

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101322\
 Data File : BG055159.D
 Acq On : 13 Oct 2022 11:28
 Operator : CG/JU
 Sample : PB148265BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB148265BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/14/2022
 Supervised By : Jagrut Upadhyay 10/14/2022

Quant Time: Oct 14 01:20:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 13 05:14:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.305	152	26872	20.000	ng	0.00
21) Naphthalene-d8	11.131	136	129666	20.000	ng	0.00
39) Acenaphthene-d10	14.915	164	109473	20.000	ng	0.00
64) Phenanthrene-d10	17.647	188	293885	20.000	ng	0.00
76) Chrysene-d12	21.930	240	283440	20.000	ng	0.00
86) Perylene-d12	25.349	264	297042	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.819	112	211824	154.223	ng	0.00
7) Phenol-d6	7.435	99	296440	145.903	ng	0.00
23) Nitrobenzene-d5	9.474	82	185375	76.663	ng	0.00
42) 2,4,6-Tribromophenol	16.389	330	203366	129.285	ng	0.00
45) 2-Fluorobiphenyl	13.546	172	491973	73.412	ng	0.00
79) Terphenyl-d14	20.220	244	1119208	79.527	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.634	88	18632	32.059	ng	95
3) Pyridine	4.057	79	54137	29.540	ng	97
4) n-Nitrosodimethylamine	3.963	42	36510	37.542	ng	98
6) Aniline	7.611	93	110649	42.411	ng	97
8) 2-Chlorophenol	7.858	128	77650	46.362	ng	99
9) Benzaldehyde	7.423	77	48982	35.847	ng	97
10) Phenol	7.465	94	104252	46.185	ng	99
11) bis(2-Chloroethyl)ether	7.705	93	66353	41.560	ng	97
12) 1,3-Dichlorobenzene	8.193	146	76048	39.066	ng	96
13) 1,4-Dichlorobenzene	8.340	146	78145	39.760	ng	99
14) 1,2-Dichlorobenzene	8.663	146	76928	40.872	ng	97
15) Benzyl Alcohol	8.534	79	81180	44.693	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.828	45	111464	41.763	ng	98
17) 2-Methylphenol	8.728	107	73093	45.006	ng	92
18) Hexachloroethane	9.403	117	26029	39.067	ng	99
19) n-Nitroso-di-n-propyla...	9.104	70	66619	40.628	ng	99
20) 3+4-Methylphenols	9.057	107	102370	43.932	ng	96
22) Acetophenone	9.127	105	124185	38.681	ng	# 100
24) Nitrobenzene	9.515	77	94064	37.601	ng	99
25) Isophorone	10.038	82	190798	40.631	ng	98
26) 2-Nitrophenol	10.232	139	46980	39.637	ng	99
27) 2,4-Dimethylphenol	10.273	122	86267	47.858	ng	98
28) bis(2-Chloroethoxy)met...	10.514	93	99548	42.776	ng	97
29) 2,4-Dichlorophenol	10.767	162	88510	41.084	ng	96
30) 1,2,4-Trichlorobenzene	10.990	180	83320	35.853	ng	99
31) Naphthalene	11.184	128	253874	37.229	ng	100
32) Benzoic acid	10.396	122	52977	35.526	ng	94
33) 4-Chloroaniline	11.278	127	101950	34.781	ng	96
34) Hexachlorobutadiene	11.460	225	52372	35.344	ng	98
35) Caprolactam	12.042	113	38088m	43.036	ng	
36) 4-Chloro-3-methylphenol	12.371	107	107103	44.085	ng	95
37) 2-Methylnaphthalene	12.764	142	194925	38.172	ng	97
38) 1-Methylnaphthalene	12.982	142	188895	37.914	ng	99
40) 1,2,4,5-Tetrachloroben...	13.123	216	110650	35.836	ng	99
41) Hexachlorocyclopentadiene	13.105	237	137289	88.986	ng	98
43) 2,4,6-Trichlorophenol	13.352	196	84331	37.577	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101322\
 Data File : BG055159.D
 Acq On : 13 Oct 2022 11:28
 Operator : CG/JU
 Sample : PB148265BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB148265BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/14/2022
 Supervised By : Jagrut Upadhyay 10/14/2022

Quant Time: Oct 14 01:20:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 13 05:14:13 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.422	196	94947	38.127	ng	99
46) 1,1'-Biphenyl	13.751	154	276557	38.051	ng	99
47) 2-Chloronaphthalene	13.798	162	215332	36.626	ng	100
48) 2-Nitroaniline	13.992	65	81627	38.917	ng	96
49) Acenaphthylene	14.638	152	370008	37.490	ng	99
50) Dimethylphthalate	14.356	163	335616	38.139	ng	99
51) 2,6-Dinitrotoluene	14.474	165	72499	40.035	ng	98
52) Acenaphthene	14.979	154	270299	42.336	ng	98
53) 3-Nitroaniline	14.803	138	70575	33.615	ng	98
54) 2,4-Dinitrophenol	15.009	184	92920	84.352	ng	91
55) Dibenzofuran	15.308	168	375571	37.704	ng	98
56) 4-Nitrophenol	15.097	139	139499	85.325	ng	95
57) 2,4-Dinitrotoluene	15.255	165	108041	40.734	ng	97
58) Fluorene	15.955	166	310842	38.005	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.526	232	92760	37.746	ng	98
60) Diethylphthalate	15.702	149	356431	38.454	ng	99
61) 4-Chlorophenyl-phenylether	15.937	204	168596	38.498	ng	100
62) 4-Nitroaniline	15.960	138	86009	37.945	ng	98
63) Azobenzene	16.231	77	304563	38.704	ng	98
65) 4,6-Dinitro-2-methylph...	16.013	198	66252	38.354	ng	98
66) n-Nitrosodiphenylamine	16.148	169	289368	37.323	ng	99
67) 4-Bromophenyl-phenylether	16.830	248	118196	36.893	ng	96
68) Hexachlorobenzene	16.948	284	134140	36.429	ng	99
69) Atrazine	17.083	200	131899	43.152	ng	99
70) Pentachlorophenol	17.288	266	168325	76.724	ng	96
71) Phenanthrene	17.688	178	560594	36.775	ng	99
72) Anthracene	17.776	178	568848	38.185	ng	100
73) Carbazole	18.040	167	537632	38.761	ng	99
74) Di-n-butylphthalate	18.581	149	649009	38.009	ng	99
75) Fluoranthene	19.674	202	711966	37.363	ng	98
77) Benzidine	19.844	184	426317	67.071	ng	99
78) Pyrene	20.038	202	704957	37.309	ng	98
80) Butylbenzylphthalate	20.902	149	283514	39.087	ng	99
81) Benzo(a)anthracene	21.906	228	677523	37.360	ng	99
82) 3,3'-Dichlorobenzidine	21.812	252	230910	37.102	ng	99
83) Chrysene	21.977	228	636951	37.233	ng	100
84) Bis(2-ethylhexyl)phtha...	21.789	149	400306	38.744	ng	100
85) Di-n-octyl phthalate	23.070	149	679138	39.426	ng	99
87) Indeno(1,2,3-cd)pyrene	29.286	276	829408	37.695	ng	98
88) Benzo(b)fluoranthene	24.257	252	727254	41.077	ng	99
89) Benzo(k)fluoranthene	24.333	252	707398	39.769	ng	99
90) Benzo(a)pyrene	25.191	252	639483	42.240	ng	98
91) Dibenzo(a,h)anthracene	29.357	278	724147	39.771	ng	99
92) Benzo(g,h,i)perylene	30.526	276	712601	39.953	ng	97

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

Manual Integrations

Quant Time: Oct 14 01:20:51 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 13 05:14:13 2022
Response via : Initial Calibration

