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

Data File : BN017265.D

Acq On : 03 Nov 2021 06:37

Operator : CG/JU

Sample : SSTDCCC020EC

Misc

ALS Vial : 32 Sample Multiplier: 1

Quant Time: Nov 03 10:03:36 2021

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

Quant Title : SVOA CALIBRATION

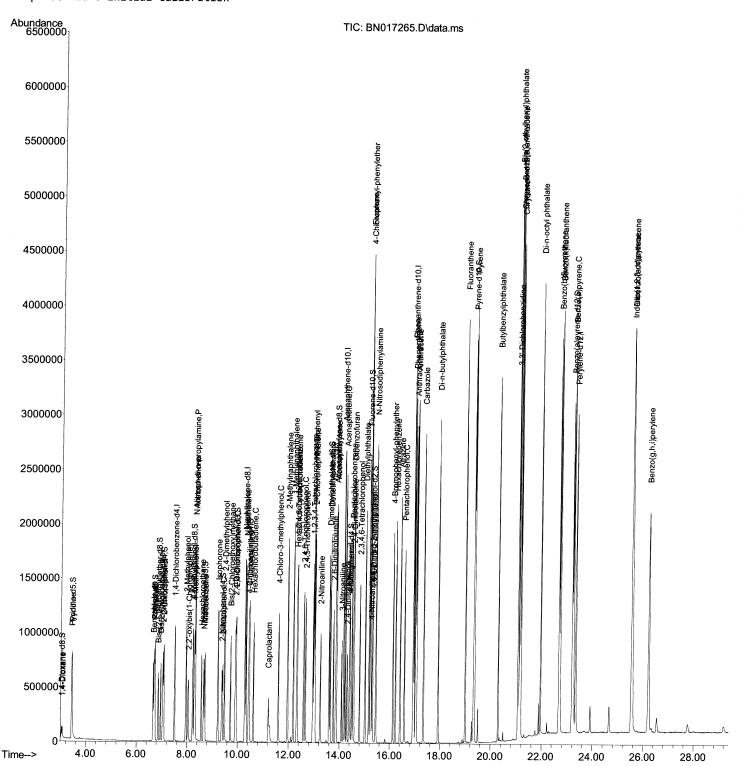
QLast Update : Tue Nov 02 15:59:34 2021

Response via : Initial Calibration



Manual IntegrationsAPPROVED

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



SFAM-EPA-BN110221.M Wed Nov 03 10:05:37 2021

Quantitation Report (Qedit)

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

Data File: BN017265.D

: 03 Nov 2021 06:37 Acq On

Operator : CG/JU Sample

: SSTDCCC020EC

Misc

ALS Vial : 32 Sample Multiplier: 1

Quant Time: Nov 03 10:03:36 2021

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

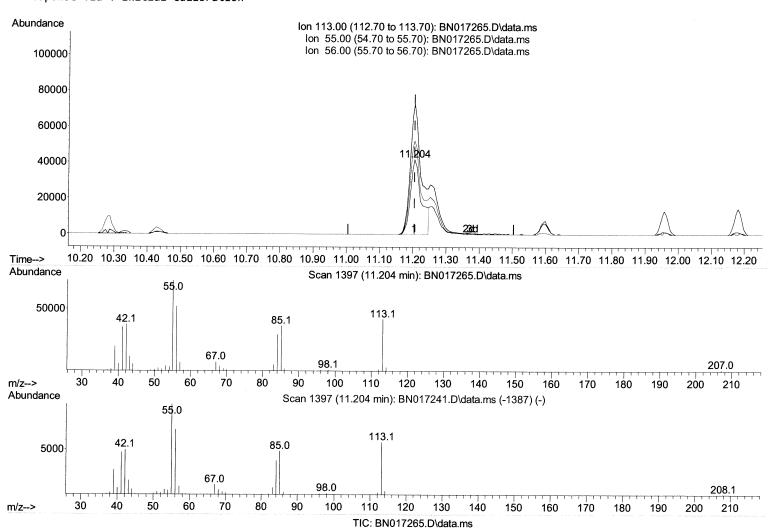
Quant Title : SVOA CALIBRATION

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



Caprolactam

11.204min (-0.000) 12.32 ng/ul

response	96021	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	172.30	174.47
56.00	123.70	125.20
0.00	0.00	0.00

Quantitation Report (Qedit)

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

Data File : BN017265.D

: 03 Nov 2021 06:37

Operator : CG/JU Sample

: SSTDCCC020EC

Misc

Sample Multiplier: 1 ALS Vial : 32

Quant Time: Nov 03 10:03:36 2021

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

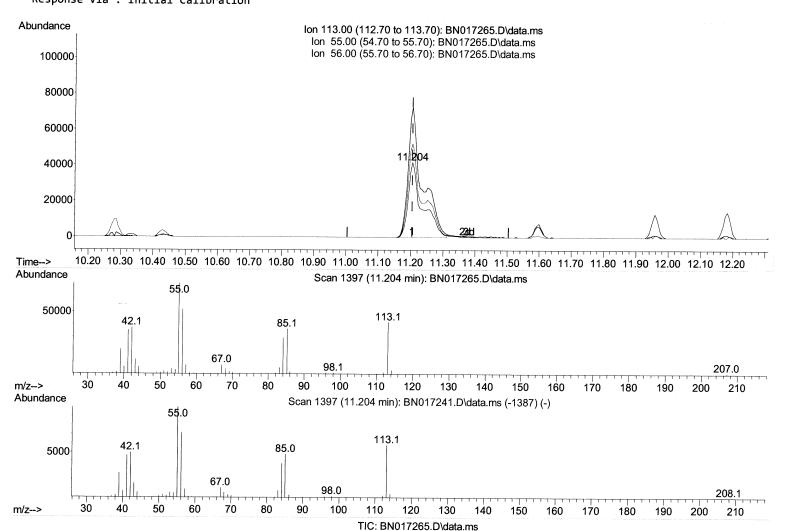
QLast Update : Tue Nov 02 15:59:34 2021 Response via : Initial Calibration

Quant Title : SVOA CALIBRATION



Manual Integrations APPROVED

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



(34) Caprolactam

11.204min (-0.000) 16.19 ng/ul m 1104/104

response	126249	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	172.30	174.47
56.00	123.70	125.20
0.00	0.00	0.00

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

Data File : BN017265.D

Acq On : 03 Nov 2021 06:37

Operator : CG/JU Sample : SSTDCCC020EC

Misc :

ALS Vial : 32 Sample Multiplier: 1

Quant Time: Nov 03 10:03:36 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
LabSampleId :
SSTDCCC020EC

Manual IntegrationsAPPROVED

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

Compound		QIon	Response	Conc Units I	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.510	152	271545	20.000 ng/u	ul 0.00
20) Naphthalene-d8	10.281		1375591	20.000 ng/u	
38) Acenaphthene-d10	14.169		924475	20.000 ng/u	
64) Phenanthrene-d10	16.922		1936944	20.000 ng/t	
79) Chrysene-d12	21.139		1898433	20.000 ng/t	
88) Perylene-d12	23.333		1824254	20.000 ng/t	
,,,	23.333	204	1024254	20.000 116/0	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.034	96	50728	7.383 ng/ι	ıL 0.00
4) Pyridine-d5	3.428	84	347384	18.121 ng/u	
7) Phenol-d5	6.699	99	452068	18.036 ng/u	
<pre>9) Bis-(2-Chloroethyl)eth</pre>	6.858	67	274269	18.332 ng/u	
11) 2-Chlorophenol-d4	7.046	132	368810	18.516 ng/u	
15) 4-Methylphenol-d8	8.228	113	376895	18.238 ng/u	
21) Nitrobenzene-d5	8.663	128	185223	17.354 ng/u	
24) 2-Nitrophenol-d4	9.381	143	210174	17.655 ng/u	
28) 2,4-Dichlorophenol-d3	9.916	165	380815	17.792 ng/u	
31) 4-Chloroaniline-d4	10.428	131	558101	17.732 ng/u	
46) Dimethylphthalate-d6	13.592	166	1134361	16.549 ng/u	
49) Acenaphthylene-d8	13.857	160	1493250	17.320 ng/u	
54) 4-Nitrophenol-d4	14.392	143		•	
60) Fluorene-d10	15.169	176	227468	16.907 ng/u	
65) 4,6-Dinitro-2-methylph	15.298		994642	17.005 ng/u	
73) Anthracene-d10		200	204710	16.522 ng/u	
81) Pyrene-d10	17.022	188 212	1591845	17.415 ng/u	
92) Benzo(a)pyrene-d12	19.328 23.198	264	1876206 2035197	18.264 ng/u 20.650 ng/u	
,(u/py/ elle u22	23.130	204	2033137	20.030 lig/u	1 0.00
Target Compounds					Qvalue
2) 1,4-Dioxane	3.070	88	49204	7.317 ng/u	L 94
5) Pyridine	3.446	79	354495	18.152 ng/u	
6) Benzaldehyde	6.658	77	236317	18.577 ng/u	1 96
8) Phenol	6.722	94	463485	18.250 ng/u	
<pre>10) Bis(2-Chloroethyl)ether</pre>	6.952	93	366812	18.226 ng/u	
12) 2-Chlorophenol	7.075	128	377616	18.546 ng/u	
13) 2-Methylphenol	7.963	108	354845	18.184 ng/u	
14) 2,2'-oxybis(1-Chloropr	8.052	45	543048	18.222 ng/u	
16) Acetophenone	8.328	105	583097	18.758 ng/u	
17) N-Nitroso-di-n-propyla	8.322	70	291774	18.427 ng/u	
18) 4-Methylphenol	8.293	108	397690	18.403 ng/u	
19) Hexachloroethane	8.569		145455	18.288 ng/u	
22) Nitrobenzene	8.705	77	412553	17.195 ng/u	
23) Isophorone	9.222	82	820133	17.196 ng/u	
25) 2-Nitrophenol	9.410	139	227021	17.691 ng/u	
26) 2,4-Dimethylphenol	9.487	107	434725	17.263 ng/u	
27) Bis(2-Chloroethoxy)met	9.722	93	519933	17.203 ng/u.	
29) 2,4-Dichlorophenol	9.940	162	370892	17.515 ng/u	
30) Naphthalene	10.334	128	1281995	17.342 ng/u	
32) 4-Chloroaniline	10.354	127	563624	17.542 ng/u.	
33) Hexachlorobutadiene	10.622	225	216625	•	
34) Caprolactam	11.204	113		16.750 ng/ul 16.193 ng/ul	
35) 4-Chloro-3-methylphenol	11.598	107			
55, 4 Chitor O-5-methyrphenor	11.330	TO	413031	17.392 ng/u]	L 99

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

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ALS Vial : 32 Sample Multiplier: 1

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

Instrument :
BNA_N
LabSampleId :
SSTDCCC020EC

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	883314	17.222 ng/ul	99
37) 1-Methylnaphthalene	12.181	142	915612	17.376 ng/ul	97
39) 1,2,4,5-Tetrachloroben	12.334	216	452311	17.084 ng/ul	97
40) Hexachlorocyclopentadiene	12.316	237	352363	20.886 ng/ul	99
41) 2,4,6-Trichlorophenol	12.587	196	302938	17.227 ng/ul	98
42) 2,4,5-Trichlorophenol	12.657	196	321475	16.671 ng/ul	96
43) 1,1'-Biphenyl	12.993	154	1193242	17.163 ng/ul	98
44) 2-Chloronaphthalene	13.028	162	899030	16.767 ng/ul	100
45) 2-Nitroaniline	13.245	65	258095	17.297 ng/ul	97
47) Dimethylphthalate	13.640	163	1140884	16.714 ng/ul	99
48) 2,6-Dinitrotoluene	13.757	165	233014	17.398 ng/ul	96
50) Acenaphthylene	13.887	152	1522225	17.174 ng/ul	99
51) 3-Nitroaniline	14.087	138	230103	16.348 ng/ul	96
52) Acenaphthene	14.234	153	973256	16.942 ng/ul	99
53) 2,4-Dinitrophenol	14.292	184	164453	18.847 ng/ul	97
55) 4-Nitrophenol	14.410	109	161531	17.286 ng/ul	94
56) Dibenzofuran	14.569	168	1387512	17.022 ng/ul	99
57) 2,4-Dinitrotoluene	14.551	165	344279	17.373 ng/ul	95
58) 2,3,4,6-Tetrachlorophenol	14.804	232	277214	16.927 ng/ul	96
59) Diethylphthalate	15.016	149	1165675	16.840 ng/ul	99
61) Fluorene	15.222	166	1106254	16.990 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.228	204	545278	16.834 ng/ul	98
63) 4-Nitroaniline	15.257	138	240465	17.171 ng/ul	97
66) 4,6-Dinitro-2-methylph	15.316	198	205985	16.747 ng/ul	98
67) N-Nitrosodiphenylamine	15.445	169	1001148	17.464 ng/ul	99
68) 4-Bromophenyl-phenylether	16.122	248	343179	17.130 ng/ul	97
69) Hexachlorobenzene	16.228	284	396446	17.021 ng/ul	96
70) Atrazine	16.410	200	360733	17.142 ng/ul	97
71) Pentachlorophenol	16.575	266	290666	20.293 ng/ul	94
72) Phenanthrene	16.963	178	1812330	17.134 ng/ul	99
74) Anthracene	17.057	178	1864178	17.508 ng/ul	99
75) 1,2,3,4-Tetrachloroben	12.951	216	465632	17.160 ng/uL	95
76) Pentachlorobenzene	14.492	250	478829	16.979 ng/uL	99
77) Carbazole	17.333	167	1700475	17.822 ng/ul	98
78) Di-n-butylphthalate	17.916	149	2000960	17.409 ng/ul	99
80) Fluoranthene	18.992	202	2161191	17.866 ng/ul	95
82) Pyrene	19.357	202	2233341	18.026 ng/ul	99
83) Butylbenzylphthalate	20.292	149	945300	18.343 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.063	252	794271	18.178 ng/ul	98
85) Benzo(a)anthracene	21.122	228	2221197	17.843 ng/ul	97
86) Bis(2-ethylhexyl)phtha	21.080	149	1472919	18.754 ng/ul	100
87) Chrysene	21.174	228	2167007	17.923 ng/ul	98
89) Di-n-octyl phthalate	21.951	149	2529383	21.042 ng/ul	100
90) Benzo(b)fluoranthene	22.674	252	2429500	19.986 ng/ul	99
91) Benzo(k)fluoranthene	22.721	252	2337766	20.413 ng/ul	98
93) Benzo(a)pyrene	23.239	252	2408016	20.388 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	25.545	276	2778526	19.959 ng/ul	97
95) Dibenzo(a,h)anthracene 96) Benzo(g,h,i)perylene	25.557	278	2344979	19.872 ng/ul	98
50) Belizo(g,ll,1)perytene	26.221	276	2291181	19.450 ng/ul	99

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