

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010323\
 Data File : BG056182.D
 Acq On : 3 Jan 2023 12:07
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 01/04/2023
 Supervised By :Yogesh Patel 01/05/2023

Quant Time: Jan 04 02:11:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 04 02:08:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.338	152	49950	20.000	ng	0.00
21) Naphthalene-d8	11.170	136	192663	20.000	ng	# 0.00
39) Acenaphthene-d10	14.948	164	161410	20.000	ng	0.00
64) Phenanthrene-d10	17.686	188	404357	20.000	ng	0.00
76) Chrysene-d12	21.987	240	368698	20.000	ng	0.00
86) Perylene-d12	25.448	264	394831	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.847	112	231857	82.879	ng	0.00
7) Phenol-d6	7.469	99	353490	82.279	ng	0.00
23) Nitrobenzene-d5	9.507	82	302027	80.187	ng	0.00
42) 2,4,6-Tribromophenol	16.429	330	161422	78.965	ng	0.00
45) 2-Fluorobiphenyl	13.579	172	932755	79.799	ng	0.00
79) Terphenyl-d14	20.259	244	1631879	79.273	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.673	88	51590	40.374	ng	# 90
3) Pyridine	4.090	79	160960	41.358	ng	# 93
4) n-Nitrosodimethylamine	3.996	42	32071	40.510	ng	85
6) Aniline	7.645	93	232272	41.227	ng	99
8) 2-Chlorophenol	7.892	128	119460	42.912	ng	95
9) Benzaldehyde	7.463	77	97038	37.987	ng	94
10) Phenol	7.492	94	177485	40.736	ng	97
11) bis(2-Chloroethyl)ether	7.739	93	120861	40.934	ng	95
12) 1,3-Dichlorobenzene	8.227	146	138089	40.757	ng	96
13) 1,4-Dichlorobenzene	8.373	146	141434	41.570	ng	97
14) 1,2-Dichlorobenzene	8.697	146	133711	40.730	ng	99
15) Benzyl Alcohol	8.567	79	124634	39.860	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.861	45	24077	46.145	ng	# 79
17) 2-Methylphenol	8.761	107	130225	40.733	ng	97
18) Hexachloroethane	9.431	117	41988	40.624	ng	87
19) n-Nitroso-di-n-propyla...	9.137	70	104095	39.845	ng	98
20) 3+4-Methylphenols	9.096	107	185838	41.345	ng	97
22) Acetophenone	9.161	105	217906	40.001	ng	# 99
24) Nitrobenzene	9.554	77	149096	39.944	ng	98
25) Isophorone	10.071	82	313768	39.593	ng	97
26) 2-Nitrophenol	10.265	139	77220	40.434	ng	96
27) 2,4-Dimethylphenol	10.312	122	138332	39.713	ng	98
28) bis(2-Chloroethoxy)met...	10.547	93	167395	40.641	ng	97
29) 2,4-Dichlorophenol	10.806	162	152863	39.694	ng	97
30) 1,2,4-Trichlorobenzene	11.029	180	169818	38.859	ng	97
31) Naphthalene	11.223	128	416571	41.084	ng	99
32) Benzoic acid	10.430	122	104531	40.445	ng	93
33) 4-Chloroaniline	11.317	127	198654	41.035	ng	97
34) Hexachlorobutadiene	11.499	225	121171	38.552	ng	96
35) Caprolactam	12.081	113	51847	39.765	ng	92
36) 4-Chloro-3-methylphenol	12.410	107	158620	40.592	ng	99
37) 2-Methylnaphthalene	12.804	142	336125	39.260	ng	99
38) 1-Methylnaphthalene	13.015	142	317828	39.977	ng	95
40) 1,2,4,5-Tetrachloroben...	13.162	216	239018	39.510	ng	99
41) Hexachlorocyclopentadiene	13.138	237	134057	38.868	ng	98
43) 2,4,6-Trichlorophenol	13.391	196	159551	39.885	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010323\
 Data File : BG056182.D
 Acq On : 3 Jan 2023 12:07
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 01/04/2023
 Supervised By :Yogesh Patel 01/05/2023

Quant Time: Jan 04 02:11:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 04 02:08:15 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.462	196	189285	40.001	ng	98
46) 1,1'-Biphenyl	13.791	154	474370	41.037	ng	98
47) 2-Chloronaphthalene	13.838	162	370251	40.579	ng	98
48) 2-Nitroaniline	14.026	65	81565	41.957	ng	98
49) Acenaphthylene	14.678	152	600751	40.460	ng	100
50) Dimethylphthalate	14.390	163	518570	39.798	ng	98
51) 2,6-Dinitrotoluene	14.513	165	112774	40.618	ng	95
52) Acenaphthene	15.013	154	388836	40.265	ng	98
53) 3-Nitroaniline	14.842	138	109060	41.274	ng	100
54) 2,4-Dinitrophenol	15.048	184	77070	38.856	ng	97
55) Dibenzofuran	15.348	168	636948	39.857	ng	96
56) 4-Nitrophenol	15.130	139	82988	40.836	ng	90
57) 2,4-Dinitrotoluene	15.295	165	159994	41.205	ng	91
58) Fluorene	15.988	166	504772	39.271	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.559	232	173672	39.088	ng	95
60) Diethylphthalate	15.735	149	482454	39.482	ng	99
61) 4-Chlorophenyl-phenyle...	15.970	204	313150	39.039	ng	98
62) 4-Nitroaniline	16.000	138	113486	41.928	ng	88
63) Azobenzene	16.264	77	376366	39.913	ng	97
65) 4,6-Dinitro-2-methylph...	16.053	198	113091	41.512	ng	95
66) n-Nitrosodiphenylamine	16.182	169	464468	41.400	ng	98
67) 4-Bromophenyl-phenylether	16.863	248	207073	40.216	ng	97
68) Hexachlorobenzene	16.993	284	196909	40.989	ng	99
69) Atrazine	17.116	200	186043	40.013	ng	98
70) Pentachlorophenol	17.328	266	152900	41.433	ng	94
71) Phenanthrene	17.727	178	844606	40.306	ng	99
72) Anthracene	17.821	178	873362	40.402	ng	100
73) Carbazole	18.080	167	729027	40.122	ng	98
74) Di-n-butylphthalate	18.614	149	743390	41.042	ng	99
75) Fluoranthene	19.719	202	1051765	38.478	ng	99
77) Benzidine	19.883	184	487407	37.723	ng	98
78) Pyrene	20.077	202	1059609	40.774	ng	99
80) Butylbenzylphthalate	20.935	149	317931	42.643	ng	99
81) Benzo(a)anthracene	21.963	228	1036837	40.038	ng	99
82) 3,3'-Dichlorobenzidine	21.863	252	365456	39.203	ng	97
83) Chrysene	22.034	228	967631	39.891	ng	99
84) Bis(2-ethylhexyl)phtha...	21.828	149	449322	41.831	ng	99
85) Di-n-octyl phthalate	23.121	149	762311	42.040	ng	100
87) Indeno(1,2,3-cd)pyrene	29.437	276	1174117	40.889	ng	# 97
88) Benzo(b)fluoranthene	24.343	252	1024163m	40.826	ng	
89) Benzo(k)fluoranthene	24.413	252	1011155	40.477	ng	100
90) Benzo(a)pyrene	25.283	252	998884	40.824	ng	# 99
91) Dibenzo(a,h)anthracene	29.502	278	971488	40.354	ng	99
92) Benzo(g,h,i)perylene	30.688	276	951410	40.753	ng	97

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

Manual Integrations

Quant Time: Jan 04 02:11:14 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 04 02:08:15 2023
Response via : Initial Calibration

