

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051322\
 Data File : BG053558.D
 Acq On : 14 May 2022 00:57
 Operator : CG/JU
 Sample : N2829-01MSD 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-02-051222MSD

Quant Time: May 14 03:15:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 11 06:00:46 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 05/16/2022
 Supervised By : Jagrut Upadhyay 05/16/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.088	152	42340	20.000	ng	0.00
21) Naphthalene-d8	10.896	136	162828	20.000	ng	0.00
39) Acenaphthene-d10	14.714	164	110677	20.000	ng	0.00
64) Phenanthrene-d10	17.458	188	229887	20.000	ng	0.00
76) Chrysene-d12	21.740	240	206293	20.000	ng	0.00
86) Perylene-d12	25.030	264	209497	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.662	112	28327	11.354	ng	0.01
7) Phenol-d6	7.248	99	43845	11.580	ng	0.00
23) Nitrobenzene-d5	9.245	82	29347	7.672	ng	0.00
42) 2,4,6-Tribromophenol	16.200	330	18397	11.270	ng	0.00
45) 2-Fluorobiphenyl	13.334	172	68644	9.063	ng	0.00
79) Terphenyl-d14	20.060	244	96876	8.746	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.535	88	6912	5.280	ng	# 84
3) Pyridine	3.941	79	16303m	5.100	ng	
4) n-Nitrosodimethylamine	3.852	42	10978	7.050	ng	# 96
6) Aniline	7.406	93	26151	6.236	ng	97
8) 2-Chlorophenol	7.659	128	22604	8.667	ng	95
9) Benzaldehyde	7.224	77	14751m	6.133	ng	
10) Phenol	7.283	94	30618	8.268	ng	95
11) bis(2-Chloroethyl)ether	7.495	93	21287	7.708	ng	97
12) 1,3-Dichlorobenzene	7.976	146	23917	7.773	ng	96
13) 1,4-Dichlorobenzene	8.123	146	24674	7.947	ng	96
14) 1,2-Dichlorobenzene	8.446	146	24505	8.405	ng	92
15) Benzyl Alcohol	8.323	79	24137m	7.625	ng	
16) 2,2'-oxybis(1-Chloropr...	8.617	45	36275	8.070	ng	100
17) 2-Methylphenol	8.528	107	20023	8.000	ng	94
18) Hexachloroethane	9.169	117	8977	7.614	ng	99
19) n-Nitroso-di-n-propyla...	8.881	70	20440	7.524	ng	90
20) 3+4-Methylphenols	8.857	107	26912	7.613	ng	97
22) Acetophenone	8.904	105	35633	7.996	ng	# 97
24) Nitrobenzene	9.292	77	31784	8.585	ng	99
25) Isophorone	9.809	82	54908	8.529	ng	99
26) 2-Nitrophenol	10.009	139	12231	8.157	ng	98
27) 2,4-Dimethylphenol	10.068	122	21177	9.385	ng	95
28) bis(2-Chloroethoxy)met...	10.285	93	29701	8.022	ng	94
29) 2,4-Dichlorophenol	10.555	162	23575	8.656	ng	95
30) 1,2,4-Trichlorobenzene	10.761	180	26818	8.667	ng	88
31) Naphthalene	10.949	128	70657	8.531	ng	97
32) Benzoic acid	10.226	122	8888m	13.673	ng	
33) 4-Chloroaniline	11.054	127	20178	5.817	ng	99
34) Hexachlorobutadiene	11.231	225	19515	8.505	ng	99
35) Caprolactam	11.806	113	6757	6.827	ng	# 91
36) 4-Chloro-3-methylphenol	12.176	107	26320	8.079	ng	96
37) 2-Methylnaphthalene	12.547	142	50071	8.101	ng	98
38) 1-Methylnaphthalene	12.764	142	48245	8.000	ng	98
40) 1,2,4,5-Tetrachloroben...	12.917	216	33324	9.552	ng	98
41) Hexachlorocyclopentadiene	12.893	237	1974	10.072	ng	93
43) 2,4,6-Trichlorophenol	13.157	196	20173	8.859	ng	95

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051322\
 Data File : BG053558.D
 Acq On : 14 May 2022 00:57
 Operator : CG/JU
 Sample : N2829-01MSD 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-02-051222MSD

Quant Time: May 14 03:15:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 11 06:00:46 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/16/2022
 Supervised By : Jagrut Upadhyay 05/16/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.234	196	21807	8.998	ng	96
46) 1,1'-Biphenyl	13.545	154	68126	9.030	ng	99
47) 2-Chloronaphthalene	13.592	162	55132	9.328	ng	96
48) 2-Nitroaniline	13.792	65	18565	8.429	ng	94
49) Acenaphthylene	14.438	152	86079	9.126	ng	99
50) Dimethylphthalate	14.156	163	66324	7.886	ng	96
51) 2,6-Dinitrotoluene	14.279	165	13784	8.044	ng	97
52) Acenaphthene	14.779	154	49111	8.737	ng	100
53) 3-Nitroaniline	14.614	138	11989	6.476	ng	89
54) 2,4-Dinitrophenol	14.843	184	1099	7.176	ng	# 75
55) Dibenzofuran	15.108	168	84196	8.339	ng	98
56) 4-Nitrophenol	14.943	139	18925	14.312	ng	97
57) 2,4-Dinitrotoluene	15.073	165	18354	7.390	ng	# 86
58) Fluorene	15.760	166	67417	8.321	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.343	232	18450	7.503	ng	# 98
60) Diethylphthalate	15.513	149	66484	7.653	ng	97
61) 4-Chlorophenyl-phenyle...	15.742	204	34879	7.715	ng	96
62) 4-Nitroaniline	15.772	138	14022	7.302	ng	94
63) Azobenzene	16.036	77	67129	7.546	ng	98
65) 4,6-Dinitro-2-methylph...	15.842	198	1713	1.251	ng	85
66) n-Nitrosodiphenylamine	15.960	169	56708	9.395	ng	95
67) 4-Bromophenyl-phenylether	16.641	248	23794	8.862	ng	96
68) Hexachlorobenzene	16.770	284	26401	8.979	ng	95
69) Atrazine	16.899	200	22568	8.882	ng	96
70) Pentachlorophenol	17.117	266	22226m	17.870	ng	
71) Phenanthrene	17.499	178	124817	10.678	ng	97
72) Anthracene	17.593	178	110332	9.603	ng	98
73) Carbazole	17.857	167	92578	8.222	ng	97
74) Di-n-butylphthalate	18.403	149	103542	8.176	ng	97
75) Fluoranthene	19.508	202	159490	10.537	ng	99
77) Benzidine	19.684	184	27518m	6.189	ng	
78) Pyrene	19.872	202	154858	11.596	ng	99
80) Butylbenzylphthalate	20.741	149	42314	8.641	ng	92
81) Benzo(a)anthracene	21.722	228	128422	9.760	ng	99
82) 3,3'-Dichlorobenzidine	21.628	252	31923	9.567	ng	99
83) Chrysene	21.787	228	117814	9.587	ng	97
84) Bis(2-ethylhexyl)phtha...	21.605	149	60297	8.921	ng	# 99
85) Di-n-octyl phthalate	22.833	149	100824	8.822	ng	# 93
87) Indeno(1,2,3-cd)pyrene	28.789	276	95301	6.531	ng	# 95
88) Benzo(b)fluoranthene	23.978	252	123864	9.989	ng	99
89) Benzo(k)fluoranthene	24.049	252	112290	9.215	ng	98
90) Benzo(a)pyrene	24.865	252	115161	10.911	ng	# 99
91) Dibenzo(a,h)anthracene	28.830	278	82073	6.826	ng	98
92) Benzo(g,h,i)perylene	29.964	276	74444	6.247	ng	96

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

