthene	24.577	252	352944	44.850	ng	98
90) Benzo(a)pyrene	25.470	252	318643	41.424	ng	99
91) Dibenzo(a,h)anthracene	29.800	278	360706	45.399	ng	99
92) Benzo(g,h,i)perylene	31.017	276	347024	44.699	ng	97

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

