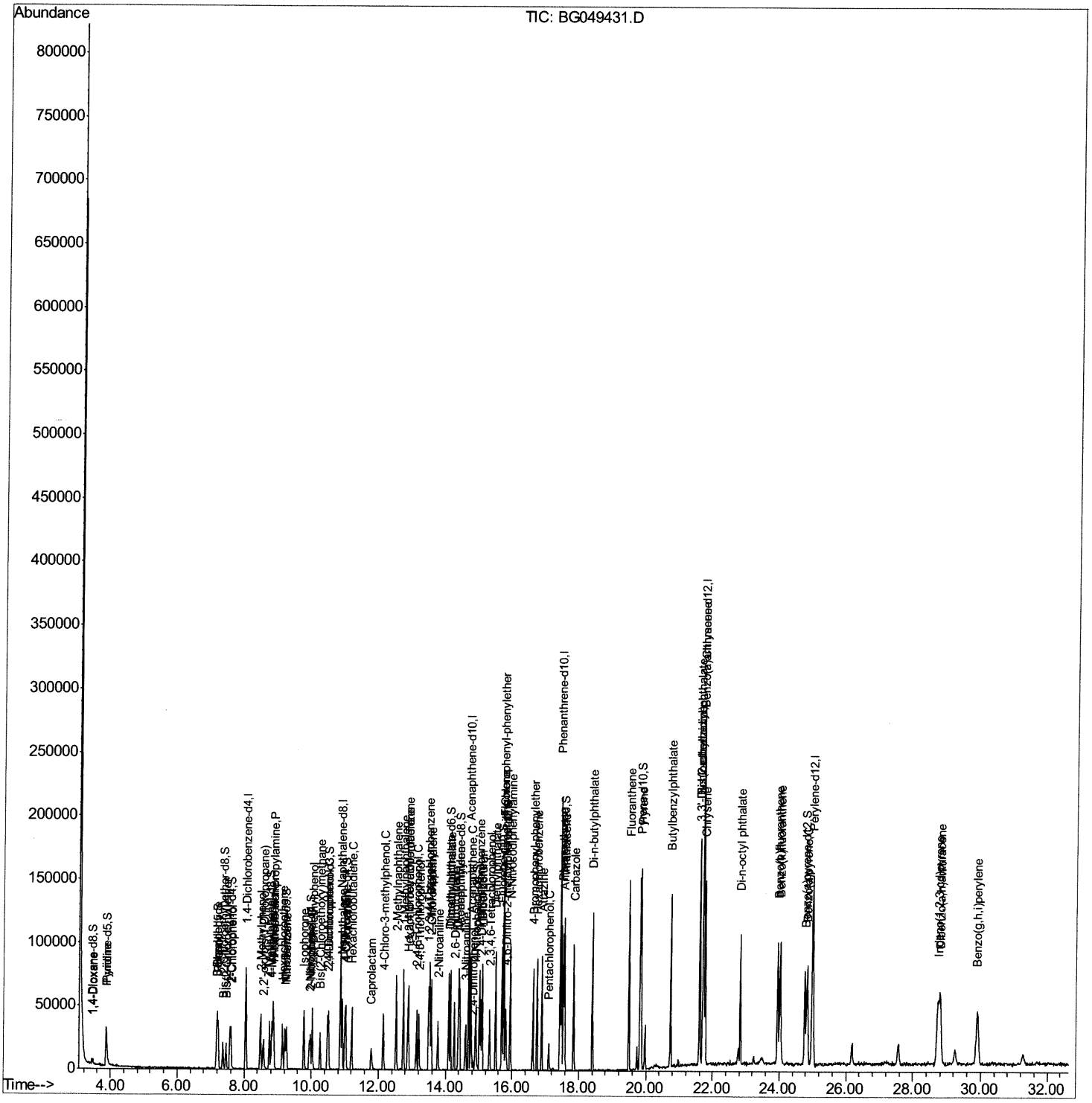
Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 24 06:07:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
7/26/2021 1:28:58 PM



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
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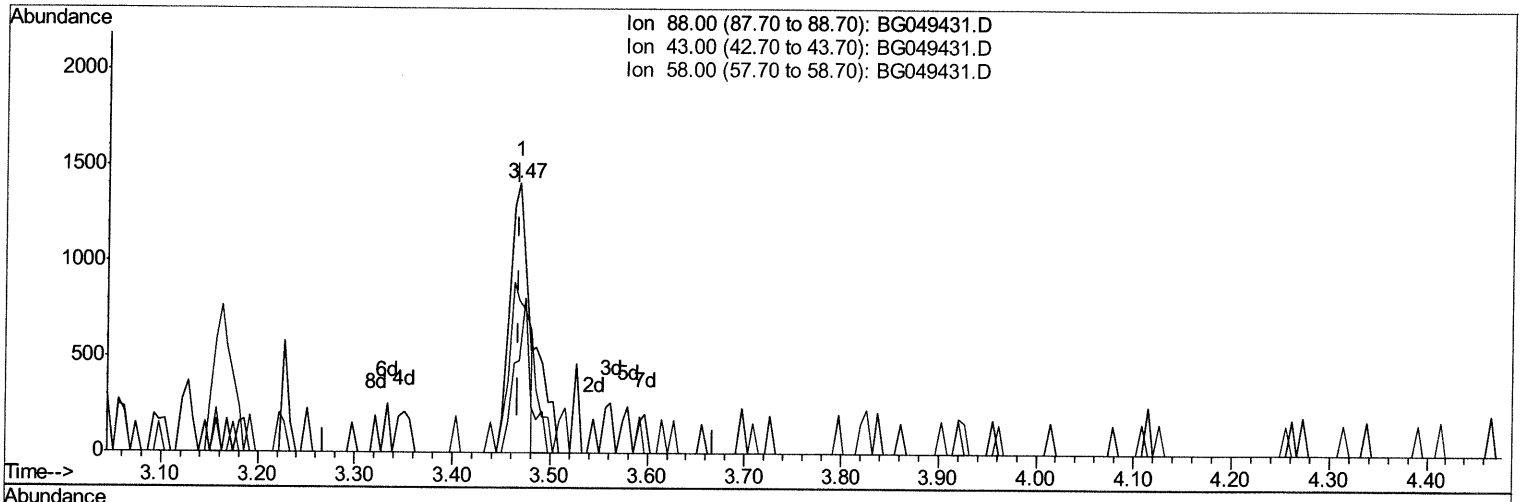
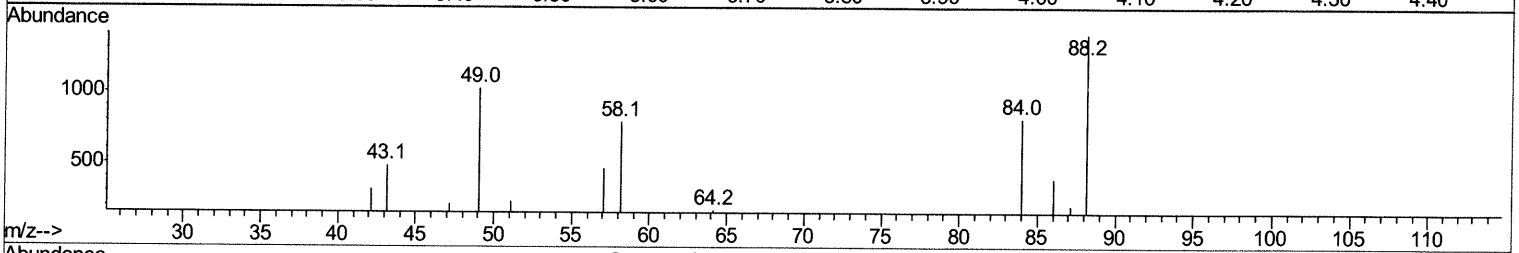
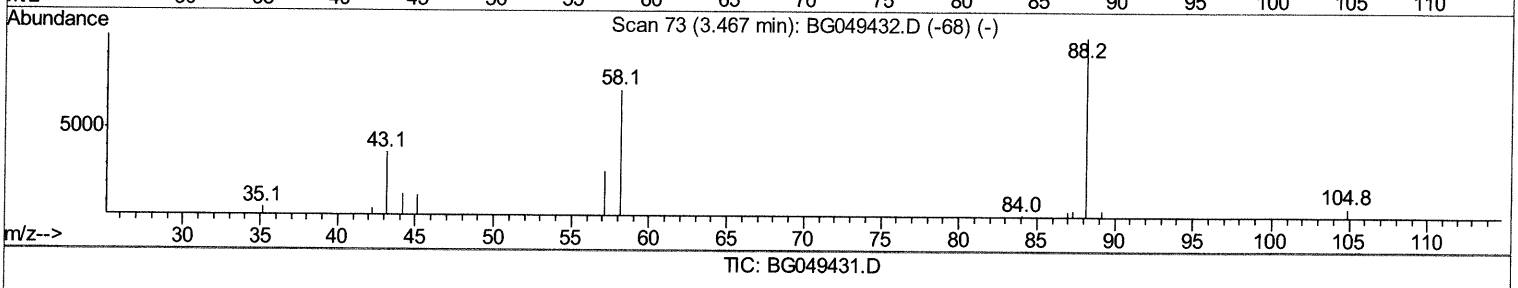






TIC: BG049431.D

(2) 1,4-Dioxane

3.468min (+0.001) 3.36ng/uL

response 1786

Ion	Exp%	Act%
88.00	100	100
43.00	36.90	34.21
58.00	71.00	56.21#
0.00	0.00	0.00

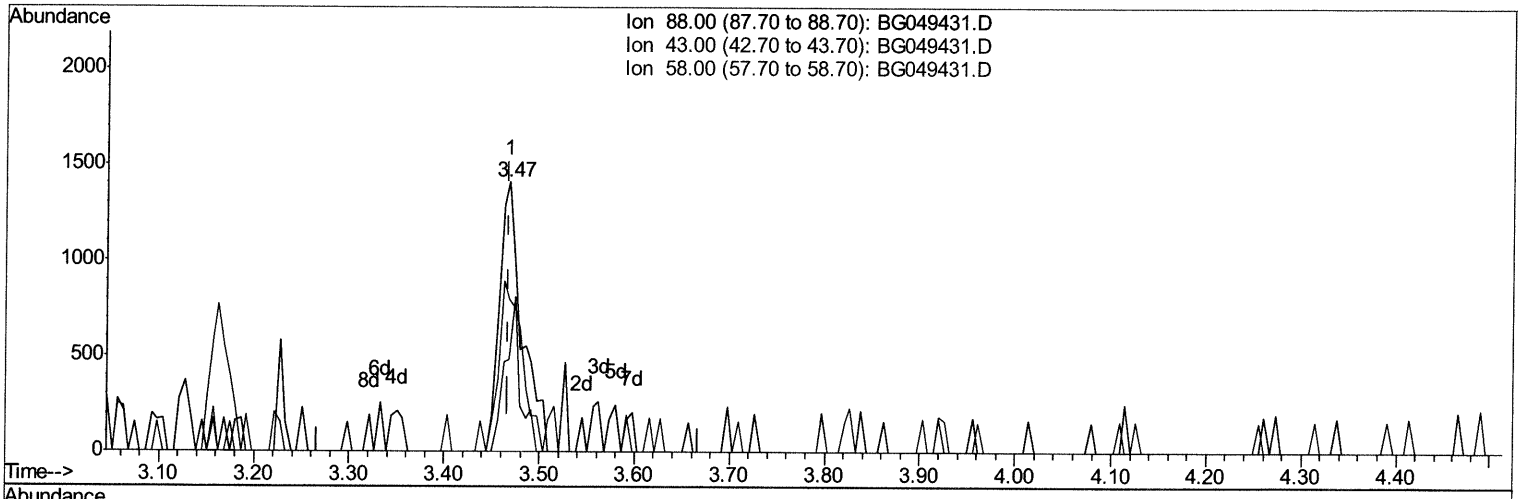
Quantitation Report (Qedit)

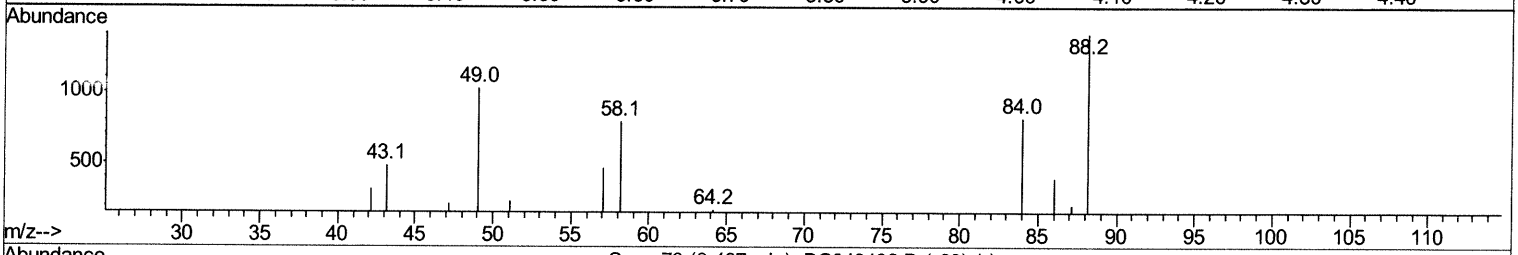
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

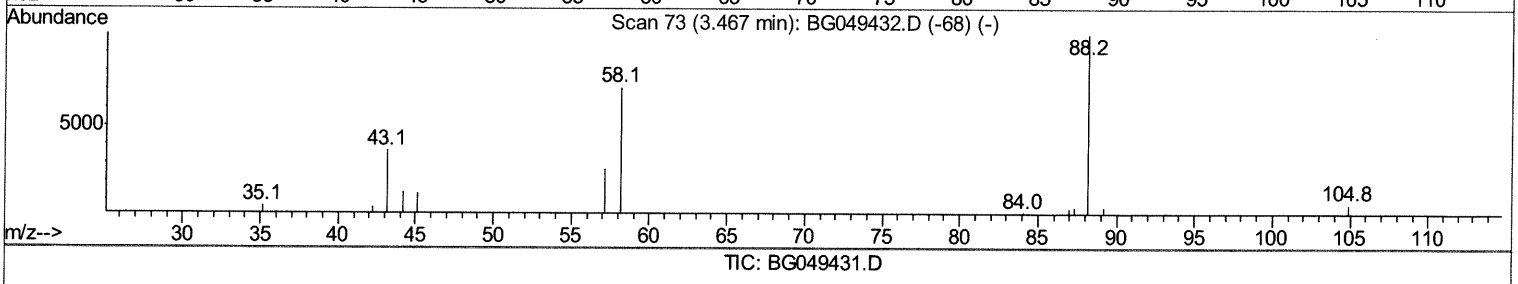
Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
7/26/2021 1:28:58 PM







TIC: BG049431.D

(2) 1,4-Dioxane

3.468min (+0.001) 4.39ng/uL m 07/27/21 JU

response 2334

Ion	Exp%	Act%
88.00	100	100
43.00	36.90	34.21
58.00	71.00	56.21#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
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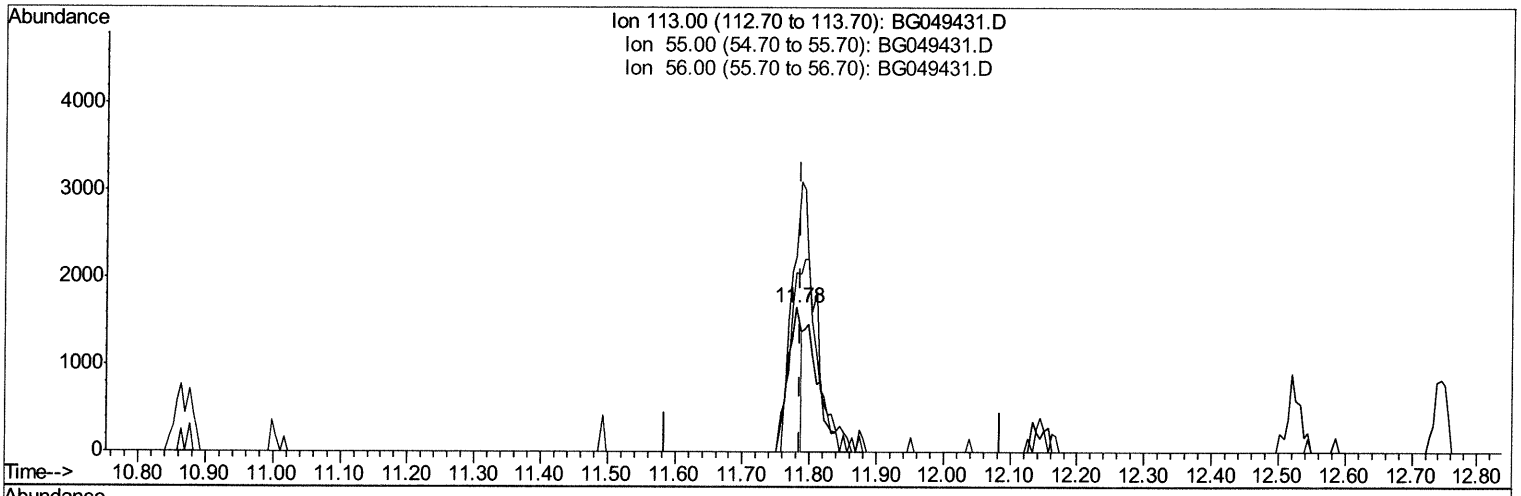

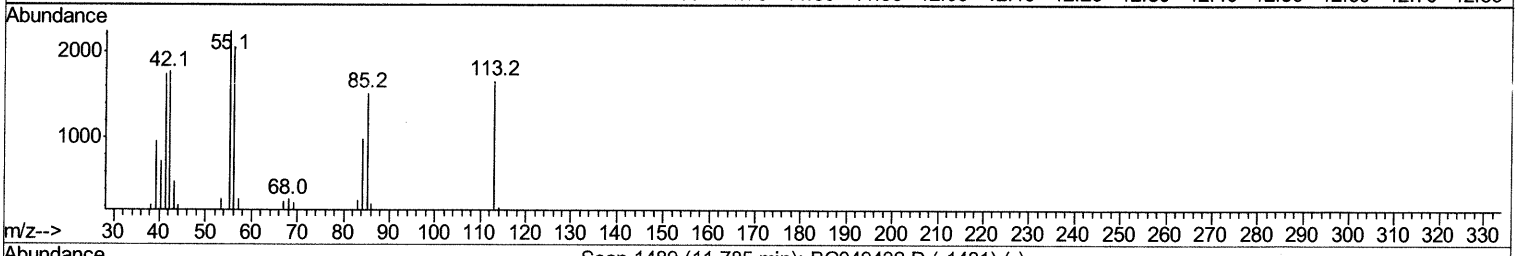

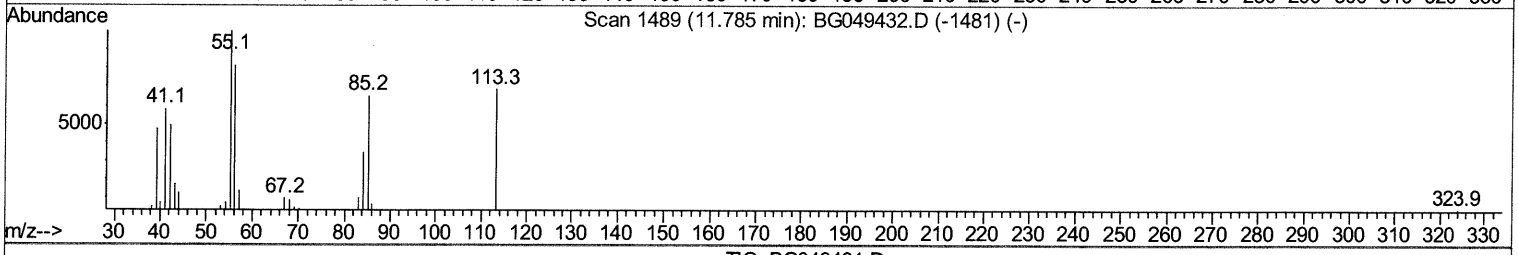

TIC: BG049431.D

(34) Caprolactam

11.780min (-0.005) 3.87ng/ul

response 2275

Ion	Exp%	Act%
113.00	100	100
55.00	145.70	135.34
56.00	118.80	123.96
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
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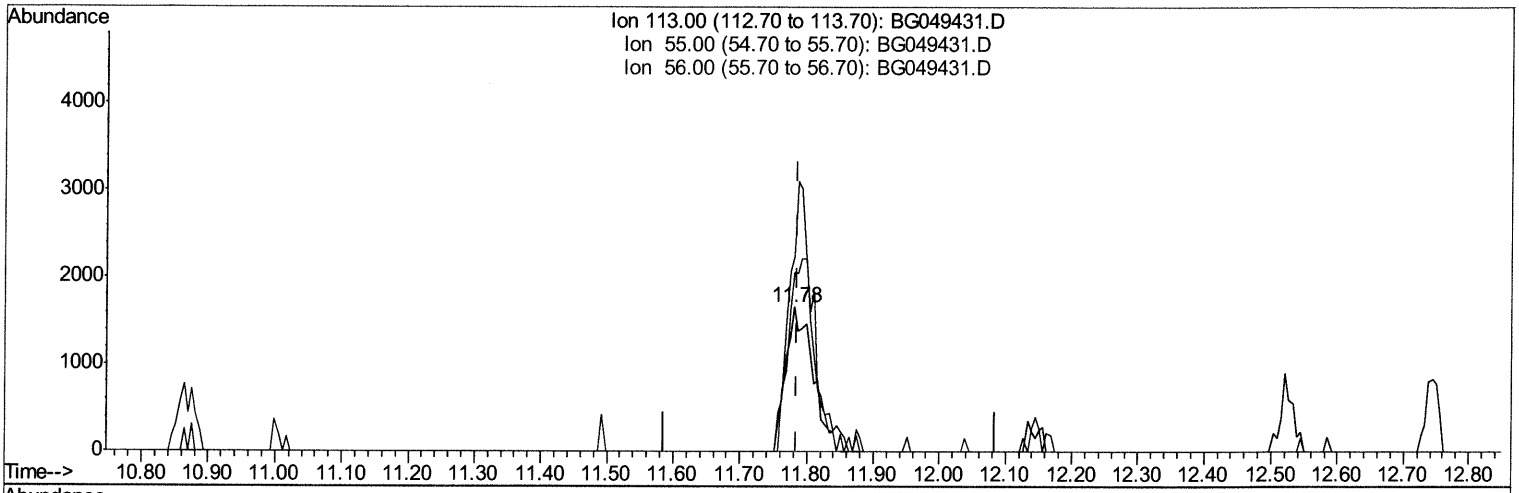

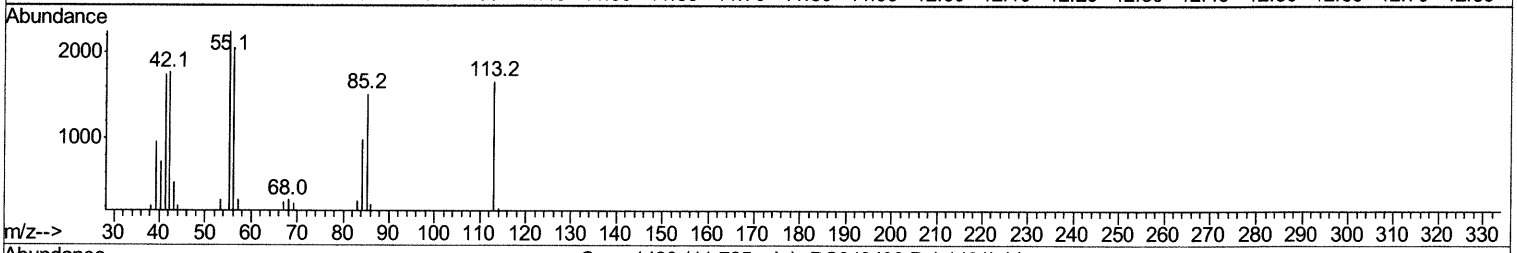

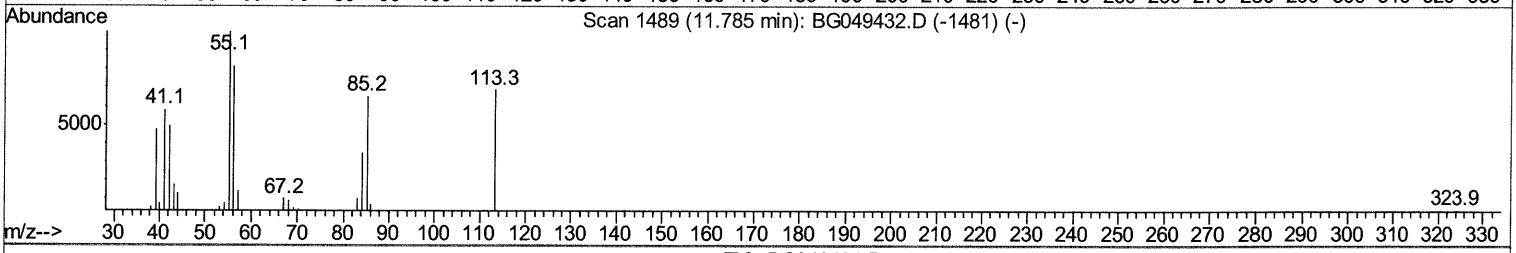

TIC: BG049431.D

(34) Caprolactam

11.780min (-0.005) 7.93ng/ul m 07/27/21 JU

response 4659

Ion	Exp%	Act%
113.00	100	100
55.00	145.70	135.34
56.00	118.80	123.96
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
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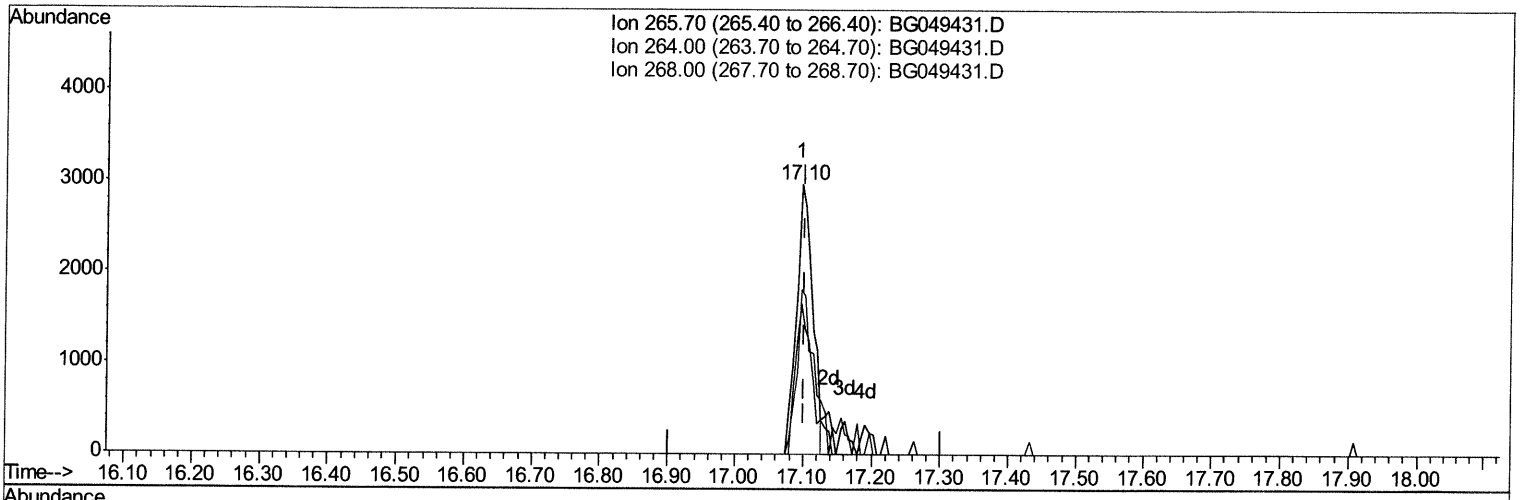

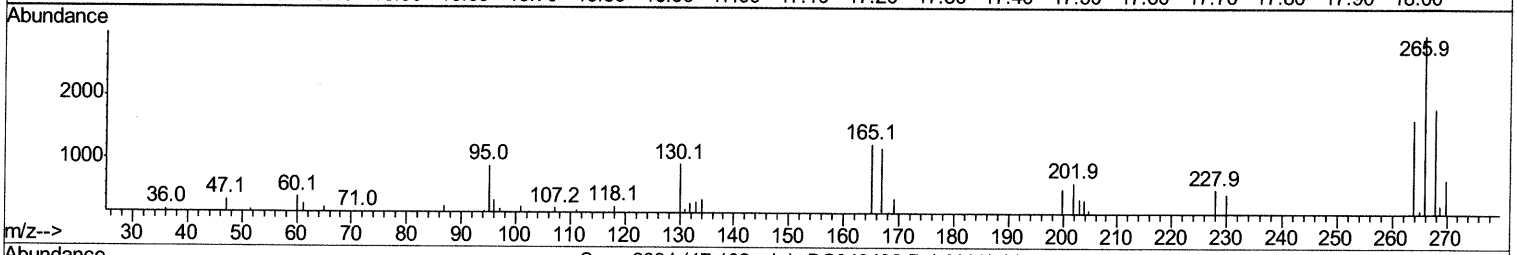

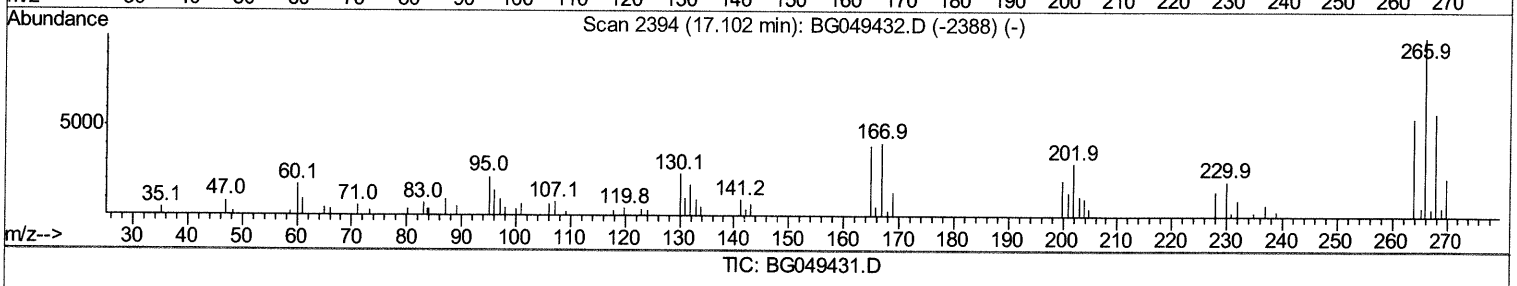

TIC: BG049431.D

(71) Pentachlorophenol (C)

17.097min (-0.005) 4.27ng/ul

response 4971

Ion	Exp%	Act%
265.70	100	100
264.00	54.70	55.55
268.00	57.80	61.06
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
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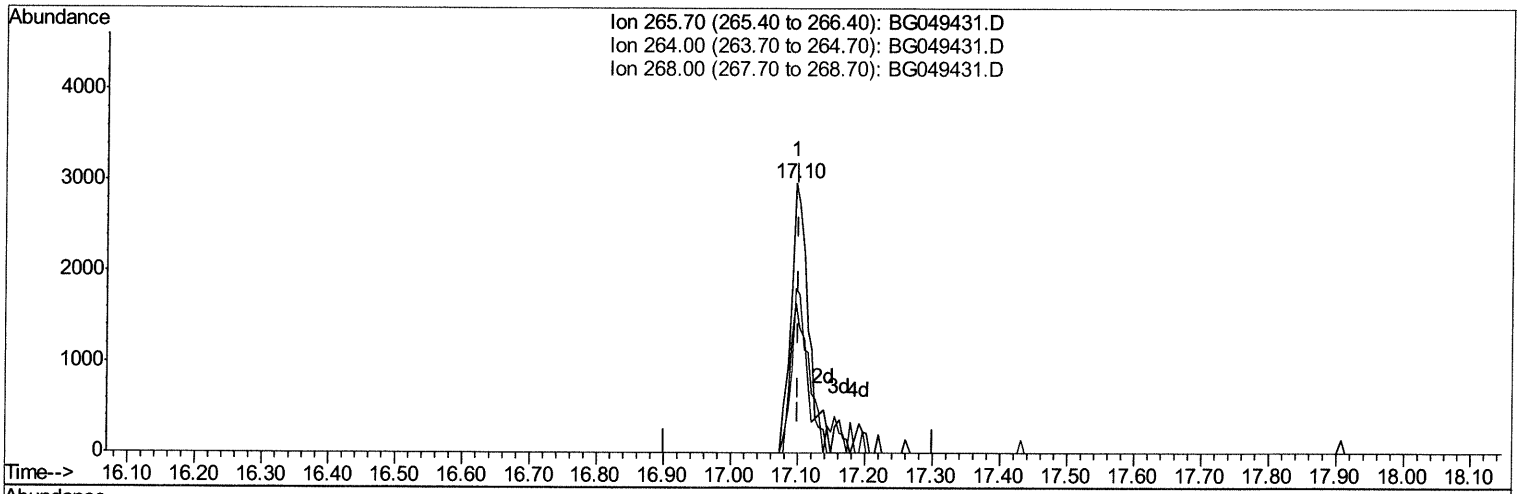

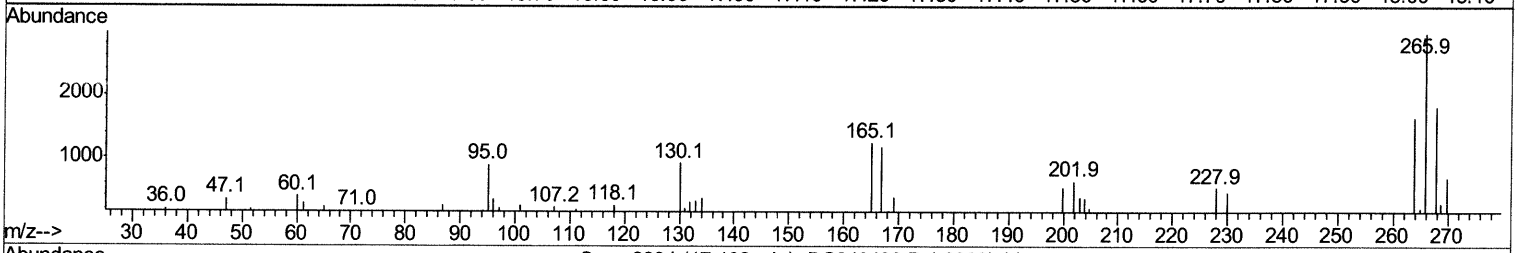

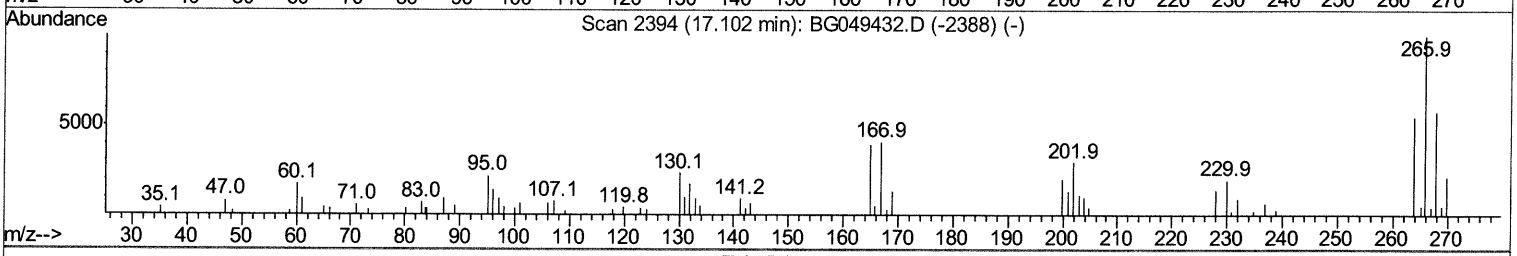

TIC: BG049431.D

(71) Pentachlorophenol (C)

17.097min (-0.005) 4.62ng/ul m 07/27/21 JU

response 5376

Ion	Exp%	Act%
265.70	100	100
264.00	54.70	55.55
268.00	57.80	61.06
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
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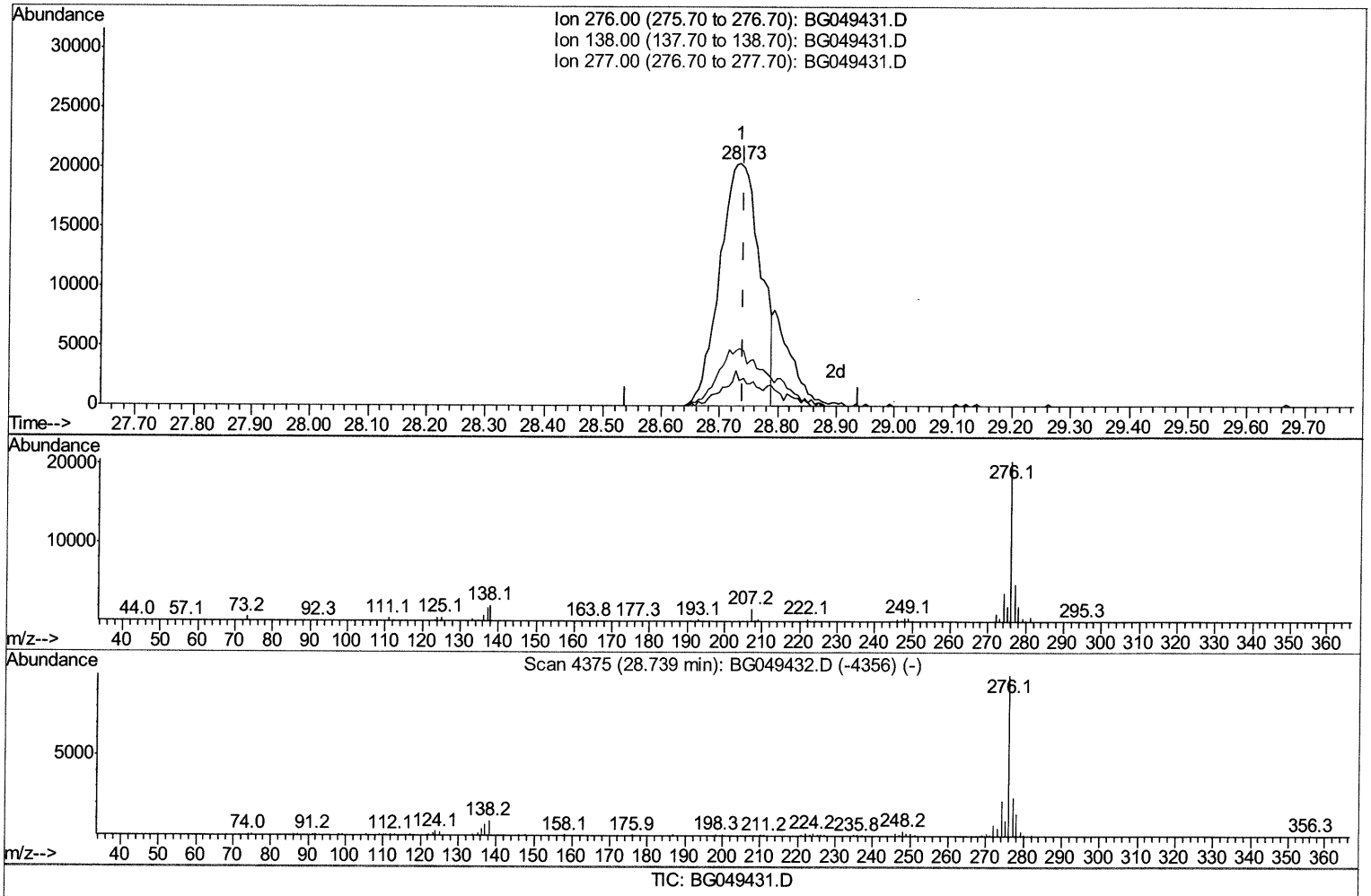

TIC: BG049431.D

(94) Indeno(1,2,3-cd)pyrene

28.734min (-0.005) 8.07ng/ul

response 97246

Ion	Exp%	Act%
276.00	100	100
138.00	9.10	10.29
277.00	24.00	23.80
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
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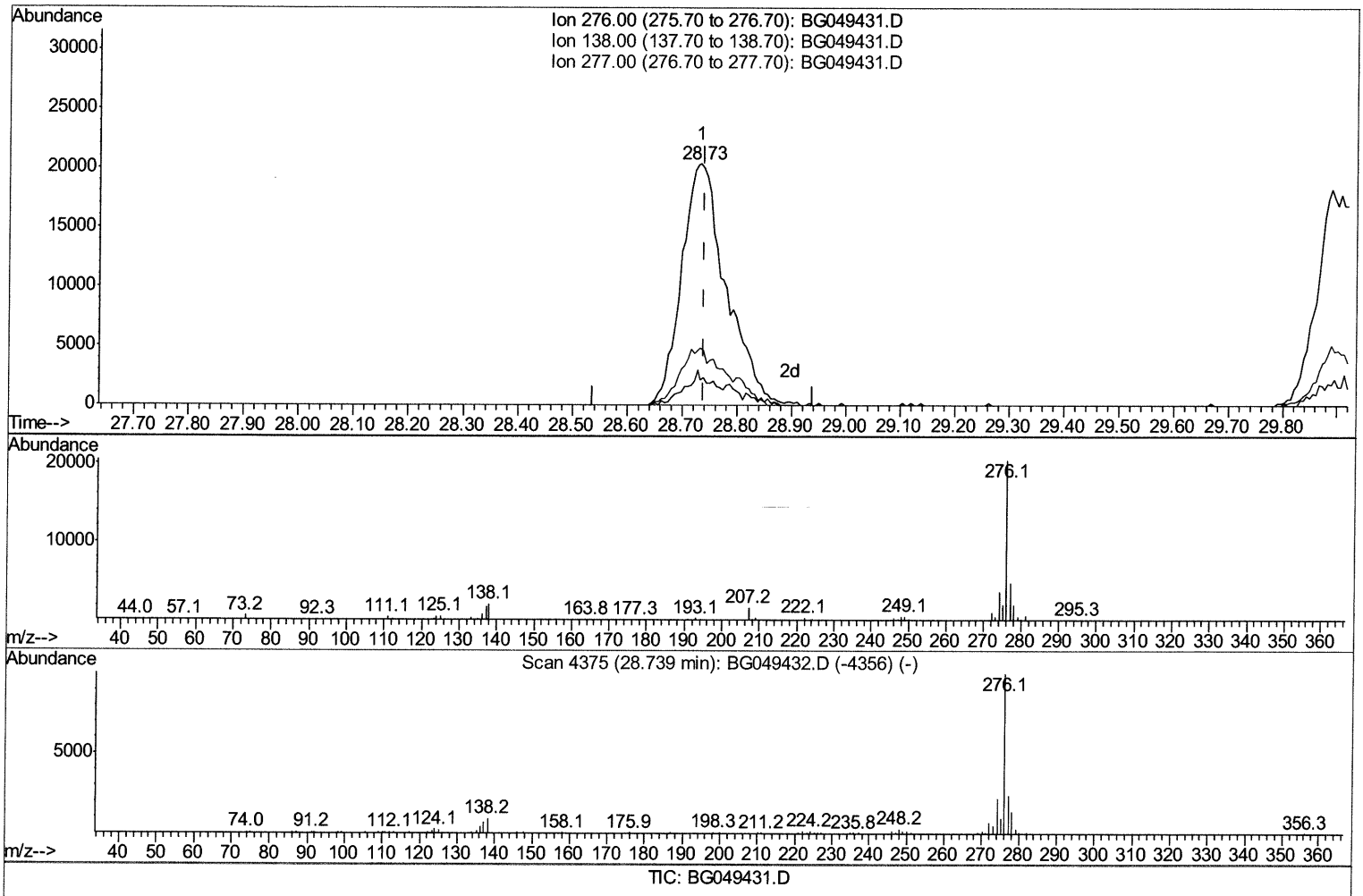


TIC: BG049431.D

(94) Indeno(1,2,3-cd)pyrene

28.734min (-0.005) 9.56ng/ul m 07/27/21 JU

response 115216

Ion	Exp%	Act%
276.00	100	100
138.00	9.10	10.29
277.00	24.00	23.80
0.00	0.00	0.00

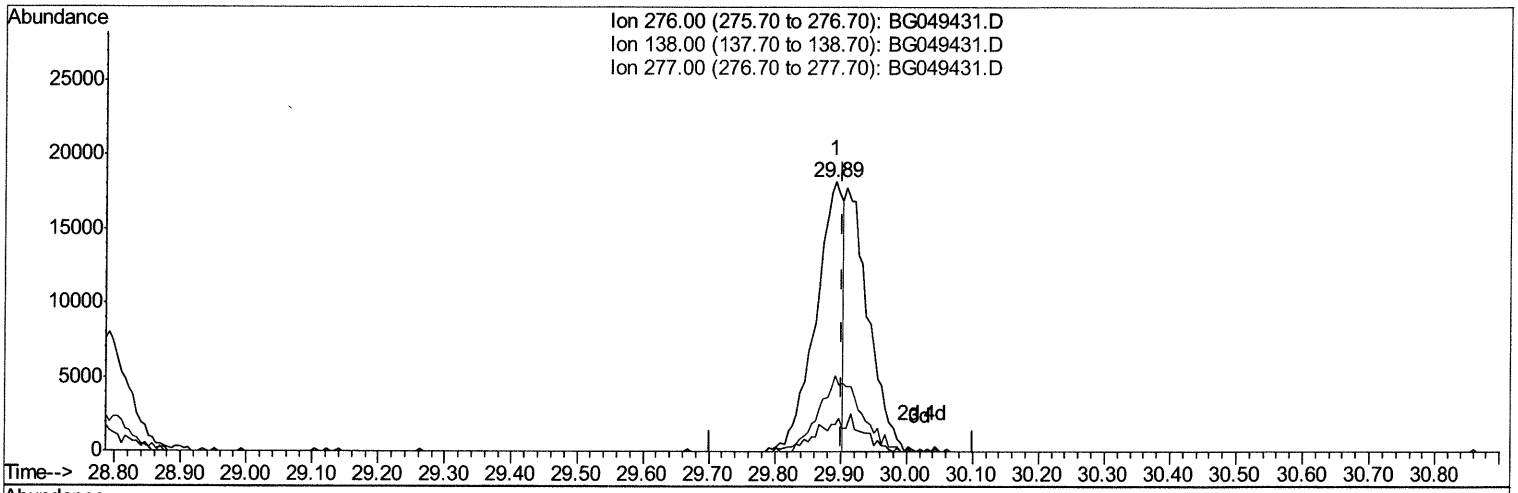
Quantitation Report (Qedit)

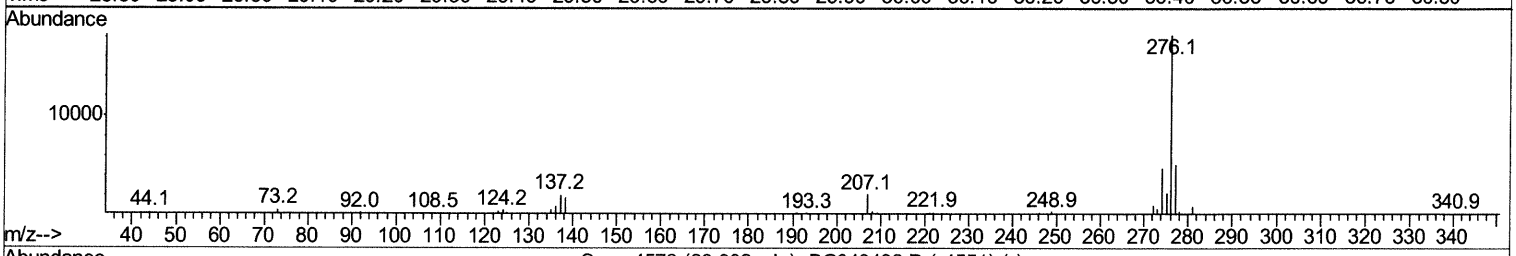
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

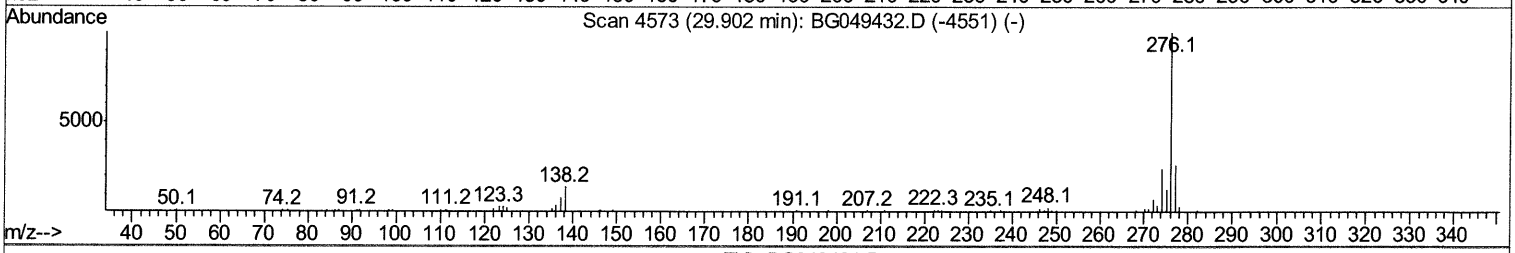
Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
7/26/2021 1:28:58 PM







TIC: BG049431.D

(96) Benzo(g,h,i)perylene

29.891min (-0.011) 5.43ng/ul

response 52986

Ion	Exp%	Act%
276.00	100	100
138.00	13.50	9.92#
277.00	25.80	28.16
0.00	0.00	0.00

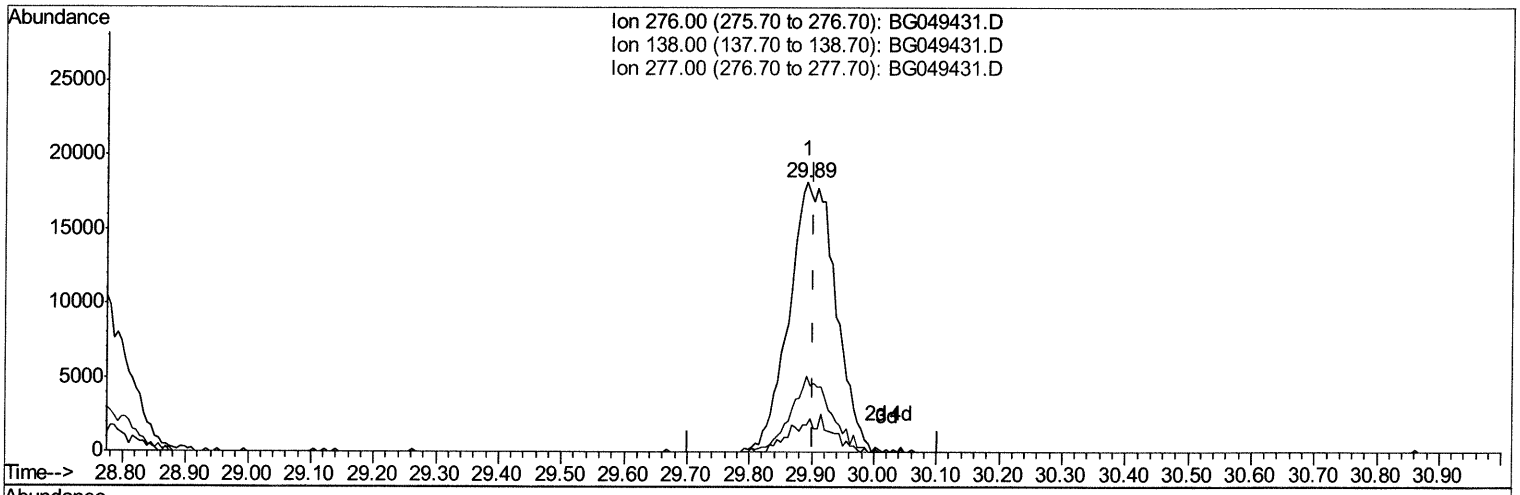
Quantitation Report (Qedit)

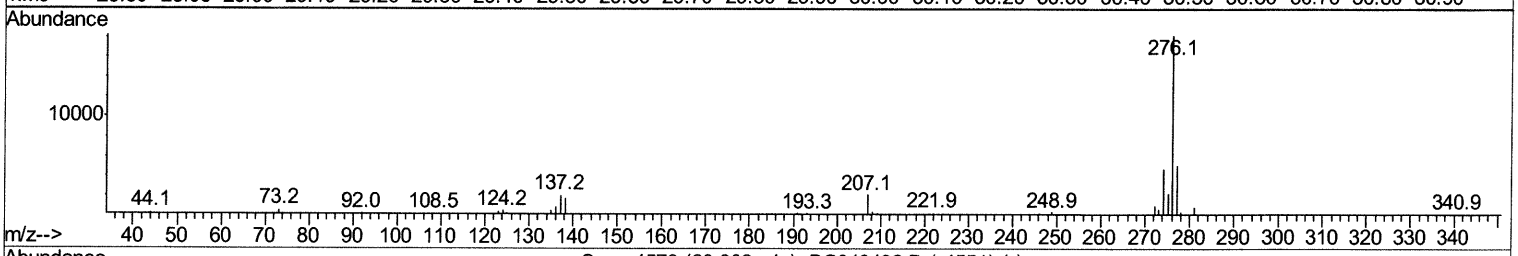
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

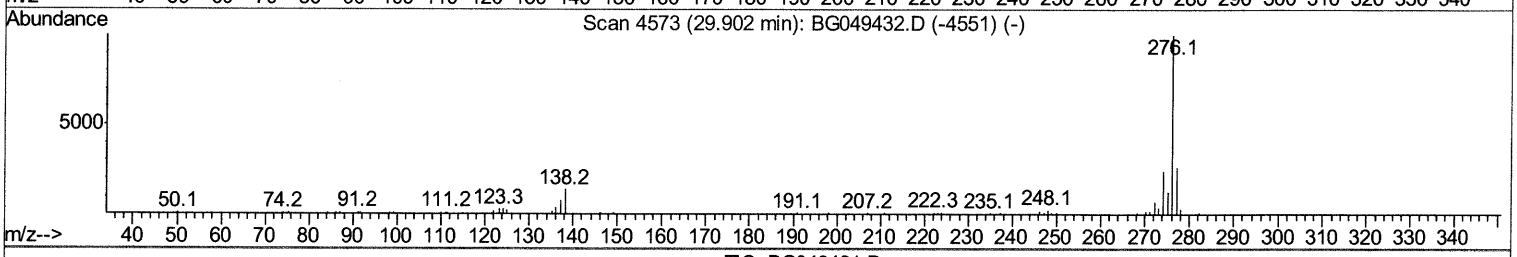
Quant Time: Jul 23 23:55:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
7/26/2021 1:28:58 PM







TIC: BG049431.D

(96) Benzo(g,h,i)perylene

29.891min (-0.011) 9.73ng/ul m 07/27/21 JU

response 95080

Ion	Exp%	Act%
276.00	100	100
138.00	13.50	9.92#
277.00	25.80	28.16
0.00	0.00	0.00

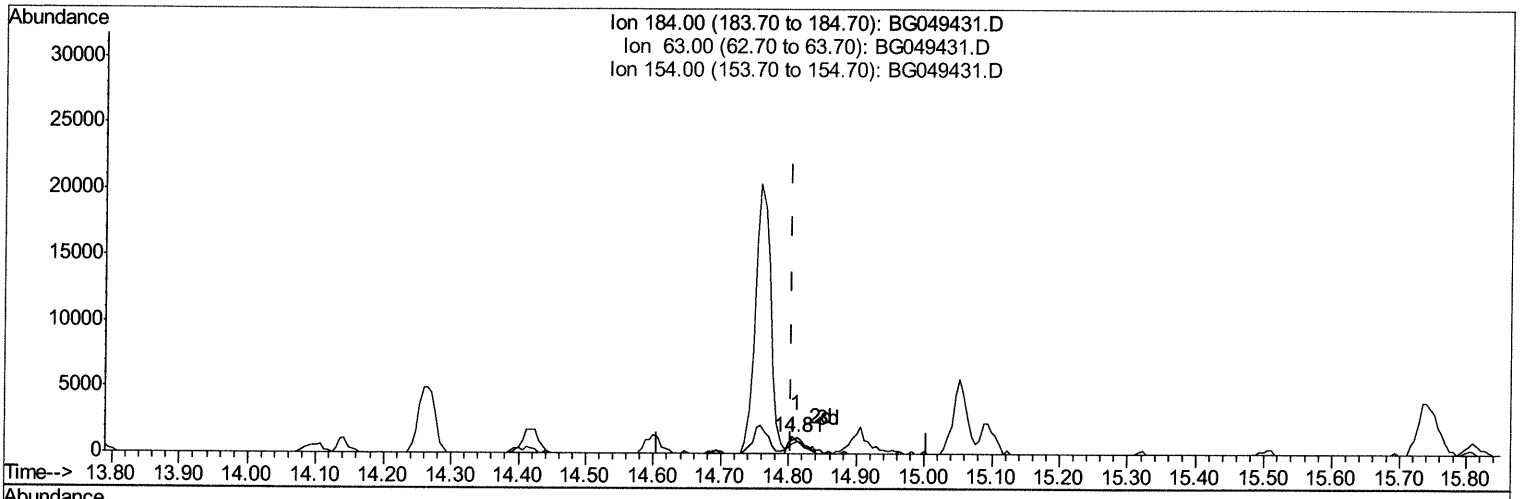
Quantitation Report (Qedit)

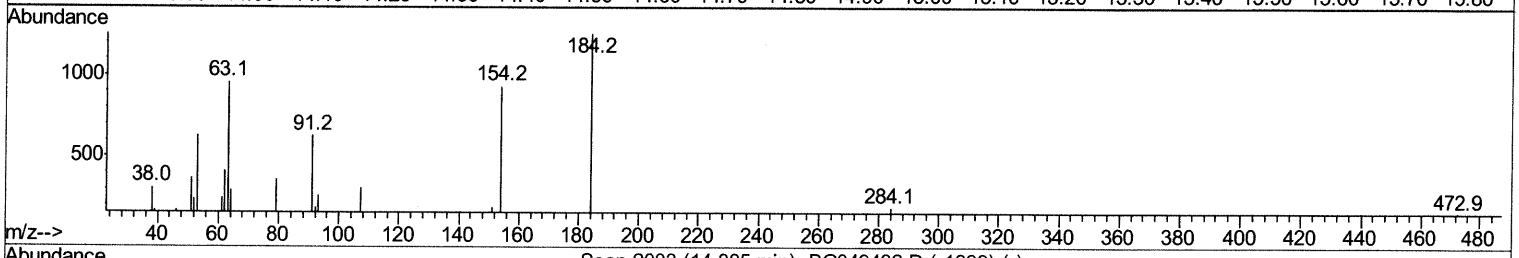
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

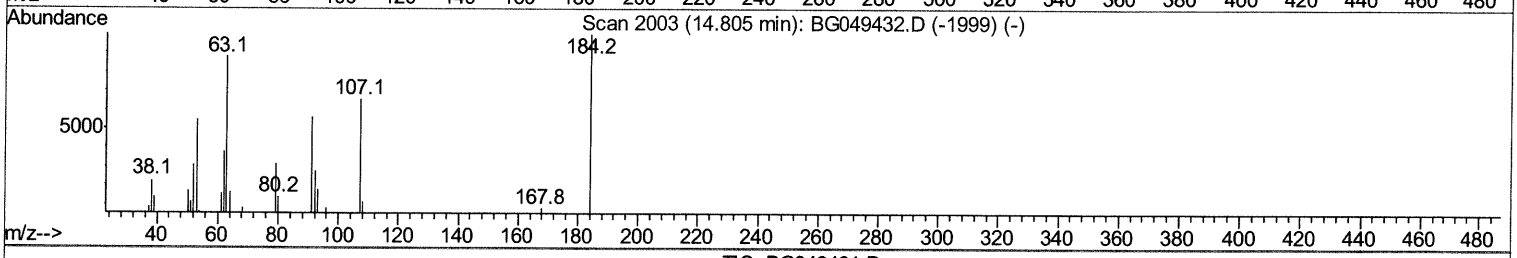
Quant Time: Jul 24 06:07:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
mohammad
7/26/2021 1:28:58 PM







TIC: BG049431.D

(53) 2,4-Dinitrophenol

14.812min (+0.007) 3.94ng/ul m 07/27/21 JU

response 2184

Ion	Exp%	Act%
184.00	100	100
63.00	106.90	76.00#
154.00	94.70	74.24#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 24 00:29:47 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
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Manual Integrations APPROVED
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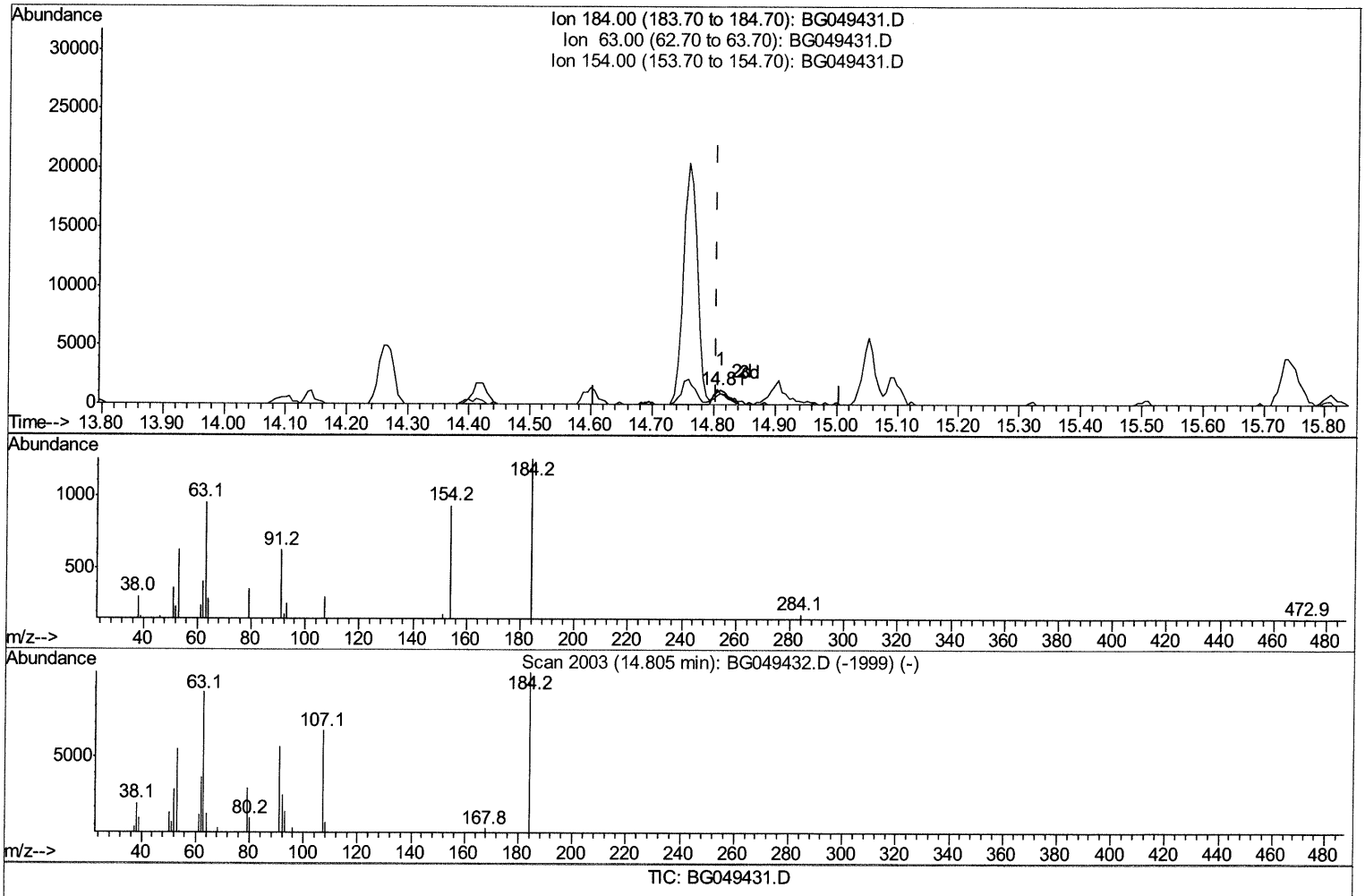

TIC: BG049431.D

(53) 2,4-Dinitrophenol

14.812min (+0.007) 3.77ng/ul

response 2087

Ion	Exp%	Act%
184.00	100	100
63.00	106.90	76.00#
154.00	94.70	74.24#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG072421\
Data File : BG049431.D
Acq On : 23 Jul 2021 18:30
Operator : CG/JU
Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
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7/26/2021 1:28:58 PM

Quant Time: Jul 24 06:07:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	22930	20.00	ng/ul	0.00
20) Naphthalene-d8	10.87	136	88631	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.69	164	58543	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.45	188	137162	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	152927	20.00	ng/ul	0.00
88) Perylene-d12	25.00	264	161215	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.43	96	1927	3.64	ng/uL	0.00
4) Pyridine-d5	3.86	84	13442	8.30	ng/ul	0.00
7) Phenol-d5	7.20	99	18945	9.51	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.37	67	10951	8.90	ng/ul	0.00
11) 2-Chlorophenol-d4	7.57	132	14377	10.39	ng/ul	0.00
15) 4-Methylphenol-d8	8.75	113	14445	8.84	ng/ul	0.00
21) Nitrobenzene-d5	9.21	128	7353	11.98	ng/ul	0.00
24) 2-Nitrophenol-d4	9.94	143	7916	11.78	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.49	165	14240	8.62	ng/ul	0.00
31) 4-Chloroaniline-d4	11.00	131	20232	9.73	ng/ul	0.00
46) Dimethylphthalate-d6	14.09	166	47684	9.90	ng/ul	0.00
49) Acenaphthylene-d8	14.39	160	55058	10.41	ng/ul	0.00
54) 4-Nitrophenol-d4	14.89	143	6380	8.90	ng/ul	0.00
60) Fluorene-d10	15.69	176	41168	9.68	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.80	200	7374	10.52	ng/ul	0.00
73) Anthracene-d10	17.54	188	63961	10.21	ng/ul	0.00
81) Pyrene-d10	19.83	212	82616	10.56	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.77	264	84510	9.67	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.47	88	2334m >	4.395	ng/uL >	07/27/21 JU
5) Pyridine	3.88	79	14568	9.002	ng/ul	96
6) Benzaldehyde	7.18	77	10667	8.480	ng/ul	86
8) Phenol	7.23	94	18267	8.674	ng/ul	93
10) Bis(2-Chloroethyl)ether	7.46	93	12651	8.237	ng/ul#	86
12) 2-Chlorophenol	7.60	128	14078	9.581	ng/ul	88
13) 2-Methylphenol	8.49	108	13671	8.677	ng/ul	88
14) 2,2'-oxybis(1-Chloropropan	8.57	45	16076	8.860	ng/ul#	85
16) Acetophenone	8.87	105	22613	8.441	ng/ul#	85
17) N-Nitroso-di-n-propylamine	8.85	70	12539	7.894	ng/ul#	92
18) 4-Methylphenol	8.81	108	13627	8.239	ng/ul	96
19) Hexachloroethane	9.14	117	6392	9.335	ng/ul#	83
22) Nitrobenzene	9.26	77	21108	10.433	ng/ul#	98
23) Isophorone	9.78	82	34269	8.195	ng/ul	97
25) 2-Nitrophenol	9.98	139	7569	9.892	ng/ul#	90
26) 2,4-Dimethylphenol	10.03	107	17189	8.538	ng/ul#	93
27) Bis(2-Chloroethoxy)methane	10.26	93	17254	8.196	ng/ul	92
29) 2,4-Dichlorophenol	10.51	162	14647	9.497	ng/ul#	90
30) Naphthalene	10.92	128	47016	10.038	ng/ul	95
32) 4-Chloroaniline	11.03	127	18638	9.115	ng/ul#	83
33) Hexachlorobutadiene	11.20	225	11279	7.688	ng/ul	95
34) Caprolactam	11.78	113	4659m >	7.929	ng/ul >	07/27/21 JU

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Sample : SSTD01002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jul 24 06:07:03 2021
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Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 23 23:52:30 2021
Response via : Initial Calibration

Instrument : BNA_G
ClientSampleId : SSTD010409

Manual Integrations APPROVED
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Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
35) 4-Chloro-3-methylphenol	12.14	107	15306	8.462 ng/ul	88
36) 2-Methylnaphthalene	12.53	142	33642	9.312 ng/ul	97
37) 1-Methylnaphthalene	12.74	142	34862	9.742 ng/ul#	97
39) 1,2,4,5-Tetrachlorobenzene	12.89	216	18583	8.339 ng/ul	92
40) Hexachlorocyclopentadiene	12.87	237	8346	5.157 ng/ul#	91
41) 2,4,6-Trichlorophenol	13.13	196	10972	8.001 ng/ul#	85
42) 2,4,5-Trichlorophenol	13.21	196	11336	7.701 ng/ul	94
43) 1,1'-Biphenyl	13.53	154	44212	10.206 ng/ul#	92
44) 2-Chloronaphthalene	13.57	162	34631	10.206 ng/ul#	92
45) 2-Nitroaniline	13.78	65	11516	10.986 ng/ul	92
47) Dimethylphthalate	14.14	163	46644	9.589 ng/ul#	96
48) 2,6-Dinitrotoluene	14.27	165	8756	10.510 ng/ul	95
50) Acenaphthylene	14.42	152	52512	10.355 ng/ul	96
51) 3-Nitroaniline	14.59	138	8572	10.702 ng/ul	94
52) Acenaphthene	14.76	153	38963	10.360 ng/ul	95
53) 2,4-Dinitrophenol	14.81	184	2184m >	3.940 ng/ul >	07/27/21 JU
55) 4-Nitrophenol	14.91	109	9067	9.545 ng/ul	98
56) Dibenzofuran	15.10	168	52294	9.788 ng/ul	96
57) 2,4-Dinitrotoluene	15.05	165	13009	10.855 ng/ul#	94
58) 2,3,4,6-Tetrachlorophenol	15.32	232	10187	7.321 ng/ul#	96
59) Diethylphthalate	15.50	149	49172	10.114 ng/ul	97
61) Fluorene	15.75	166	43851	9.643 ng/ul	93
62) 4-Chlorophenyl-phenylether	15.73	204	21711	7.870 ng/ul	89
63) 4-Nitroaniline	15.76	138	7673	9.200 ng/ul	95
66) 4,6-Dinitro-2-methylphenol	15.82	198	6570	8.927 ng/ul#	89
67) N-Nitrosodiphenylamine	15.95	169	36443	9.592 ng/ul	95
68) 4-Bromophenyl-phenylether	16.63	248	15095	8.653 ng/ul#	87
69) Hexachlorobenzene	16.75	284	17542	9.568 ng/ul	92
70) Atrazine	16.89	200	16331	9.552 ng/ul	92
71) Pentachlorophenol	17.10	266	5376m >	4.622 ng/ul >	07/27/21 JU
72) Phenanthrene	17.49	178	73425	10.543 ng/ul	100
74) Anthracene	17.58	178	73772	10.297 ng/ul	98
75) 1,2,3,4-Tetrachlorobenzene	13.50	216	19077	8.871 ng/uL	93
76) Pentachlorobenzene	15.02	250	19144	8.499 ng/uL	94
77) Carbazole	17.85	167	66326	10.927 ng/ul	99
78) Di-n-butylphthalate	18.40	149	81922	10.783 ng/ul	99
80) Fluoranthene	19.50	202	97287	10.025 ng/ul	98
82) Pyrene	19.86	202	98775	10.192 ng/ul	100
83) Butylbenzylphthalate	20.74	149	37403	10.792 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.62	252	35133	9.379 ng/ul#	99
85) Benzo(a)anthracene	21.71	228	97147	9.615 ng/ul	92
86) Bis(2-ethylhexyl)phthalate	21.60	149	55791	11.043 ng/ul#	98
87) Chrysene	21.77	228	93294	9.656 ng/ul	98
89) Di-n-octyl phthalate	22.83	149	91667	10.446 ng/ul	100
90) Benzo(b)fluoranthene	23.95	252	103569	9.713 ng/ul	99
91) Benzo(k)fluoranthene	24.02	252	100056	9.549 ng/ul	97
93) Benzo(a)pyrene	24.84	252	91719	9.791 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	28.73	276	115216m >	9.562 ng/ul >	07/27/21 JU
95) Dibenzo(a,h)anthracene	28.79	278	95772	9.776 ng/ul	97
96) Benzo(g,h,i)perylene	29.89	276	95080m >	9.735 ng/ul >	07/27/21 JU

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

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

Instrument :
BNA_G
ClientSampleId :
SSTD010409

Manual Integrations APPROVED
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