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

Data File : BN017250.D

: 02 Nov 2021 21:37

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 17 Sample Multiplier: 1

Quant Time: Nov 02 22:30:03 2021

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

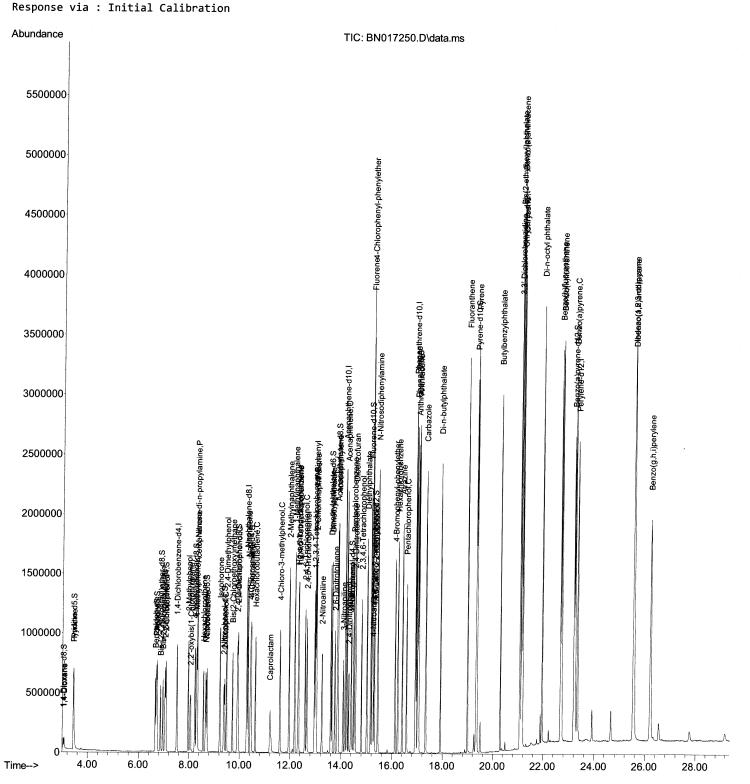
Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 15:59:34 2021

Instrument : BNA_N LabSampleId : SSTDCCC020

Manual IntegrationsAPPROVED

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



SFAM-EPA-BN110221.M Tue Nov 02 22:32:00 2021

Quantitation Report (Qedit)

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

Data File : BN017250.D

Acq On : 02 Nov 2021 21:37

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 17 Sample Multiplier: 1

Quant Time: Nov 02 22:30:03 2021

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

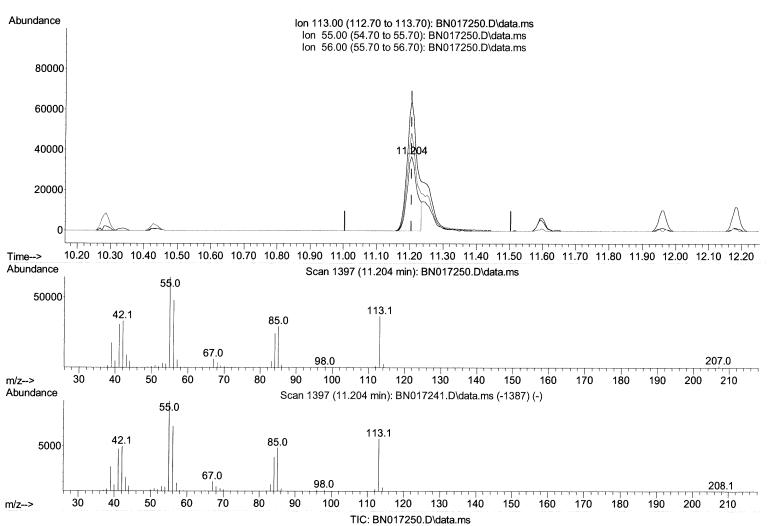
Quant Title : SVOA CALIBRATION

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



Manual Integrations APPROVED

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



(34) Caprolactam

11.204min (+ 0.000) 11.61 ng/ul

response	78028	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	172.30	175.54
56.00	123.70	131.67
0.00	0.00	0.00

Quantitation Report (Qedit)

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

Data File : BN017250.D

Acq On : 02 Nov 2021 21:37

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 17 Sample Multiplier: 1

Quant Time: Nov 02 22:30:03 2021

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

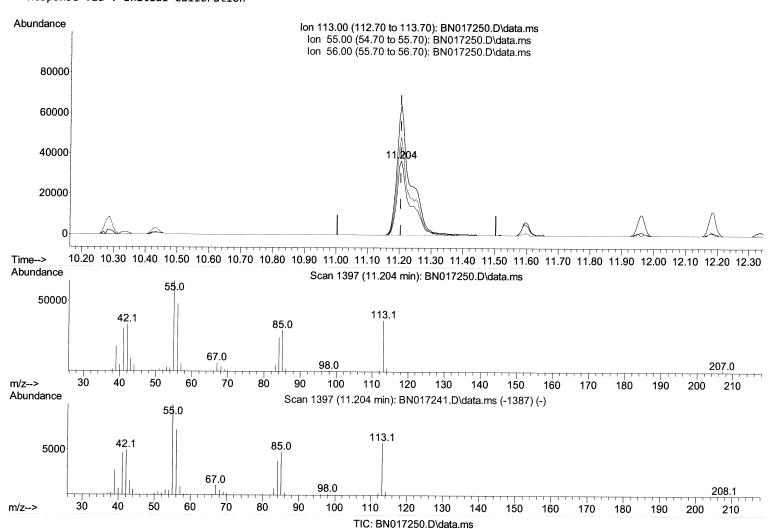
Quant Title : SVOA CALIBRATION

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



Manual IntegrationsAPPROVED

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



(34) Caprolactam

11.204min (+ 0.000) 16.22 ng/ul m WO4/2 JV

response	108992	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	172.30	175.54
56.00	123.70	131.67
0.00	0.00	0.00

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

Data File : BN017250.D

Acq On : 02 Nov 2021 21:37

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 17 Sample Multiplier: 1

Quant Time: Nov 02 22:30:03 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 **LabSampleld :** SSTDCCC020

Manual IntegrationsAPPROVED

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

Compound	R.T.		Response			
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.511	152	234762	20.000	ng/ul	0 00
20) Naphthalene-d8	10.287		1185517	20.000		
38) Acenaphthene-d10	14.169		795693	20.000		
64) Phenanthrene-d10	16.922		1663787	20.000		
79) Chrysene-d12	21.139		1701686		_	
88) Perylene-d12	23.333		1597877	20.000	-	
ob) refylene-ulz	23.333	204	139/6//	20.000	iig/ui	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.034	96	43101	7.256	ng/uL	0.00
4) Pyridine-d5	3.428	84	297570	17.954	-	
7) Phenol-d5	6.699	99	385531	17.791		
<pre>9) Bis-(2-Chloroethyl)eth</pre>	6.864	67	234851	18.157		
11) 2-Chlorophenol-d4	7.046	132	317825	18.457		
15) 4-Methylphenol-d8	8.228	113	322042	18.025	_	
21) Nitrobenzene-d5	8.663	128	159292	17.317		
24) 2-Nitrophenol-d4	9.381	143	177943	17.344	-	
28) 2,4-Dichlorophenol-d3	9.916	165	326030	17.674	-	
31) 4-Chloroaniline-d4	10.434	131	482735	17.462		
46) Dimethylphthalate-d6	13.593	166	997302	16.904	-	
49) Acenaphthylene-d8	13.857	160	1264921	17.046		
54) 4-Nitrophenol-d4	14.393	143	191400	16.529		0.00
60) Fluorene-d10	15.169	176	867112	17.224	-	
65) 4,6-Dinitro-2-methylph	15.298	200	175829	16.521	•	0.00
73) Anthracene-d10	17.022	188	1388418	17.683	_	0.00
81) Pyrene-d10	19.328	212	1615326	17.542		0.00
92) Benzo(a)pyrene-d12	23.192	264	1746694	20.234	-	0.00
					•	
Target Compounds					-	/alue
2) 1,4-Dioxane	3.070	88	42152	7.250	ng/uL	95
5) Pyridine	3.446	79	304312	18.024	ng/ul	99
Benzaldehyde	6.664	77	202313	18.396	ng/ul	99
8) Phenol	6.722	94	392266	17.866	ng/ul	99
10) Bis(2-Chloroethyl)ether	6.952	93	314656	18.084	ng/ul	100
<pre>12) 2-Chlorophenol</pre>	7.075	128	326287	18.536		98
<pre>13) 2-Methylphenol</pre>	7.963	108	301983	17.900	ng/ul	100
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.052	45	469009	18.203	ng/ul	100
16) Acetophenone	8.334	105	498788	18.560		99
17) N-Nitroso-di-n-propyla	8.322	70	249141	18.200	ng/ul	100
<pre>18) 4-Methylphenol</pre>	8.293	108	339825	18.189		100
19) Hexachloroethane	8.575	117	124305	18.077	ng/ul	94
22) Nitrobenzene	8.705	77	355033	17.170	ng/ul	99
23) Isophorone	9.228	82	697741	16.976	ng/ul	98
25) 2-Nitrophenol	9.410	139	191048	17.275	ng/ul	99
26) 2,4-Dimethylphenol	9.487	107	376361	17.341	ng/ul	99
<pre>27) Bis(2-Chloroethoxy)met</pre>	9.722	93	446851	17.029	ng/ul	99
<pre>29) 2,4-Dichlorophenol</pre>	9.940	162	322421	17.667	ng/ul	97
30) Naphthalene	10.334	128	1100048	17.267	_	100
32) 4-Chloroaniline	10.457	127	484231	17.445		100
33) Hexachlorobutadiene	10.622	225	194489	17.449		100
34) Caprolactam	11.204	113	108992m >	16.221	ng/ul	NETRIPOTO <
35) 4-Chloro-3-methylphenol	11.599	107	351041	17.152	ng/ul	100

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Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 17 Sample Multiplier: 1

Quant Time: Nov 02 22:30:03 2021

 $\label{lem:quant_Methods} Quant \ \mbox{Methods} \ \ \mbox{SFAM-EPA-BN110221.M}$

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 15:59:34 2021 Response via : Initial Calibration Instrument :
BNA_N
LabSampleId :
SSTDCCC020

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	11.963	142	760368	17.202 ng/ul	 95
37) 1-Methylnaphthalene	12.181	142		17.311 ng/ul	96
39) 1,2,4,5-Tetrachloroben	12.340	216	391709	17.189 ng/ul	97
40) Hexachlorocyclopentadiene	12.316	237		21.449 ng/ul	100
41) 2,4,6-Trichlorophenol	12.587	196	255529	16.883 ng/ul	99
42) 2,4,5-Trichlorophenol	12.657	196	279047	16.813 ng/ul	93
43) 1,1'-Biphenyl	12.999	154	1024221	17.117 ng/ul	99
44) 2-Chloronaphthalene	13.028	162	778475	16.868 ng/ul	99
45) 2-Nitroaniline	13.251	65	219039	17.055 ng/ul	96
47) Dimethylphthalate	13.640	163	989129	16.836 ng/ul	99
48) 2,6-Dinitrotoluene	13.763	165	197600	17.141 ng/ul	99
50) Acenaphthylene	13.887	152	1315339	17.241 ng/ul	100
51) 3-Nitroaniline	14.087	138	194019	16.016 ng/ul	98
52) Acenaphthene	14.234	153	849123	17.174 ng/ul	100
53) 2,4-Dinitrophenol	14.298	184	136685	18.200 ng/ul	94
55) 4-Nitrophenol	14.410	109	135876	16.894 ng/ul	95
56) Dibenzofuran	14.569	168	1211245	17.265 ng/ul	100
57) 2,4-Dinitrotoluene	14.551	165	294811	17.285 ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	14.804	232	235835	16.731 ng/ul	99
59) Diethylphthalate	15.022	149	994140	16.687 ng/ul	99
61) Fluorene	15.222	166	975537	17.407 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.228	204	474261	17.011 ng/ul	98
63) 4-Nitroaniline	15.257	138	198202	16.444 ng/ul	98
66) 4,6-Dinitro-2-methylph	15.316	198	177356	16.787 ng/ul	99
67) N-Nitrosodiphenylamine	15.445	169	848826	17.238 ng/ul	99
68) 4-Bromophenyl-phenylether	16.128	248	290370	16.874 ng/ul	98
69) Hexachlorobenzene	16.228	284	348204	17.404 ng/ul	98
70) Atrazine	16.410	200	308009	17.039 ng/ul	98
71) Pentachlorophenol	16.575	266	247029	20.078 ng/ul	96
72) Phenanthrene	16.963	178	1582473	17.417 ng/ul	99
74) Anthracene	17.057	178	1579450	17.269 ng/ul	99
75) 1,2,3,4-Tetrachloroben76) Pentachlorobenzene	12.957	216	397752	17.065 ng/uL	98
77) Carbazole	14.493	250	406158	16.766 ng/uL	97
78) Di-n-butylphthalate	17.334	167	1466029	17.887 ng/ul	99
80) Fluoranthene	17.916 18.992	149 202	1726750	17.490 ng/ul	100
82) Pyrene	19.357	202	1883803	17.374 ng/ul	97 100
83) Butylbenzylphthalate	20.292	149	1942712 799959	17.493 ng/ul 17.318 ng/ul	100
84) 3,3'-Dichlorobenzidine	21.069	252	690418	_	98 98
85) Benzo(a)anthracene	21.122	228	1983191	17.628 ng/ul 17.773 ng/ul	98
86) Bis(2-ethylhexyl)phtha	21.080	149	1256296	17.773 ng/ul 17.845 ng/ul	98
87) Chrysene	21.175	228	1927445	17.784 ng/ul	98
89) Di-n-octyl phthalate	21.951	149	2163165	20.545 ng/ul	100
90) Benzo(b)fluoranthene	22.674	252	2164180	20.325 ng/ul	97
91) Benzo(k)fluoranthene	22.722	252	2070878	20.644 ng/ul	98
93) Benzo(a)pyrene	23.239	252	2114837	20.443 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	25.545	276	2497018	20.478 ng/ul	97
95) Dibenzo(a,h)anthracene	25.563	278	2102218	20.338 ng/ul	98
96) Benzo(g,h,i)perylene	26.215	276	2103957	20.391 ng/ul	97

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