

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040822\
 Data File : BG053088.D
 Acq On : 8 Apr 2022 17:41
 Operator : CG/JU
 Sample : PB143883BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB143883BS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 04/11/2022
 Supervised By : Jagrut Upadhyay 04/11/2022

Quant Time: Apr 09 07:03:40 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Apr 08 15:31:24 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.124	152	29406	20.000	ng	0.00
21) Naphthalene-d8	10.932	136	121601	20.000	ng	# 0.00
39) Acenaphthene-d10	14.745	164	92296	20.000	ng	0.00
64) Phenanthrene-d10	17.488	188	222085	20.000	ng	0.00
76) Chrysene-d12	21.776	240	228243	20.000	ng	# 0.00
86) Perylene-d12	25.078	264	233024	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.687	112	215240	125.471	ng	0.00
7) Phenol-d6	7.284	99	315567	119.289	ng	0.00
23) Nitrobenzene-d5	9.282	82	227495	75.962	ng	0.00
42) 2,4,6-Tribromophenol	16.231	330	165895	112.977	ng	0.00
45) 2-Fluorobiphenyl	13.370	172	500506	74.776	ng	0.00
79) Terphenyl-d14	20.091	244	1028709	76.120	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.566	88	22683	27.688	ng	# 91
3) Pyridine	3.965	79	63389	29.338	ng	# 92
4) n-Nitrosodimethylamine	3.871	42	43400	40.562	ng	80
6) Aniline	7.443	93	97098	31.671	ng	99
8) 2-Chlorophenol	7.684	128	83500	45.087	ng	96
9) Benzaldehyde	7.255	77	48126m	30.435	ng	
10) Phenol	7.308	94	110343	41.904	ng	88
11) bis(2-Chloroethyl)ether	7.531	93	75917	39.995	ng	91
12) 1,3-Dichlorobenzene	8.013	146	84633	39.327	ng	95
13) 1,4-Dichlorobenzene	8.160	146	84471	38.502	ng	96
14) 1,2-Dichlorobenzene	8.477	146	84660	40.015	ng	95
15) Benzyl Alcohol	8.354	79	95400	40.572	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.641	45	126437	39.897	ng	99
17) 2-Methylphenol	8.565	107	76125	42.721	ng	92
18) Hexachloroethane	9.211	117	32067	38.977	ng	93
19) n-Nitroso-di-n-propyla...	8.923	70	76160	39.324	ng	# 93
20) 3+4-Methylphenols	8.888	107	103987	41.339	ng	89
22) Acetophenone	8.941	105	128958	38.226	ng	93
24) Nitrobenzene	9.329	77	116994	40.163	ng	93
25) Isophorone	9.851	82	209445	41.145	ng	98
26) 2-Nitrophenol	10.039	139	46769	38.395	ng	90
27) 2,4-Dimethylphenol	10.098	122	82343	47.201	ng	89
28) bis(2-Chloroethoxy)met...	10.327	93	109573	38.315	ng	98
29) 2,4-Dichlorophenol	10.580	162	90880	41.439	ng	99
30) 1,2,4-Trichlorobenzene	10.797	180	94239	37.966	ng	98
31) Naphthalene	10.985	128	253133	39.018	ng	99
32) Benzoic acid	10.239	122	42716	38.408	ng	90
33) 4-Chloroaniline	11.085	127	55511	19.981	ng	# 93
34) Hexachlorobutadiene	11.273	225	69896	37.908	ng	97
35) Caprolactam	11.855	113	30668m	39.672	ng	
36) 4-Chloro-3-methylphenol	12.201	107	101318	39.682	ng	92
37) 2-Methylnaphthalene	12.583	142	186313	38.811	ng	93
38) 1-Methylnaphthalene	12.800	142	179269m	38.257	ng	
40) 1,2,4,5-Tetrachloroben...	12.947	216	118110	37.678	ng	98
41) Hexachlorocyclopentadiene	12.930	237	158129	97.719	ng	93
43) 2,4,6-Trichlorophenol	13.182	196	83726	38.739	ng	93

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040822\
 Data File : BG053088.D
 Acq On : 8 Apr 2022 17:41
 Operator : CG/JU
 Sample : PB143883BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB143883BS

Quant Time: Apr 09 07:03:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 15:31:24 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/11/2022
 Supervised By : Jagrut Upadhyay 04/11/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.259	196	91503	39.731	ng	99
46) 1,1'-Biphenyl	13.582	154	255341	38.417	ng	98
47) 2-Chloronaphthalene	13.629	162	202156	38.123	ng	99
48) 2-Nitroaniline	13.823	65	79299	39.405	ng	87
49) Acenaphthylene	14.469	152	347217	41.828	ng	98
50) Dimethylphthalate	14.193	163	283527	38.407	ng	98
51) 2,6-Dinitrotoluene	14.310	165	62346	40.483	ng	86
52) Acenaphthene	14.809	154	250094m	44.427	ng	
53) 3-Nitroaniline	14.639	138	41659	25.582	ng	# 88
54) 2,4-Dinitrophenol	14.851	184	79959	81.598	ng	# 86
55) Dibenzofuran	15.144	168	328510	37.269	ng	97
56) 4-Nitrophenol	14.950	139	101100	81.402	ng	# 75
57) 2,4-Dinitrotoluene	15.097	165	89231	40.697	ng	# 89
58) Fluorene	15.790	166	270352	37.975	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.368	232	84049	37.579	ng	96
60) Diethylphthalate	15.550	149	292606	37.830	ng	99
61) 4-Chlorophenyl-phenyle...	15.779	204	153523	38.041	ng	97
62) 4-Nitroaniline	15.802	138	62486	36.363	ng	# 82
63) Azobenzene	16.072	77	290251	37.220	ng	97
65) 4,6-Dinitro-2-methylph...	15.867	198	56928	36.741	ng	99
66) n-Nitrosodiphenylamine	15.990	169	246031	38.524	ng	97
67) 4-Bromophenyl-phenylether	16.672	248	107280	37.699	ng	95
68) Hexachlorobenzene	16.801	284	120156	38.990	ng	94
69) Atrazine	16.936	200	106276	42.601	ng	97
70) Pentachlorophenol	17.142	266	143123	84.977	ng	93
71) Phenanthrene	17.529	178	473525	39.035	ng	98
72) Anthracene	17.623	178	472463	39.314	ng	98
73) Carbazole	17.888	167	438522	37.538	ng	99
74) Di-n-butylphthalate	18.440	149	515730	38.926	ng	98
75) Fluoranthene	19.538	202	619628	39.849	ng	99
77) Benzidine	19.709	184	251467	38.615	ng	100
78) Pyrene	19.897	202	613198	37.723	ng	99
80) Butylbenzylphthalate	20.772	149	230833	38.089	ng	97
81) Benzo(a)anthracene	21.753	228	604708	38.749	ng	97
82) 3,3'-Dichlorobenzidine	21.659	252	147967	25.782	ng	99
83) Chrysene	21.823	228	546515	37.735	ng	99
84) Bis(2-ethylhexyl)phtha...	21.647	149	321807	39.284	ng	99
85) Di-n-octyl phthalate	22.887	149	532816	38.891	ng	97
87) Indeno(1,2,3-cd)pyrene	28.867	276	663578	38.330	ng	# 98
88) Benzo(b)fluoranthene	24.026	252	612893	41.686	ng	# 98
89) Benzo(k)fluoranthene	24.097	252	580057	40.143	ng	99
90) Benzo(a)pyrene	24.919	252	589661	47.078	ng	# 97
91) Dibenzo(a,h)anthracene	28.926	278	578891	40.627	ng	# 96
92) Benzo(g,h,i)perylene	30.042	276	566100	40.308	ng	100

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040822\
 Data File : BG053088.D
 Acq On : 8 Apr 2022 17:41
 Operator : CG/JU
 Sample : PB143883BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB143883BS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 04/11/2022
 Supervised By : Jagrut Upadhyay 04/11/2022

Quant Time: Apr 09 07:03:40 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Apr 08 15:31:24 2022

Response via : Initial Calibration

