

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012712.D
 Acq On : 23 Nov 2022 16:06
 Operator : CG/JU
 Sample : N5740-02MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-12MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/25/2022
 Supervised By : Jagrut Upadhyay 11/25/2022

Quant Time: Nov 24 02:29:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	343864	20.000	ng	0.00
21) Naphthalene-d8	10.852	136	1274444	20.000	ng	0.00
39) Acenaphthene-d10	14.669	164	733710	20.000	ng	0.00
64) Phenanthrene-d10	17.422	188	1555333	20.000	ng	0.00
76) Chrysene-d12	21.510	240	1752850	20.000	ng	0.00
86) Perylene-d12	24.033	264	1871678	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2710984	170.799	ng	0.00
7) Phenol-d6	7.181	99	3119522	137.044	ng	0.00
23) Nitrobenzene-d5	9.199	82	1903668	83.306	ng	0.00
42) 2,4,6-Tribromophenol	16.157	330	1300666	136.494	ng	0.00
45) 2-Fluorobiphenyl	13.293	172	4516020	87.689	ng	0.00
79) Terphenyl-d14	19.969	244	7881491	95.565	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.458	88	396644	59.299	ng	# 44
3) Pyridine	3.858	79	957923	45.074	ng	96
4) n-Nitrosodimethylamine	3.764	42	479758	51.619	ng	# 88
6) Aniline	7.352	93	1104097	38.368	ng	98
8) 2-Chlorophenol	7.593	128	1077049	52.609	ng	99
9) Benzaldehyde	7.164	77	500032	33.341	ng	98
10) Phenol	7.211	94	1177371	45.390	ng	97
11) bis(2-Chloroethyl)ether	7.452	93	1067730	47.143	ng	99
12) 1,3-Dichlorobenzene	7.922	146	1101937	43.314	ng	99
13) 1,4-Dichlorobenzene	8.069	146	1120227	42.873	ng	99
14) 1,2-Dichlorobenzene	8.387	146	1083881	42.693	ng	98
15) Benzyl Alcohol	8.269	79	842525m	52.060	ng	
16) 2,2'-oxybis(1-Chloropr...	8.569	45	1432806	42.692	ng	99
17) 2-Methylphenol	8.469	107	853252	48.162	ng	98
18) Hexachloroethane	9.122	117	370662	38.832	ng	94
19) n-Nitroso-di-n-propyla...	8.840	70	724241	45.074	ng	96
20) 3+4-Methylphenols	8.799	107	1108689	46.475	ng	98
22) Acetophenone	8.858	105	1552414	49.899	ng	98
24) Nitrobenzene	9.240	77	1074263	45.941	ng	98
25) Isophorone	9.769	82	2067809	49.351	ng	99
26) 2-Nitrophenol	9.957	139	450801	55.723	ng	98
27) 2,4-Dimethylphenol	10.010	122	961336	66.910	ng	97
28) bis(2-Chloroethoxy)met...	10.252	93	1293357	52.191	ng	98
29) 2,4-Dichlorophenol	10.487	162	962421	48.821	ng	98
30) 1,2,4-Trichlorobenzene	10.710	180	997537	43.577	ng	98
31) Naphthalene	10.899	128	3065040	44.313	ng	100
32) Benzoic acid	10.069	122	137399	24.578	ng	99
33) 4-Chloroaniline	11.005	127	311150	12.345	ng	98
34) Hexachlorobutadiene	11.187	225	599403	41.727	ng	99
35) Caprolactam	11.775	113	271180m	49.270	ng	
36) 4-Chloro-3-methylphenol	12.116	107	958201	50.025	ng	98
37) 2-Methylnaphthalene	12.504	142	2102847	42.676	ng	99
38) 1-Methylnaphthalene	12.722	142	1994674	41.227	ng	99
40) 1,2,4,5-Tetrachloroben...	12.863	216	1128940	48.338	ng	99
41) Hexachlorocyclopentadiene	12.851	237	991483	89.260	ng	98
43) 2,4,6-Trichlorophenol	13.099	196	770546	49.055	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012712.D
 Acq On : 23 Nov 2022 16:06
 Operator : CG/JU
 Sample : N5740-02MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-12MS

Quant Time : Nov 24 02:29:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/25/2022
 Supervised By : Jagrut Upadhyay 11/25/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.163	196	840991	55.638	ng	98
46) 1,1'-Biphenyl	13.504	154	2738904	49.346	ng	100
47) 2-Chloronaphthalene	13.546	162	2105459	47.902	ng	99
48) 2-Nitroaniline	13.746	65	451396	40.305	ng	96
49) Acenaphthylene	14.393	152	3318045	47.537	ng	99
50) Dimethylphthalate	14.122	163	2694395	49.908	ng	100
51) 2,6-Dinitrotoluene	14.240	165	514098	45.311	ng	96
52) Acenaphthene	14.734	154	2065672	46.697	ng	97
53) 3-Nitroaniline	14.563	138	385998	30.939	ng	# 99
54) 2,4-Dinitrophenol	14.763	184	135738	33.111	ng	95
55) Dibenzofuran	15.069	168	3214008	46.713	ng	100
56) 4-Nitrophenol	14.863	139	1000670	88.531	ng	97
57) 2,4-Dinitrotoluene	15.022	165	691564	41.549	ng	97
58) Fluorene	15.716	166	2566817	45.376	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.287	232	758104	51.259	ng	98
60) Diethylphthalate	15.487	149	2567184	50.281	ng	100
61) 4-Chlorophenyl-phenylether	15.710	204	1276946	45.469	ng	99
62) 4-Nitroaniline	15.728	138	557128	43.146	ng	98
63) Azobenzene	16.004	77	2386787	44.354	ng	96
65) 4,6-Dinitro-2-methylph...	15.781	198	177502	30.639	ng	99
66) n-Nitrosodiphenylamine	15.922	169	2210127	50.842	ng	99
67) 4-Bromophenyl-phenylether	16.610	248	790690	49.703	ng	98
68) Hexachlorobenzene	16.722	284	947345	48.194	ng	99
69) Atrazine	16.875	200	789948	57.474	ng	98
70) Pentachlorophenol	17.063	266	1258967	110.761	ng	100
71) Phenanthrene	17.469	178	4223910	48.010	ng	99
72) Anthracene	17.557	178	4241585	50.738	ng	99
73) Carbazole	17.822	167	4077812	53.991	ng	100
74) Di-n-butylphthalate	18.375	149	4635023	56.310	ng	98
75) Fluoranthene	19.422	202	5165723	49.769	ng	99
77) Benzidine	19.598	184	2976976	288.224	ng	99
78) Pyrene	19.775	202	5490495	47.444	ng	99
80) Butylbenzylphthalate	20.645	149	1484430	76.917	ng	95
81) Benzo(a)anthracene	21.492	228	5707050	48.176	ng	99
82) 3,3'-Dichlorobenzidine	21.422	252	1018899	37.393	ng	99
83) Chrysene	21.551	228	5381865	47.462	ng	99
84) Bis(2-ethylhexyl)phtha...	21.416	149	2643710	71.080	ng	99
85) Di-n-octyl phthalate	22.386	149	5004937	69.977	ng	# 87
87) Indeno(1,2,3-cd)pyrene	26.692	276	7159568	54.616	ng	100
88) Benzo(b)fluoranthene	23.263	252	6126394	48.917	ng	100
89) Benzo(k)fluoranthene	23.316	252	6019363	46.891	ng	100
90) Benzo(a)pyrene	23.921	252	5351422	53.598	ng	99
91) Dibenzo(a,h)anthracene	26.715	278	6159364	55.121	ng	100
92) Benzo(g,h,i)perylene	27.504	276	5943888	58.941	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012712.D
 Acq On : 23 Nov 2022 16:06
 Operator : CG/JU
 Sample : N5740-02MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :

BNA_P

ClientSampleId :

MH-12MS

Quant Time: Nov 24 02:29:26 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Nov 23 00:38:14 2022

Response via : Initial Calibration

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 11/25/2022
 Supervised By : Jagrut Upadhyay 11/25/2022

