

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062622\
 Data File : BG054075.D
 Acq On : 26 Jun 2022 16:37
 Operator : CG/JU
 Sample : N3488-12MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-7-16-18-20220623MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/27/2022
 Supervised By : Jagrut Upadhyay 06/27/2022

Supervised By : Jagrut
 Upadhyay
 06/27/2022

Quant Time: Jun 27 00:29:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 00:27:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.073	152	25736	20.000	ng	0.00
21) Naphthalene-d8	10.881	136	105394	20.000	ng	0.00
39) Acenaphthene-d10	14.695	164	76555	20.000	ng	0.00
64) Phenanthrene-d10	17.438	188	197870	20.000	ng	0.00
76) Chrysene-d12	21.716	240	216304	20.000	ng	0.00
86) Perylene-d12	24.965	264	226643	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.640	112	213222	158.401	ng	0.00
7) Phenol-d6	7.233	99	281291	154.042	ng	0.00
23) Nitrobenzene-d5	9.230	82	181924	97.002	ng	0.00
42) 2,4,6-Tribromophenol	16.181	330	152804	124.232	ng	0.00
45) 2-Fluorobiphenyl	13.320	172	454469	86.325	ng	0.00
79) Terphenyl-d14	20.047	244	1000647	93.174	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.525	88	25648	43.587	ng	92
3) Pyridine	3.919	79	65264	41.563	ng	93
4) n-Nitrosodimethylamine	3.837	42	39356	56.980	ng	# 94
6) Aniline	7.391	93	54627	25.232	ng	96
8) 2-Chlorophenol	7.638	128	77217	53.724	ng	94
9) Benzaldehyde	7.203	77	52892	49.078	ng	# 78
10) Phenol	7.256	94	102492	55.382	ng	93
11) bis(2-Chloroethyl)ether	7.485	93	73791	51.949	ng	96
12) 1,3-Dichlorobenzene	7.961	146	85622	47.542	ng	91
13) 1,4-Dichlorobenzene	8.108	146	86370	47.742	ng	98
14) 1,2-Dichlorobenzene	8.425	146	83779	48.241	ng	97
15) Benzyl Alcohol	8.308	79	73134	53.586	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.596	45	120471	56.734	ng	99
17) 2-Methylphenol	8.514	107	66675	51.864	ng	90
18) Hexachloroethane	9.160	117	29678	49.451	ng	# 84
19) n-Nitroso-di-n-propyla...	8.872	70	59896	48.823	ng	95
20) 3+4-Methylphenols	8.837	107	92642	51.385	ng	96
22) Acetophenone	8.890	105	117233	45.761	ng	# 98
24) Nitrobenzene	9.272	77	88161	47.736	ng	# 88
25) Isophorone	9.794	82	164374	46.928	ng	# 94
26) 2-Nitrophenol	9.988	139	42250	51.794	ng	# 83
27) 2,4-Dimethylphenol	10.047	122	73302	55.519	ng	91
28) bis(2-Chloroethoxy)met...	10.276	93	99177	52.560	ng	# 96
29) 2,4-Dichlorophenol	10.529	162	83897	47.393	ng	93
30) 1,2,4-Trichlorobenzene	10.740	180	87550	39.942	ng	98
31) Naphthalene	10.928	128	233537	43.893	ng	99
32) Benzoic acid	10.206	122	31353m	52.977	ng	
33) 4-Chloroaniline	11.034	127	25021	11.336	ng	# 91
34) Hexachlorobutadiene	11.222	225	63168	37.484	ng	95
35) Caprolactam	11.792	113	25665m	52.665	ng	
36) 4-Chloro-3-methylphenol	12.150	107	83543	48.548	ng	87
37) 2-Methylnaphthalene	12.532	142	172233	43.753	ng	96
38) 1-Methylnaphthalene	12.750	142	166882	43.539	ng	97
40) 1,2,4,5-Tetrachloroben...	12.897	216	115843	39.271	ng	98
41) Hexachlorocyclopentadiene	12.879	237	144065	104.883	ng	98
43) 2,4,6-Trichlorophenol	13.132	196	74551	44.992	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062622\
 Data File : BG054075.D
 Acq On : 26 Jun 2022 16:37
 Operator : CG/JU
 Sample : N3488-12MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-7-16-18-20220623MS

Quant Time: Jun 27 00:29:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 00:27:54 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 06/27/2022
 Supervised By :Jagrut Upadhyay 06/27/2022

Supervised By :Jagrut
 Upadhyay
 06/27/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.208	196	84013	45.409	ng	96
46) 1,1'-Biphenyl	13.531	154	244526	45.856	ng	99
47) 2-Chloronaphthalene	13.578	162	190554	44.155	ng	96
48) 2-Nitroaniline	13.772	65	60816	52.976	ng	94
49) Acenaphthylene	14.418	152	307524	45.826	ng	99
50) Dimethylphthalate	14.148	163	259895	44.576	ng	99
51) 2,6-Dinitrotoluene	14.260	165	57358	48.693	ng	# 80
52) Acenaphthene	14.759	154	224298m	52.436	ng	
53) 3-Nitroaniline	14.595	138	21662	18.981	ng	92
54) 2,4-Dinitrophenol	14.800	184	70213	105.553	ng	# 80
55) Dibenzofuran	15.094	168	305480	44.325	ng	96
56) 4-Nitrophenol	14.906	139	95982	118.462	ng	# 86
57) 2,4-Dinitrotoluene	15.053	165	82535	49.611	ng	93
58) Fluorene	15.740	166	254492	44.974	ng	94
59) 2,3,4,6-Tetrachlorophenol	15.323	232	78199	43.220	ng	# 94
60) Diethylphthalate	15.505	149	259858	46.023	ng	98
61) 4-Chlorophenyl-phenyle...	15.729	204	143973	42.499	ng	93
62) 4-Nitroaniline	15.758	138	57301	48.146	ng	# 73
63) Azobenzene	16.022	77	229749	50.551	ng	93
65) 4,6-Dinitro-2-methylph...	15.817	198	49539	50.182	ng	92
66) n-Nitrosodiphenylamine	15.946	169	232137	45.015	ng	97
67) 4-Bromophenyl-phenylether	16.628	248	100568	40.694	ng	94
68) Hexachlorobenzene	16.751	284	109626	38.780	ng	96
69) Atrazine	16.892	200	106450	49.830	ng	94
70) Pentachlorophenol	17.092	266	124902	95.117	ng	98
71) Phenanthrene	17.485	178	453145	44.575	ng	97
72) Anthracene	17.574	178	463375	46.677	ng	100
73) Carbazole	17.838	167	424372	49.803	ng	99
74) Di-n-butylphthalate	18.396	149	492054	49.944	ng	# 98
75) Fluoranthene	19.489	202	597807	44.796	ng	96
77) Benzidine	19.671	184	260794	58.220	ng	99
78) Pyrene	19.853	202	602510	46.119	ng	97
80) Butylbenzylphthalate	20.735	149	215969	51.575	ng	94
81) Benzo(a)anthracene	21.698	228	616833	44.043	ng	98
82) 3,3'-Dichlorobenzidine	21.610	252	165668	39.592	ng	97
83) Chrysene	21.763	228	580437	43.479	ng	97
84) Bis(2-ethylhexyl)phtha...	21.604	149	316992	52.929	ng	95
85) Di-n-octyl phthalate	22.826	149	560697	54.470	ng	98
87) Indeno(1,2,3-cd)pyrene	28.696	276	734795	42.962	ng	# 94
88) Benzo(b)fluoranthene	23.937	252	652883	47.377	ng	99
89) Benzo(k)fluoranthene	24.007	252	625208m	45.823	ng	
90) Benzo(a)pyrene	24.818	252	577636	49.681	ng	# 98
91) Dibenzo(a,h)anthracene	28.755	278	633548	44.795	ng	97
92) Benzo(g,h,i)perylene	29.859	276	630536	45.175	ng	94

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

