

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
 Data File : BG055467.D
 Acq On : 5 Nov 2022 12:28
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 11/07/2022
 Supervised By : Jagrut Upadhyay 11/07/2022

Quant Time: Nov 07 00:52:37 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Nov 07 00:51:08 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	31097	20.000	ng	0.00
21) Naphthalene-d8	11.056	136	133841	20.000	ng	# 0.00
39) Acenaphthene-d10	14.846	164	98323	20.000	ng	0.00
64) Phenanthrene-d10	17.584	188	236879	20.000	ng	0.00
76) Chrysene-d12	21.855	240	230283	20.000	ng	0.00
86) Perylene-d12	25.222	264	258450	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.750	112	131507	75.945	ng	0.00
7) Phenol-d6	7.378	99	184883	76.993	ng	0.00
23) Nitrobenzene-d5	9.405	82	188998	75.258	ng	0.00
42) 2,4,6-Tribromophenol	16.332	330	111284	82.977	ng	0.00
45) 2-Fluorobiphenyl	13.471	172	500011	75.399	ng	0.00
79) Terphenyl-d14	20.157	244	877112	74.064	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.553	88	27849	35.458	ng	98
3) Pyridine	3.976	79	85569m	36.711	ng	
4) n-Nitrosodimethylamine	3.888	42	35626	38.690	ng	98
6) Aniline	7.542	93	115876	37.206	ng	98
8) 2-Chlorophenol	7.789	128	78928	37.515	ng	98
9) Benzaldehyde	7.360	77	51708	36.627	ng	98
10) Phenol	7.407	94	103483	38.074	ng	98
11) bis(2-Chloroethyl)ether	7.636	93	74762	36.766	ng	97
12) 1,3-Dichlorobenzene	8.118	146	88748	37.215	ng	99
13) 1,4-Dichlorobenzene	8.265	146	90767	37.371	ng	98
14) 1,2-Dichlorobenzene	8.588	146	87520	37.646	ng	99
15) Benzyl Alcohol	8.471	79	73996	38.637	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.753	45	92927	34.349	ng	98
17) 2-Methylphenol	8.671	107	74325	38.006	ng	98
18) Hexachloroethane	9.323	117	30667	37.012	ng	94
19) n-Nitroso-di-n-propyla...	9.035	70	68017	36.249	ng	96
20) 3+4-Methylphenols	9.005	107	106267	38.244	ng	98
22) Acetophenone	9.058	105	131610	37.516	ng	# 98
24) Nitrobenzene	9.446	77	97560	37.261	ng	98
25) Isophorone	9.969	82	184131	37.022	ng	98
26) 2-Nitrophenol	10.163	139	51830	38.444	ng	98
27) 2,4-Dimethylphenol	10.216	122	72427	38.314	ng	100
28) bis(2-Chloroethoxy)met...	10.445	93	97058	36.638	ng	99
29) 2,4-Dichlorophenol	10.709	162	90690	39.328	ng	97
30) 1,2,4-Trichlorobenzene	10.915	180	96120	38.493	ng	98
31) Naphthalene	11.109	128	286250	37.986	ng	100
32) Benzoic acid	10.392	122	42275	39.087	ng	98
33) 4-Chloroaniline	11.215	127	122118	39.001	ng	98
34) Hexachlorobutadiene	11.385	225	59143	38.104	ng	100
35) Caprolactam	11.996	113	33306	39.010	ng	94
36) 4-Chloro-3-methylphenol	12.325	107	99143	39.266	ng	99
37) 2-Methylnaphthalene	12.695	142	212787	38.443	ng	99
38) 1-Methylnaphthalene	12.913	142	206668	38.281	ng	100
40) 1,2,4,5-Tetrachloroben...	13.059	216	111141	38.266	ng	100
41) Hexachlorocyclopentadiene	13.030	237	45210	36.456	ng	95
43) 2,4,6-Trichlorophenol	13.294	196	80052	39.578	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110522\
 Data File : BG055467.D
 Acq On : 5 Nov 2022 12:28
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/07/2022
 Supervised By : Jagrut Upadhyay 11/07/2022

Quant Time: Nov 07 00:52:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 00:51:08 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.377	196	88777	39.405	ng	99
46) 1,1'-Biphenyl	13.682	154	274183	37.897	ng	99
47) 2-Chloronaphthalene	13.735	162	224150	38.557	ng	99
48) 2-Nitroaniline	13.935	65	69030	38.659	ng	99
49) Acenaphthylene	14.575	152	370182	38.780	ng	100
50) Dimethylphthalate	14.293	163	319284	38.903	ng	99
51) 2,6-Dinitrotoluene	14.423	165	68990	40.153	ng	99
52) Acenaphthene	14.910	154	218675	38.490	ng	98
53) 3-Nitroaniline	14.752	138	77461	40.112	ng	100
54) 2,4-Dinitrophenol	14.969	184	35975	37.566	ng	96
55) Dibenzofuran	15.245	168	366938	39.020	ng	99
56) 4-Nitrophenol	15.063	139	53560	40.450	ng	97
57) 2,4-Dinitrotoluene	15.204	165	100225	40.796	ng	99
58) Fluorene	15.886	166	299703	39.232	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.468	232	80969	39.970	ng	98
60) Diethylphthalate	15.639	149	328942	38.567	ng	99
61) 4-Chlorophenyl-phenylether	15.868	204	154070	39.010	ng	99
62) 4-Nitroaniline	15.915	138	81583	39.931	ng	98
63) Azobenzene	16.162	77	266996	37.659	ng	98
65) 4,6-Dinitro-2-methylph...	15.968	198	63259	41.459	ng	99
66) n-Nitrosodiphenylamine	16.085	169	267943	38.046	ng	99
67) 4-Bromophenyl-phenylether	16.761	248	107210	38.287	ng	98
68) Hexachlorobenzene	16.890	284	124399	38.957	ng	94
69) Atrazine	17.020	200	88087	37.256	ng	99
70) Pentachlorophenol	17.237	266	57947	37.378	ng	98
71) Phenanthrene	17.625	178	507403	38.081	ng	100
72) Anthracene	17.713	178	498121	38.141	ng	99
73) Carbazole	17.983	167	460833	37.926	ng	100
74) Di-n-butylphthalate	18.506	149	571791	37.678	ng	100
75) Fluoranthene	19.616	202	607066	37.878	ng	99
77) Benzidine	19.787	184	183455	38.265	ng	98
78) Pyrene	19.975	202	614520	37.827	ng	99
80) Butylbenzylphthalate	20.827	149	259899	38.373	ng	99
81) Benzo(a)anthracene	21.832	228	612011	38.715	ng	100
82) 3,3'-Dichlorobenzidine	21.738	252	216239	39.890	ng	98
83) Chrysene	21.902	228	580995	38.524	ng	99
84) Bis(2-ethylhexyl)phtha...	21.691	149	374676	38.056	ng	100
85) Di-n-octyl phthalate	22.942	149	641171	38.009	ng	99
87) Indeno(1,2,3-cd)pyrene	29.111	276	792835	38.208	ng	# 91
88) Benzo(b)fluoranthene	24.146	252	643336	39.188	ng	99
89) Benzo(k)fluoranthene	24.217	252	628616	38.103	ng	99
90) Benzo(a)pyrene	25.063	252	547210	38.549	ng	99
91) Dibenzo(a,h)anthracene	29.158	278	655737	38.303	ng	99
92) Benzo(g,h,i)perylene	30.322	276	650273	38.191	ng	99

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

SSTDCCC040

APPROVED

Reviewed By :Christian Giraldo 11/07/2022
Supervised By :Jaqrut Upadhyay 11/07/2022

Quant Time: Nov 07 00:52:37 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 07 00:51:08 2022
Response via : Initial Calibration

