Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110221\

Data File : BN017253.D

Acq On : 02 Nov 2021 23:25

Operator : CG/JU Sample : PB140353BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 02 23:58:24 2021

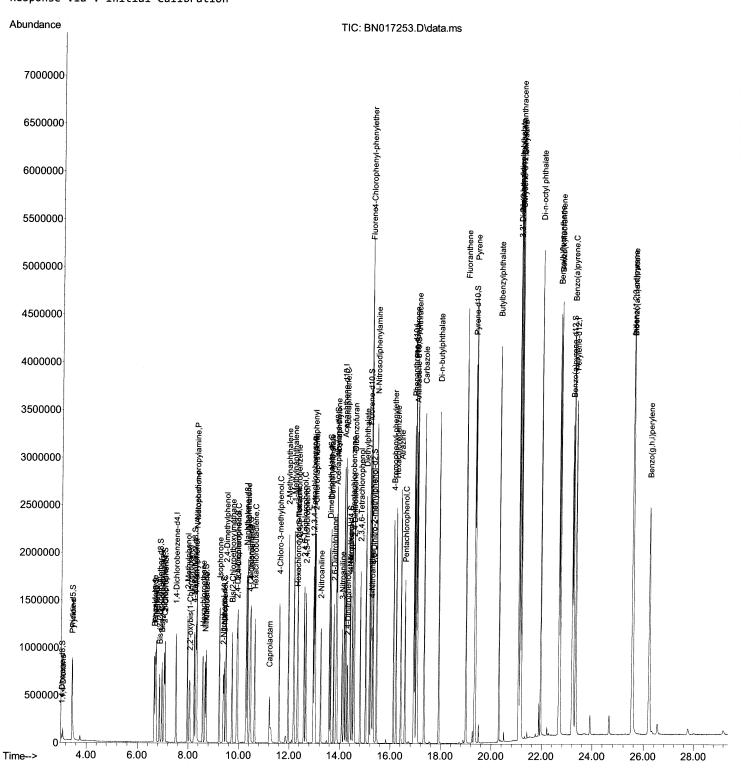
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 15:59:34 2021 Response via : Initial Calibration Instrument : BNA_N ClientSampleId : SLCS353

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/03/2021 Supervised By :mohammad ahmed 11/08/2021



SFAM-EPA-BN110221.M Wed Nov 03 00:13:24 2021

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110221\

Data File : BN017253.D

Acq On : 02 Nov 2021 23:25

Operator : CG/JU Sample : PB140353BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 02 23:58:24 2021

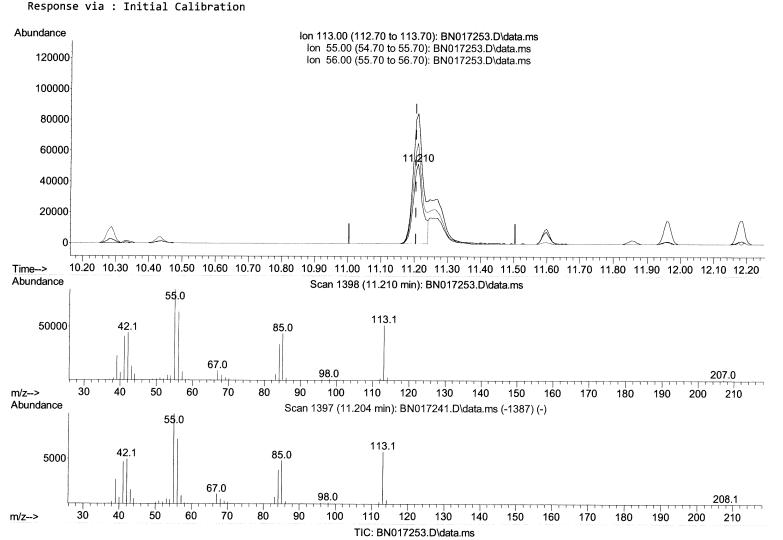
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 15:59:34 2021



Manual IntegrationsAPPROVED

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(34) Caprolactam

11.210min (+ 0.006) 12.54 ng/ul

response	101574	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	172.30	163.21
56.00	123.70	126.12
0.00	0.00	0.00

Quantitation Report (Qedit)

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Data File : BN017253.D

Acq On : 02 Nov 2021 23:25

Operator : CG/JU Sample : PB140353BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 02 23:58:24 2021

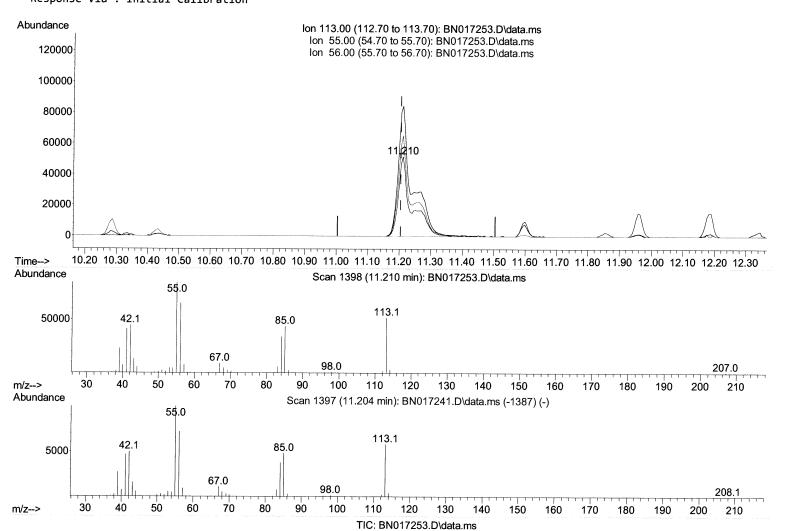
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 15:59:34 2021
Response via : Initial Calibration



Manual IntegrationsAPPROVED

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(34) Caprolactam

response	153412	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	172.30	163.21
56.00	123.70	126.12
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110221\

Data File : BN017253.D

Acq On : 02 Nov 2021 23:25

Operator : CG/JU Sample : PB140353BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 02 23:58:24 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M

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Response via : Initial Calibration

Instrument:
BNA_N
ClientSampleId:
SLCS353

Manual IntegrationsAPPROVED

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Compound		QIon	Response	Conc U	nits Dev	(Min)
Intonnal Ctandard						
Internal Standards	7 [11	150	206202	20.00		
1) 1,4-Dichlorobenzene-d4	7.511				ng/ul	0.00
20) Naphthalene-d8	10.287				ng/ul	0.00
38) Acenaphthene-d10	14.169				ng/ul	0.00
64) Phenanthrene-d10	16.922				ng/ul	0.00
79) Chrysene-d12	21.139				ng/ul	0.00
88) Perylene-d12	23.333	264	2257328	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.034	96	26485	2 522	3 ng/uL	0.00
4) Pyridine-d5	3.428	84	365319		_	
7) Phenol-d5	6.699	99	513597		ng/ul ng/ul	0.00 0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	6.864	67	299985		-	
11) 2-Chlorophenol-d4	7.046	132			ng/ul ng/ul	0.00
15) 4-Methylphenol-d8			413840			0.00
21) Nitrobenzene-d5	8.234	113	422658		ng/ul	0.00
	8.663	128	204221		ng/ul	0.00
24) 2-Nitrophenol-d4	9.381	143	225593		ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.916	165	421147		ng/ul	0.00
31) 4-Chloroaniline-d4	10.434	131	581656		ng/ul	0.00
46) Dimethylphthalate-d6	13.593	166	1301090		ng/ul	0.00
49) Acenaphthylene-d8	13.857	160	1632906		ng/ul	0.00
54) 4-Nitrophenol-d4	14.398	143	257335		ng/ul	0.00
60) Fluorene-d10	15.169	176	1132950		ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.304	200	238381		ng/ul	0.00
73) Anthracene-d10	17.022	188	1781319		ng/ul	0.00
81) Pyrene-d10	19.328	212	2125771	18.364	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.198	264	2217092	18.180	ng/ul	0.00
Target Compounds					0.45	1
2) 1,4-Dioxane	2 070		FF074	7 (14	Qva	
	3.070	88	55874		ng/uL	94
5) Pyridine	3.446	79	394981	18.536	-	99
6) Benzaldehyde	6.664	77	297329	21.421		99
8) Phenol	6.728	94	554630	20.015	-	98
10) Bis(2-Chloroethyl)ether	6.952	93	434163	19.771		100
12) 2-Chlorophenol	7.075	128	445718	20.062	-	99
13) 2-Methylphenol	7.963	108	426564	20.034		98
14) 2,2'-oxybis(1-Chloropr	8.052	45	640485	19.696	_	99
16) Acetophenone	8.334	105	689827	20.338		99
17) N-Nitroso-di-n-propyla	8.322	70	342088	19.801	-	99
18) 4-Methylphenol	8.293	108	479881	20.352		100
19) Hexachloroethane		117	173349	19.974		97
22) Nitrobenzene	8.705	77	485575	19.481	ng/ul	100
23) Isophorone	9.228	82	959110	19.357		99
25) 2-Nitrophenol	9.410	139	266663	20.002	ng/ul	97
26) 2,4-Dimethylphenol	9.487	107	519651	19.862	ng/ul	98
<pre>27) Bis(2-Chloroethoxy)met</pre>	9.722	93	619575	19.586	ng/ul	99
29) 2,4-Dichlorophenol	9.940	162	445656	20.258	ng/ul	98
30) Naphthalene	10.334	128	1511870	19.686	_	100
32) 4-Chloroaniline	10.457	127	650399	19.438		98
33) Hexachlorobutadiene	10.622	225	260928	19.420	ng/ul	99
34) Caprolactam	11.210	113	153412m>			ucielyoun
35) 4-Chloro-3-methylphenol	11.599	107	493643	20.009		99
					-	

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Misc

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QLast Update : Tue Nov 02 15:59:34 2021
Response via : Initial Calibration

Instrument:
BNA_N
ClientSampleId:
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/03/2021 Supervised By :mohammad ahmed 11/08/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	11.957	142	1048667	19.680 ng/ul	99
37) 1-Methylnaphthalene	12.181	142		19.288 ng/ul	96
39) 1,2,4,5-Tetrachloroben		216	532139	19.543 ng/ul	97
40) Hexachlorocyclopentadiene	12.316	237	303120	17.470 ng/ul	99
41) 2,4,6-Trichlorophenol	12.587	196	362596	20.049 ng/ul	99
42) 2,4,5-Trichlorophenol	12.657	196	391083	19.720 ng/ul	94
43) 1,1'-Biphenyl	12.998	154	1424665	19.926 ng/ul	99
44) 2-Chloronaphthalene	13.034	162	1068876	19.383 ng/ul	96
45) 2-Nitroaniline	13.251	65	313136	20.405 ng/ul	94
47) Dimethylphthalate	13.640	163	1392693	19.839 ng/ul	100
48) 2,6-Dinitrotoluene	13.763	165	286866	20.826 ng/ul	99
50) Acenaphthylene	13.887	152	1807762	19.831 ng/ul	98
51) 3-Nitroaniline	14.087	138	288590	19.937 ng/ul	98
52) Acenaphthene	14.234	153	1162495	19.677 ng/ul	99
53) 2,4-Dinitrophenol	14.298	184	169735	18.914 ng/ul	95
55) 4-Nitrophenol	14.410	109	194340	20.222 ng/ul	95
56) Dibenzofuran	14.569	168	1667580	19.893 ng/ul	98
57) 2,4-Dinitrotoluene	14.551	165	420202	20.618 ng/ul	96
58) 2,3,4,6-Tetrachlorophenol		232	33 0932	19.648 ng/ul	97
59) Diethylphthalate	15.022	149	1423785	20.000 ng/ul	100
61) Fluorene	15.222	166	1343253	20.059 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.228	204	649568	19.499 ng/ul	99
63) 4-Nitroaniline	15.257	138	313562	21.771 ng/ul	97
66) 4,6-Dinitro-2-methylph	15.316	198	254867	20.129 ng/ul	100
67) N-Nitrosodiphenylamine	15.445	169	1199045	20.319 ng/ul	98
68) 4-Bromophenyl-phenylether		248	416547	20.198 ng/ul	99
69) Hexachlorobenzene	16.228	284	484839	20.221 ng/ul	98
70) Atrazine	16.410	200	428758	19.792 ng/ul	97
71) Pentachlorophenol	16.581	266	286365	19.421 ng/ul	98
72) Phenanthrene	16.969	178	2200828	20.212 ng/ul	99
74) Anthracene	17.057	178	2220152	20.255 ng/ul	100
75) 1,2,3,4-Tetrachloroben	12.951	216	537544	19.244 ng/uL	94
76) Pentachlorobenzene	14.493	250	551136	18.984 ng/uL	97
77) Carbazole	17.334	167	2064172	21.015 ng/ul	98
78) Di-n-butylphthalate	17.922	149	2498641	21.117 ng/ul	99
80) Fluoranthene	18.992	202	2656593	19.490 ng/ul	96
82) Pyrene	19.357	202	2735131	19.591 ng/ul	100
83) Butylbenzylphthalate84) 3,3'-Dichlorobenzidine	20.292	149	1157702	19.937 ng/ul	98
· ·	21.069	252	936429	19.020 ng/ul	99
85) Benzo(a)anthracene86) Bis(2-ethylhexyl)phtha	21.122	228	2746812	19.582 ng/ul	97
87) Chrysene	21.080	149	1768039	19.978 ng/ul	99
89) Di-n-octyl phthalate	21.175 21.951	228	2651339	19.460 ng/ul	96
90) Benzo(b)fluoranthene	22.674	149 252	3062021	20.586 ng/ul	100
91) Benzo(k)fluoranthene			2988413	19.867 ng/ul	98
93) Benzo(a)pyrene	22.721 23.239	252 252	2822827 2875591	19.919 ng/ul	98 07
94) Indeno(1,2,3-cd)pyrene	25.545	276	3397499	19.676 ng/ul	97 97
95) Dibenzo(a,h)anthracene	25.568	278	2892881	19.723 ng/ul	97 99
96) Benzo(g,h,i)perylene	26.221	276	2788086	19.812 ng/ul 19.127 ng/ul	99 99
				_	

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