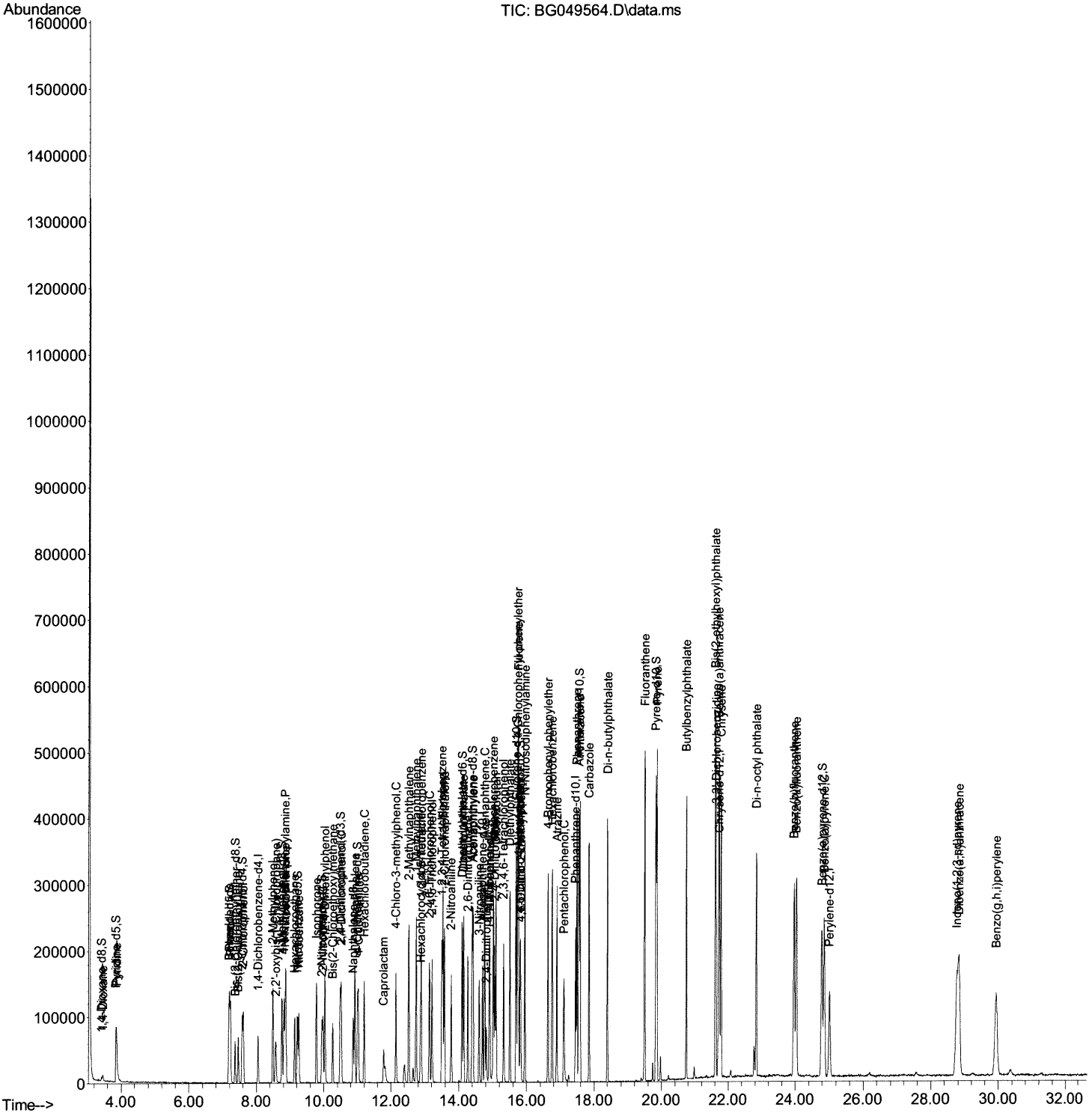
Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM



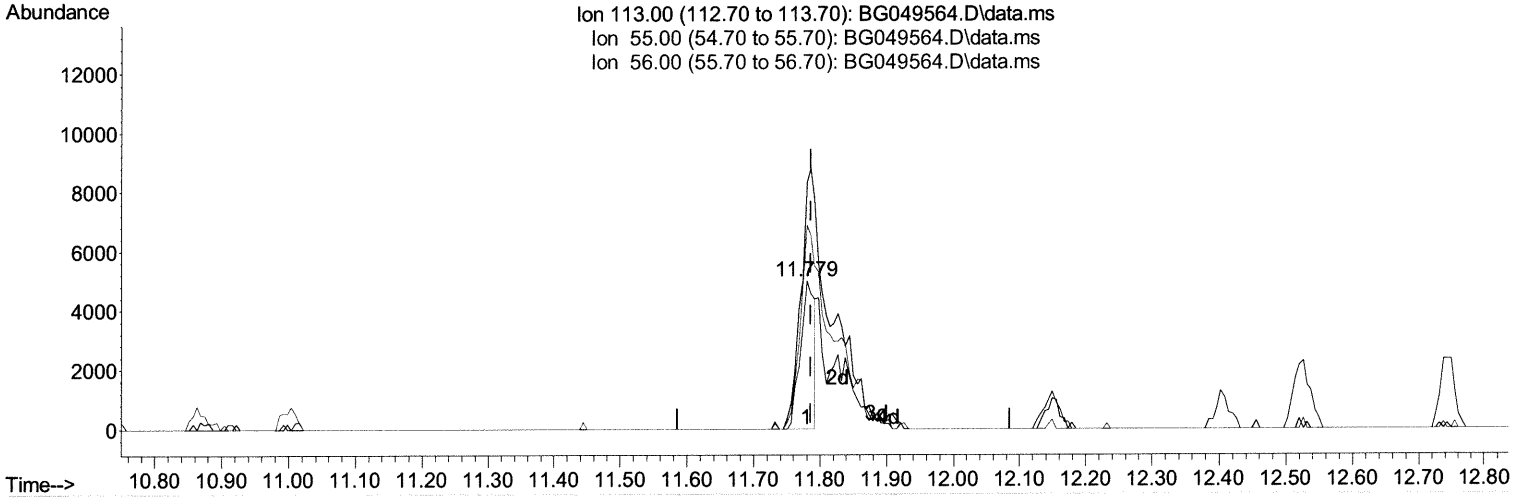
Quantitation Report (Qedit)

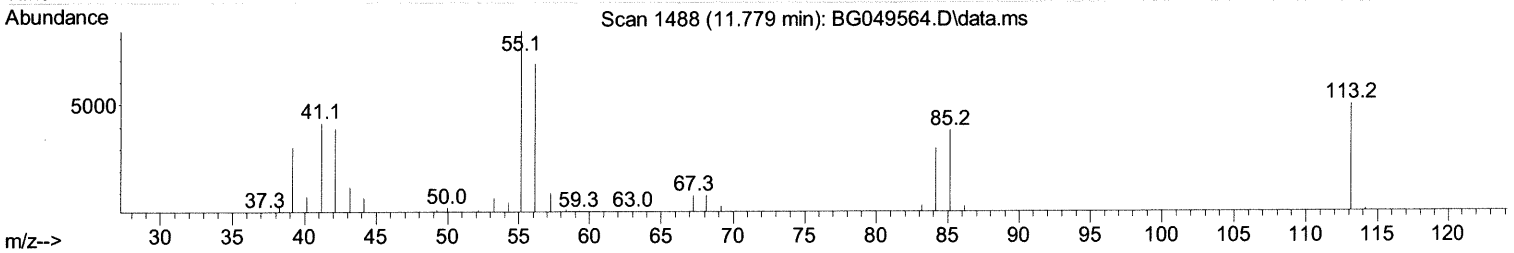
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

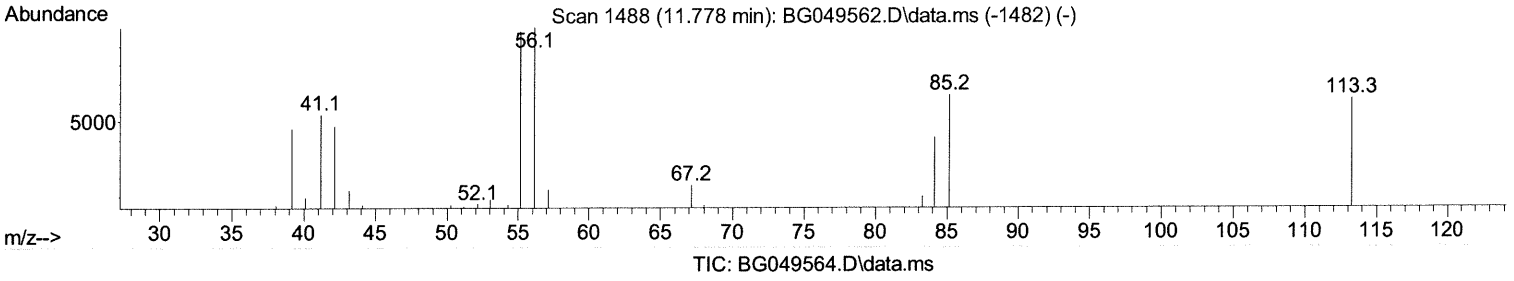
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(34) Caprolactam

11.779min (-0.006) 17.02 ng/ul

response 7411

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	145.70	166.95
56.00	118.80	137.82
0.00	0.00	0.00

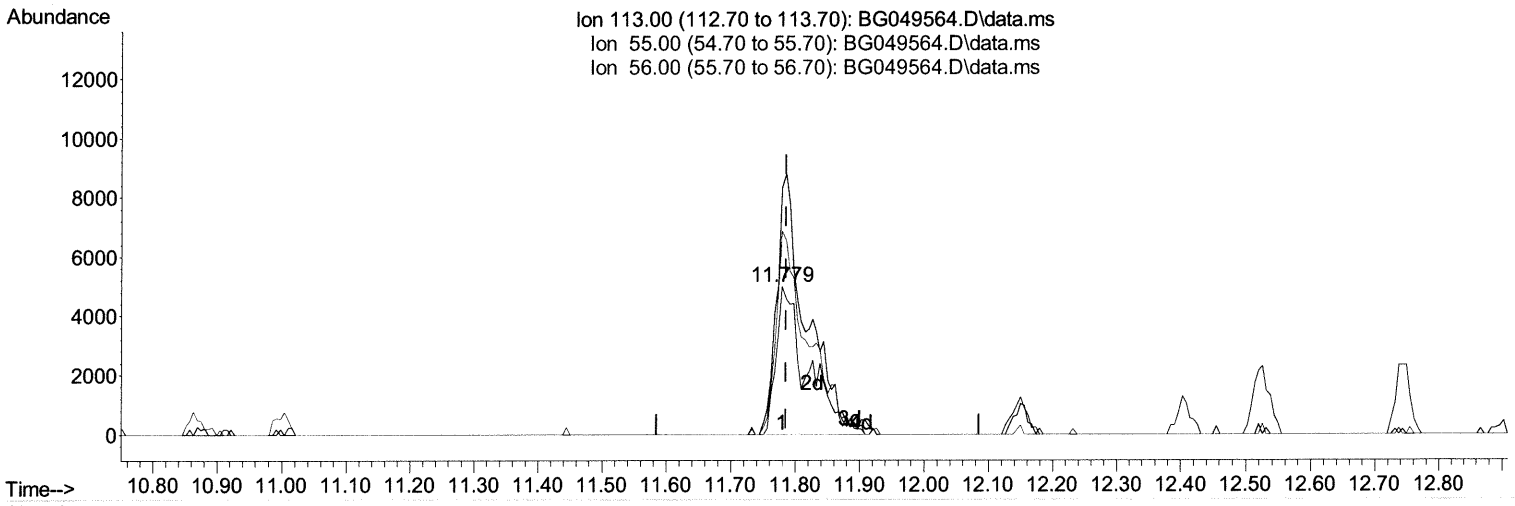
Quantitation Report (Qedit)

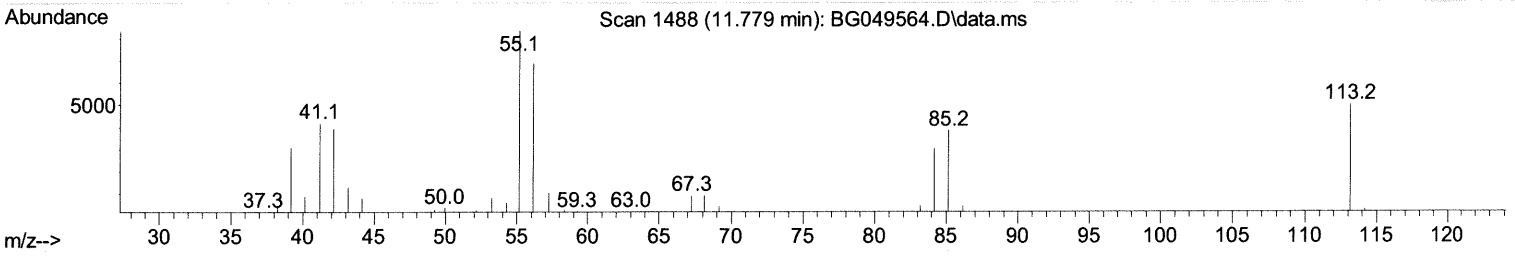
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

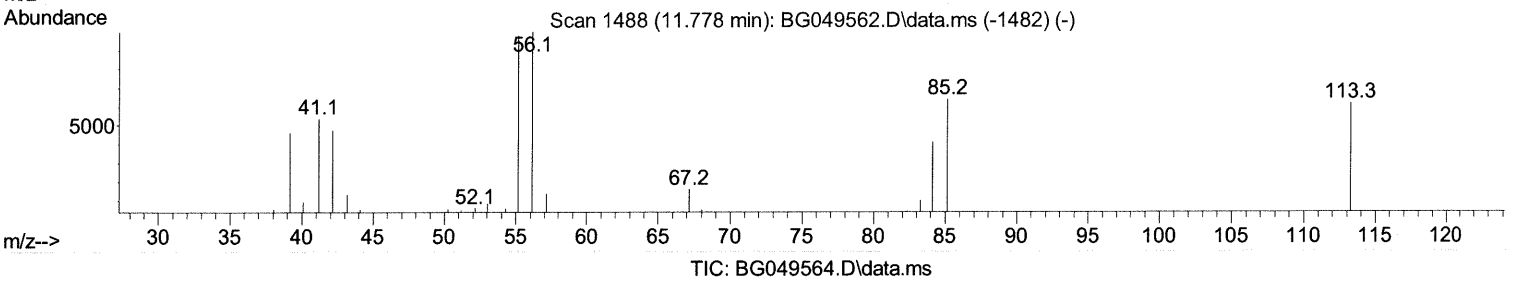
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(34) Caprolactam

11.779min (-0.006) 38.31 ng/ul m 08/02/21 JU

response 16677

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	145.70	166.95
56.00	118.80	137.82
0.00	0.00	0.00

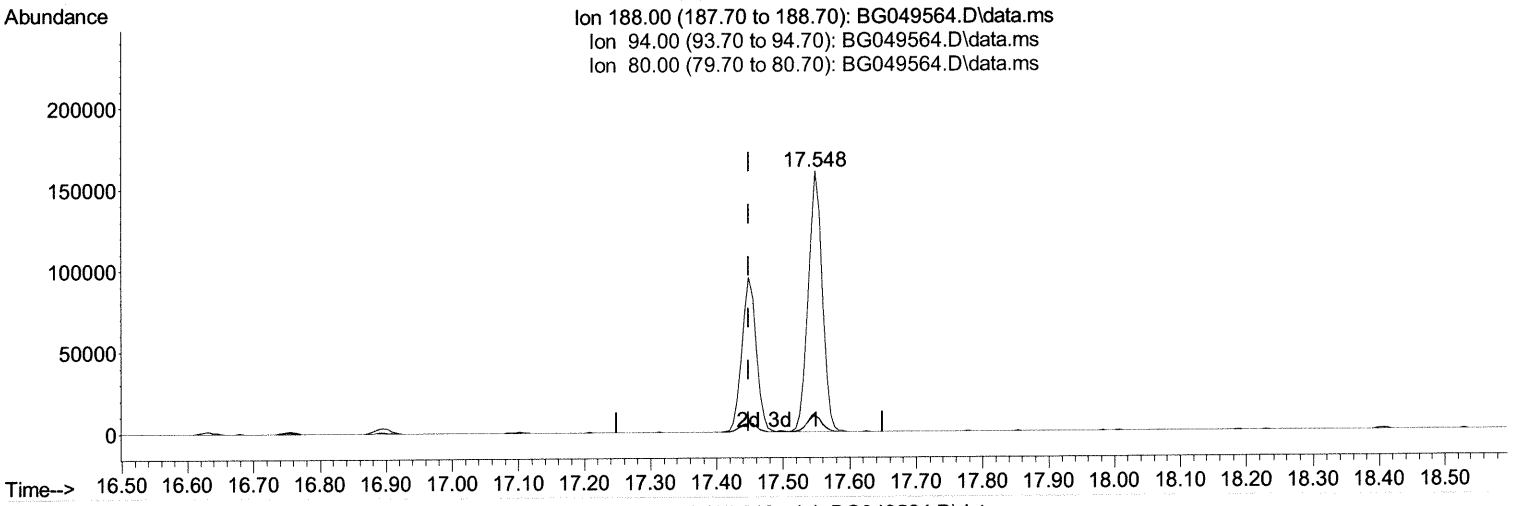
Quantitation Report (Qedit)

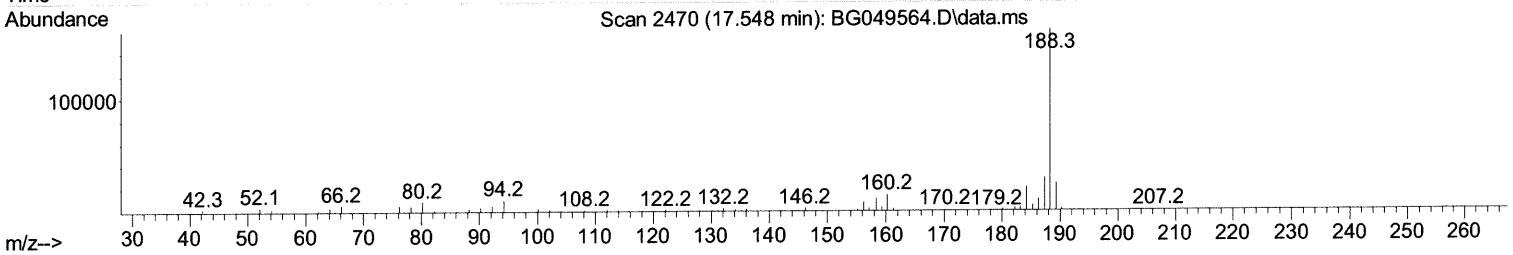
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

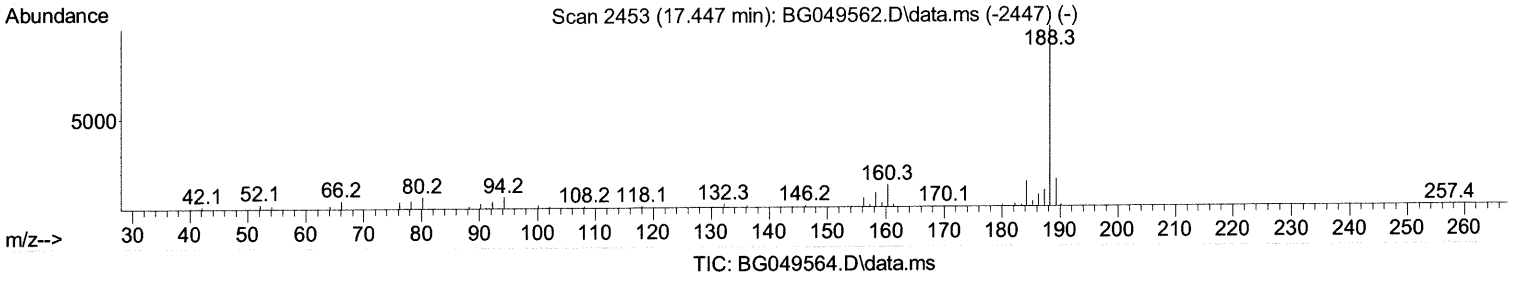
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(64) Phenanthrene-d10 (I)

17.548min (+ 0.100) 20.00 ng/ul

response 235086

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.60	6.62
80.00	7.00	5.78
0.00	0.00	0.00

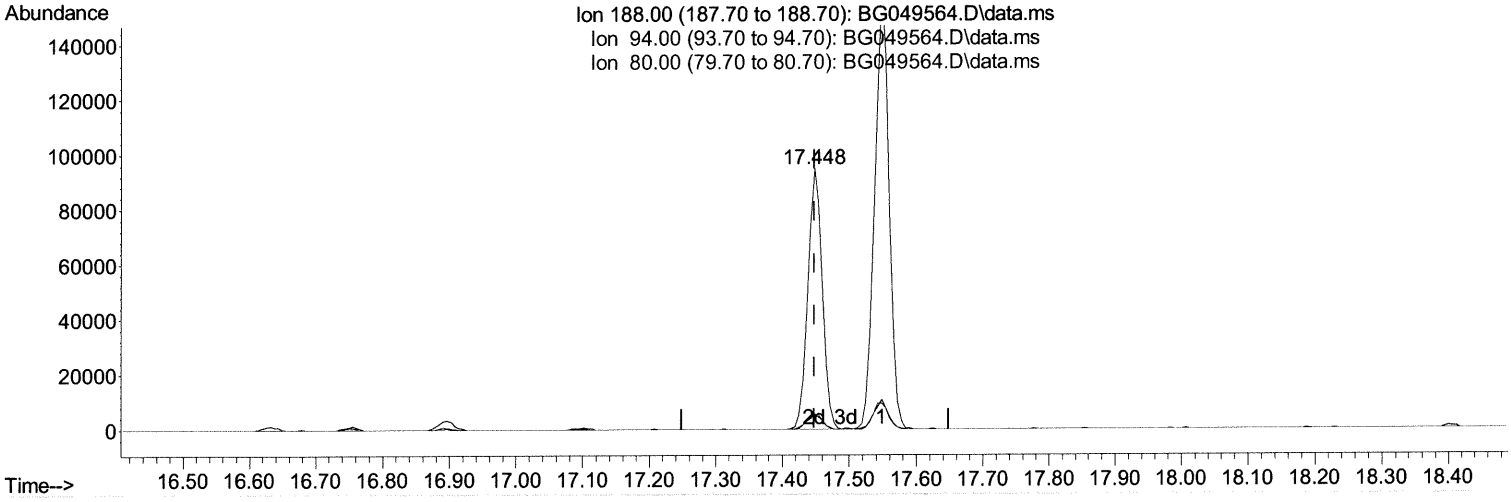
Quantitation Report (Qedit)

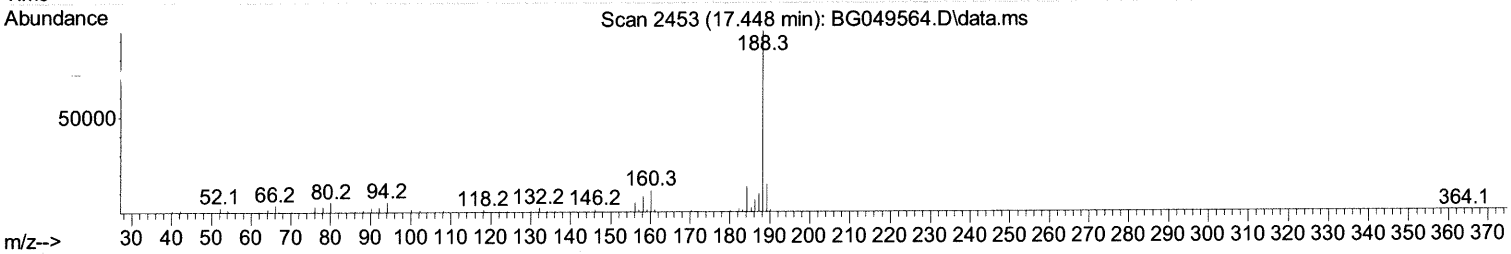
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

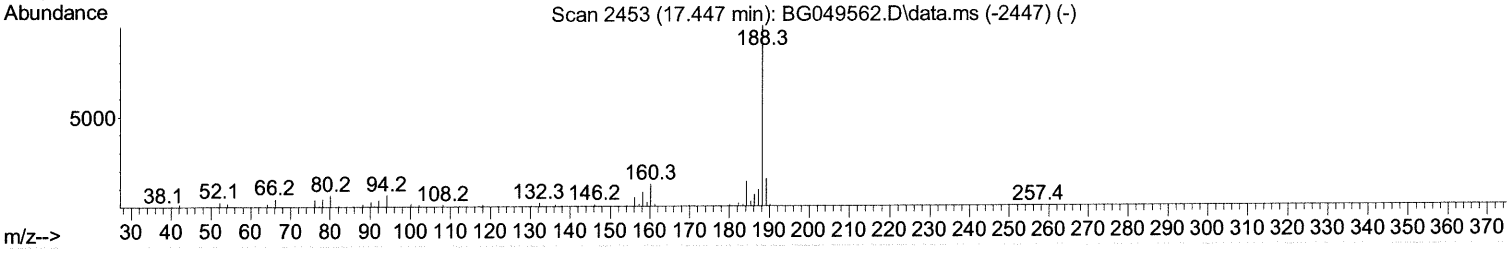
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(64) Phenanthrene-d10 (I)

17.448min (-0.000) 20.00 ng/ul m

response 141503

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.60	5.72
80.00	7.00	5.57#
0.00	0.00	0.00

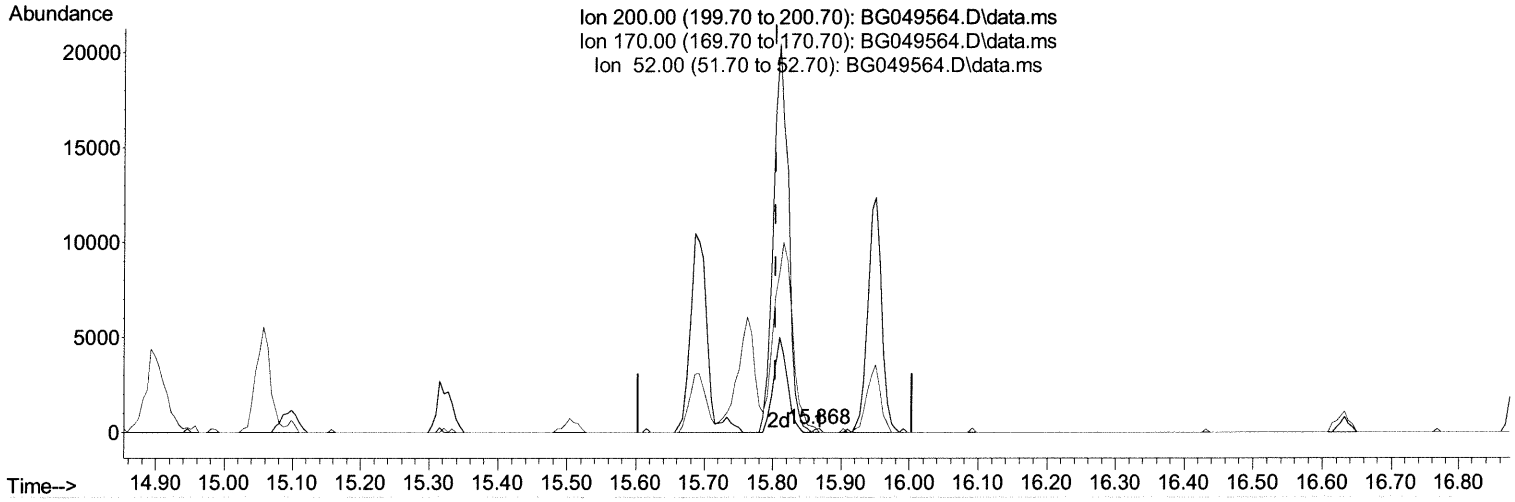
Quantitation Report (Qedit)

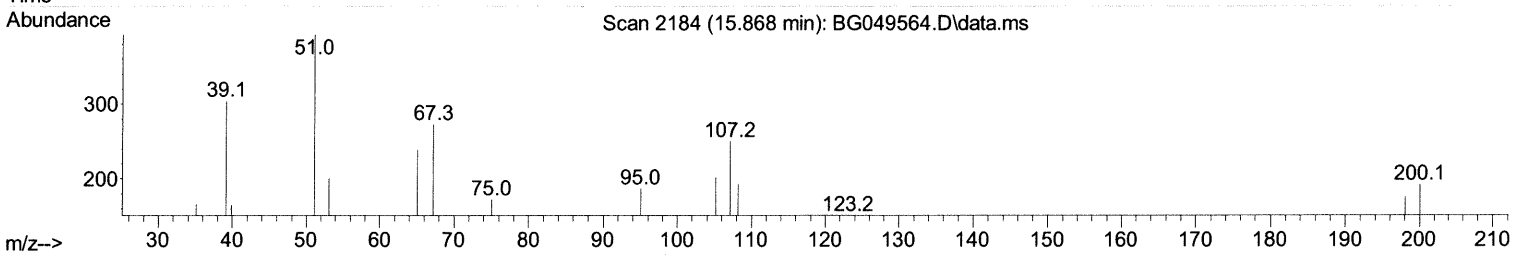
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

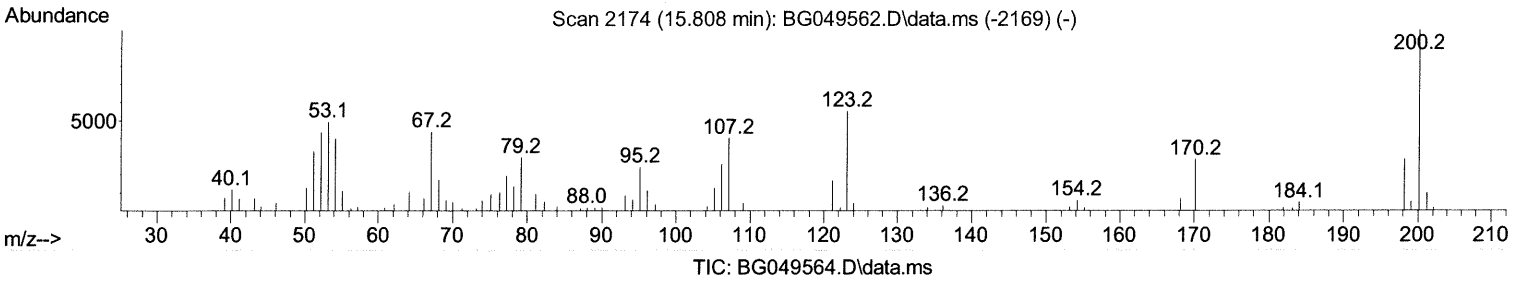
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.868min (+ 0.064) 0.14 ng/ul

response 121

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	25.70	0.00#
52.00	52.30	0.00#
0.00	0.00	0.00

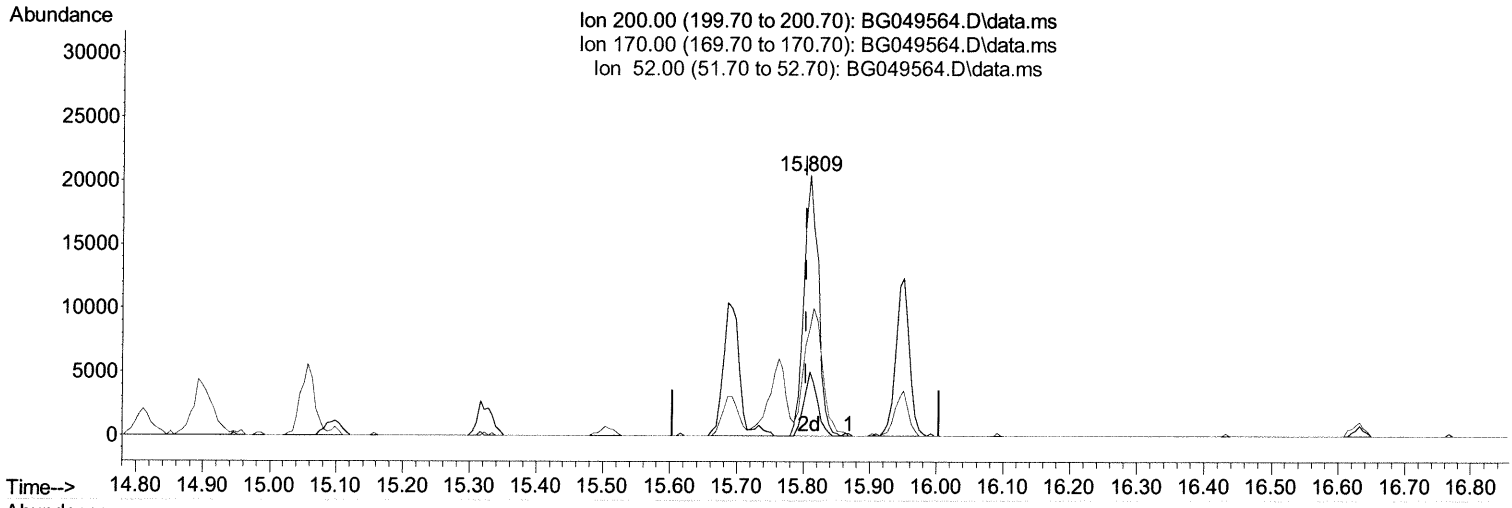
Quantitation Report (Qedit)

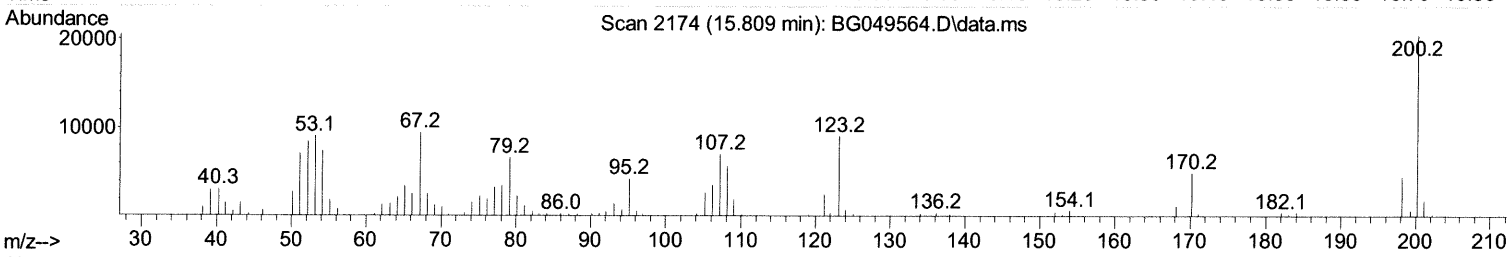
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

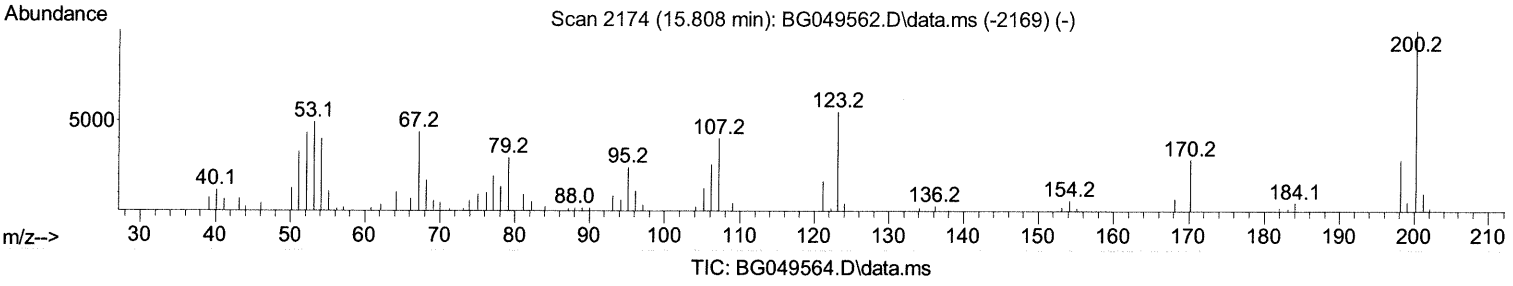
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.809min (+ 0.006) 35.37 ng/ul m 08/02/21 JU

response 31015

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	25.70	24.42
52.00	52.30	41.32#
0.00	0.00	0.00

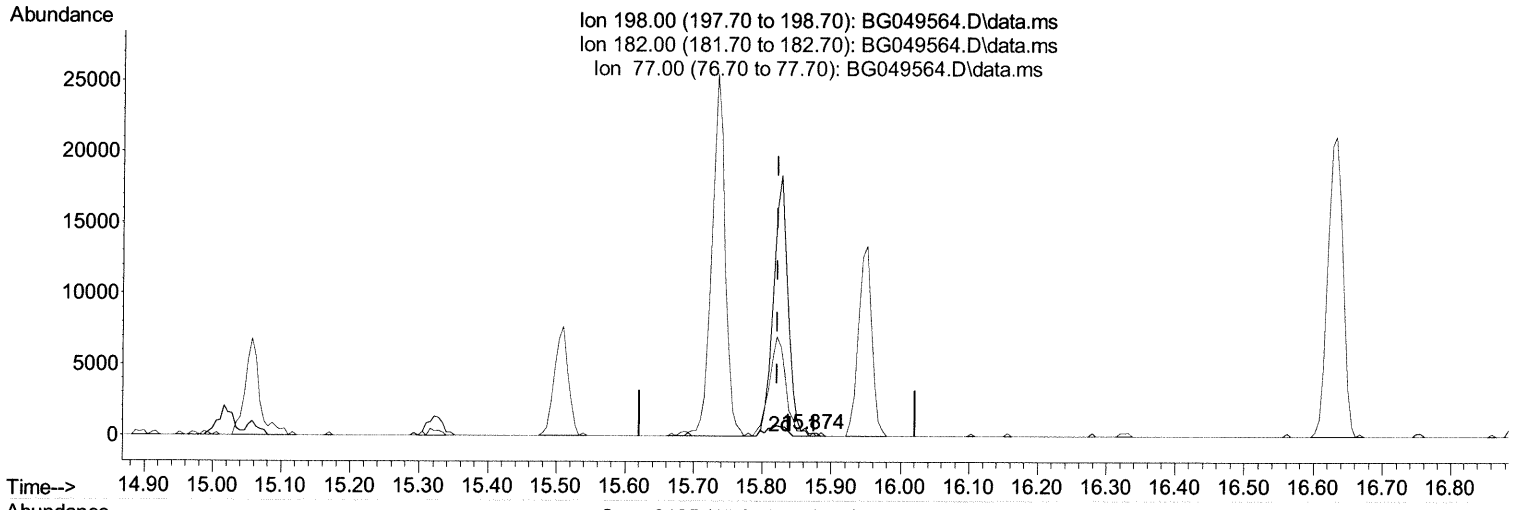
Quantitation Report (Qedit)

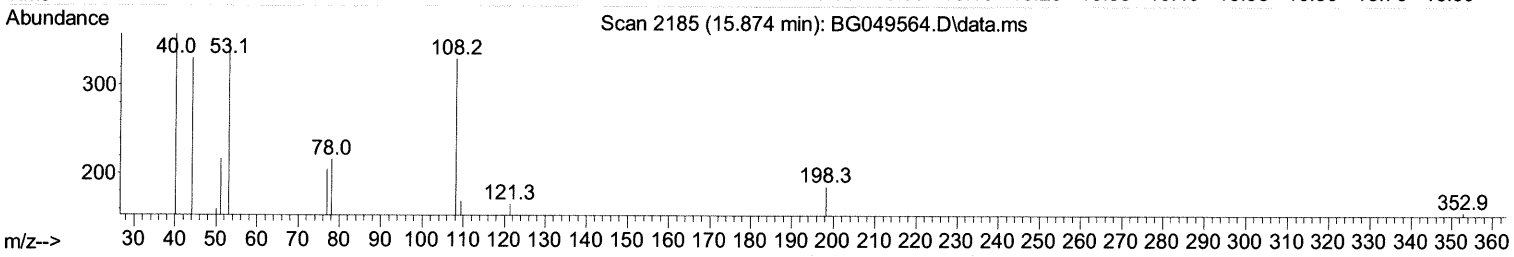
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

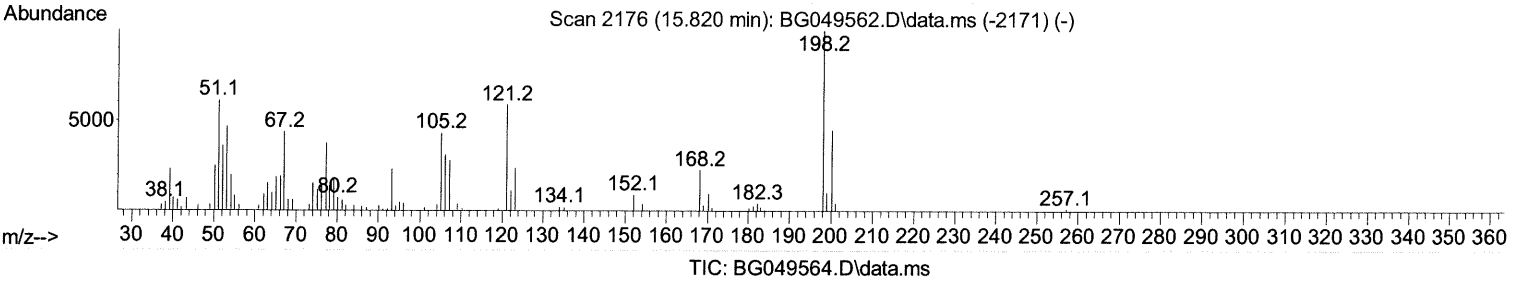
Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(66) 4,6-Dinitro-2-methylphenol

15.874min (+ 0.053) 0.16 ng/ul

response 131

Ion	Exp%	Act%
198.00	100.00	100.00
182.00	5.30	0.00#
77.00	26.00	110.22#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
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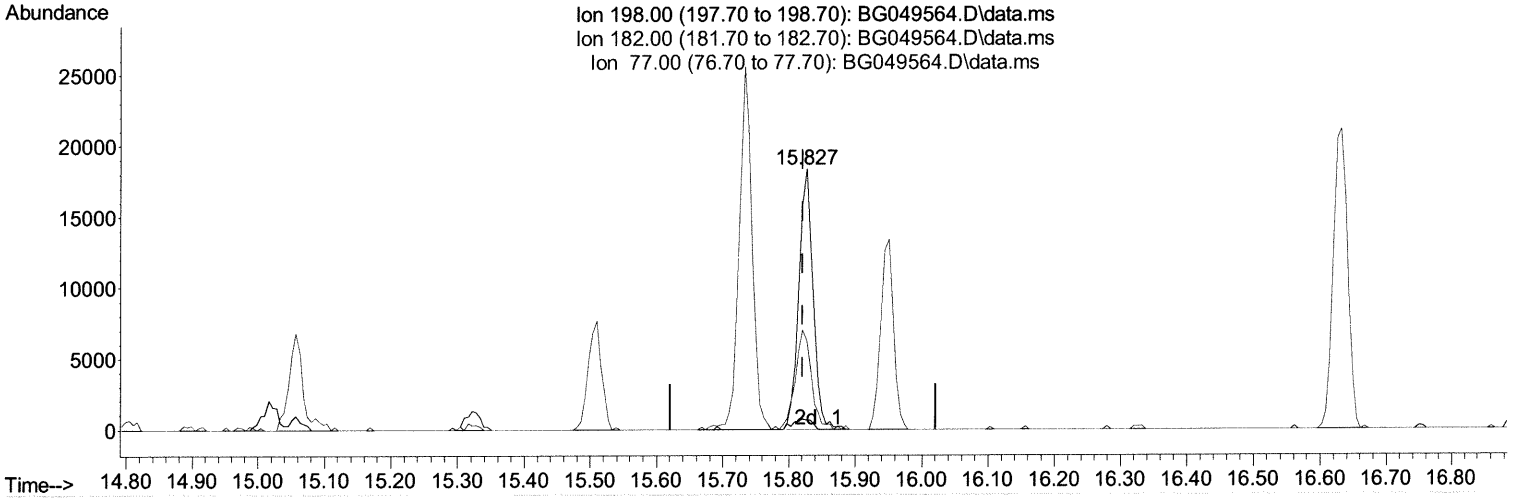
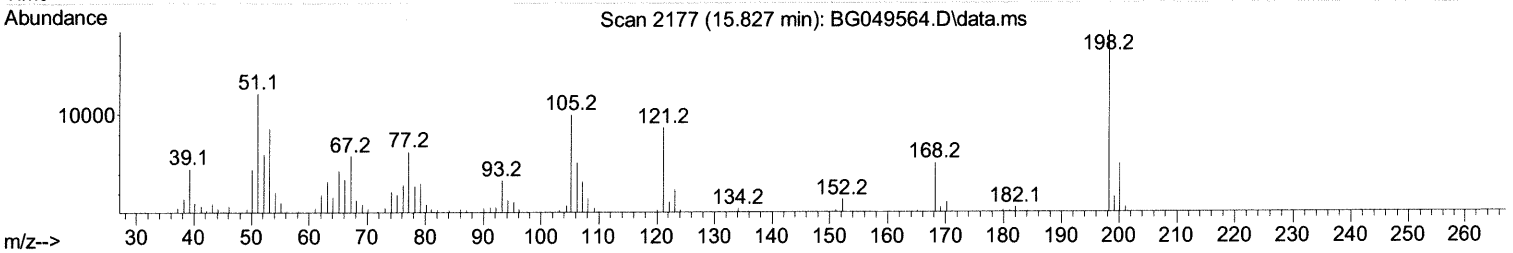
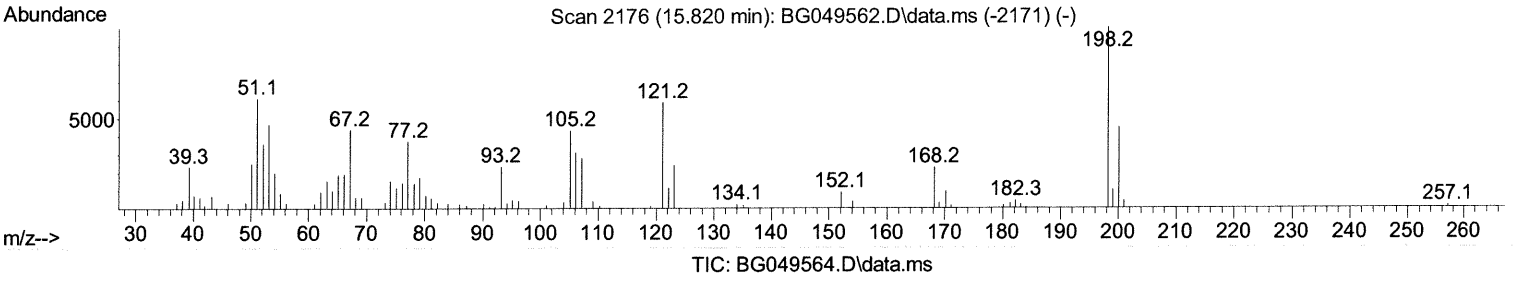
Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(66) 4,6-Dinitro-2-methylphenol

15.827min (+ 0.006) 33.94 ng/ul m 08/02/21 JU

response 27397

Ion	Exp%	Act%
198.00	100.00	100.00
182.00	5.30	3.49#
77.00	26.00	34.15#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
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Instrument :
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ClientSampleId :
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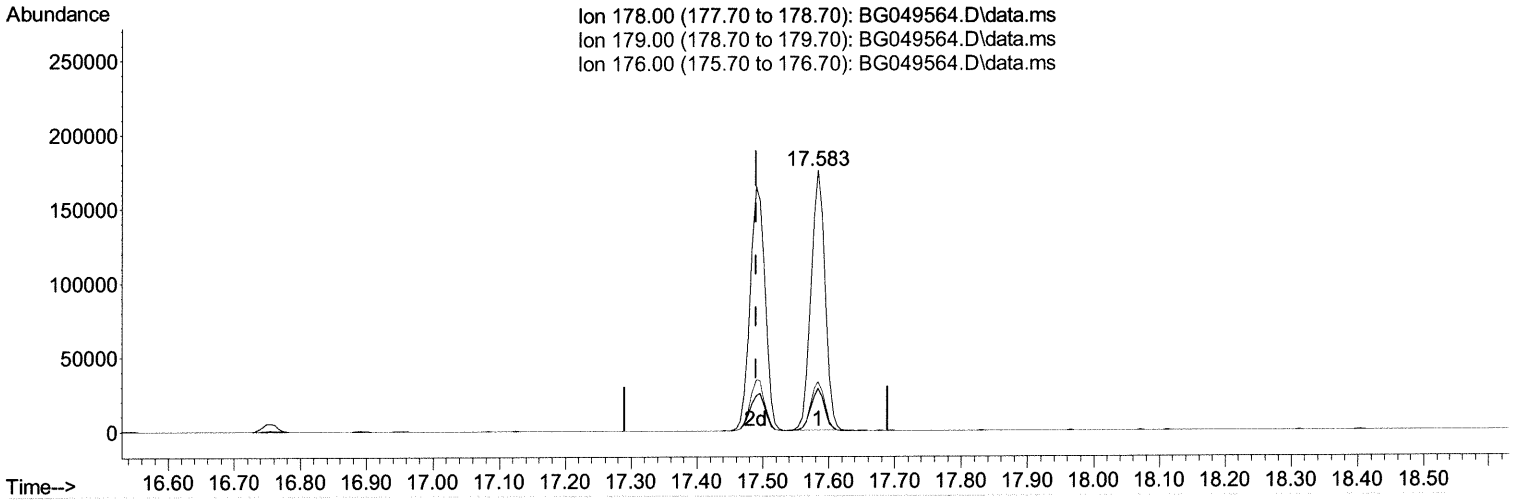
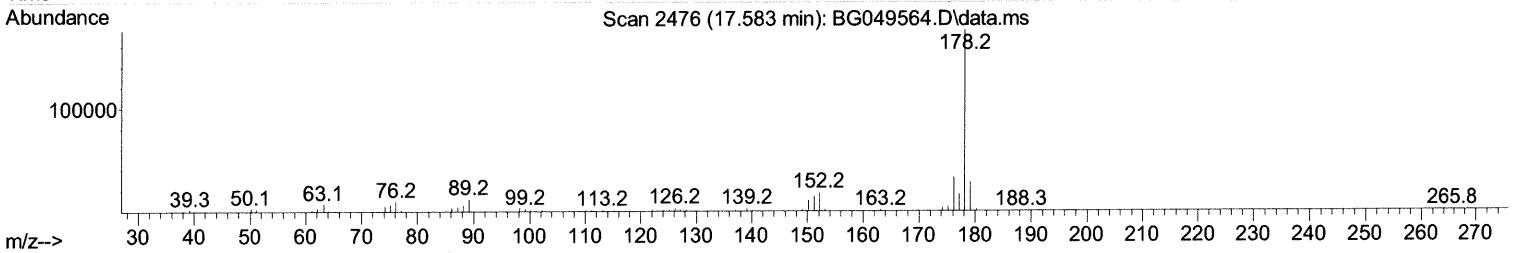
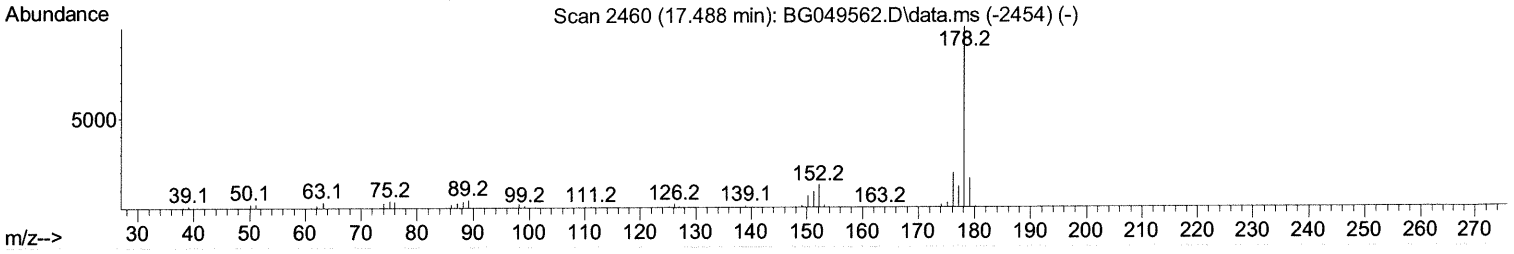
Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(72) Phenanthrene

17.583min (+ 0.094) 35.00 ng/ul

response 257864

Ion	Exp%	Act%
178.00	100.00	100.00
179.00	16.50	16.17
176.00	18.80	18.73
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
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Instrument :
BNA_G
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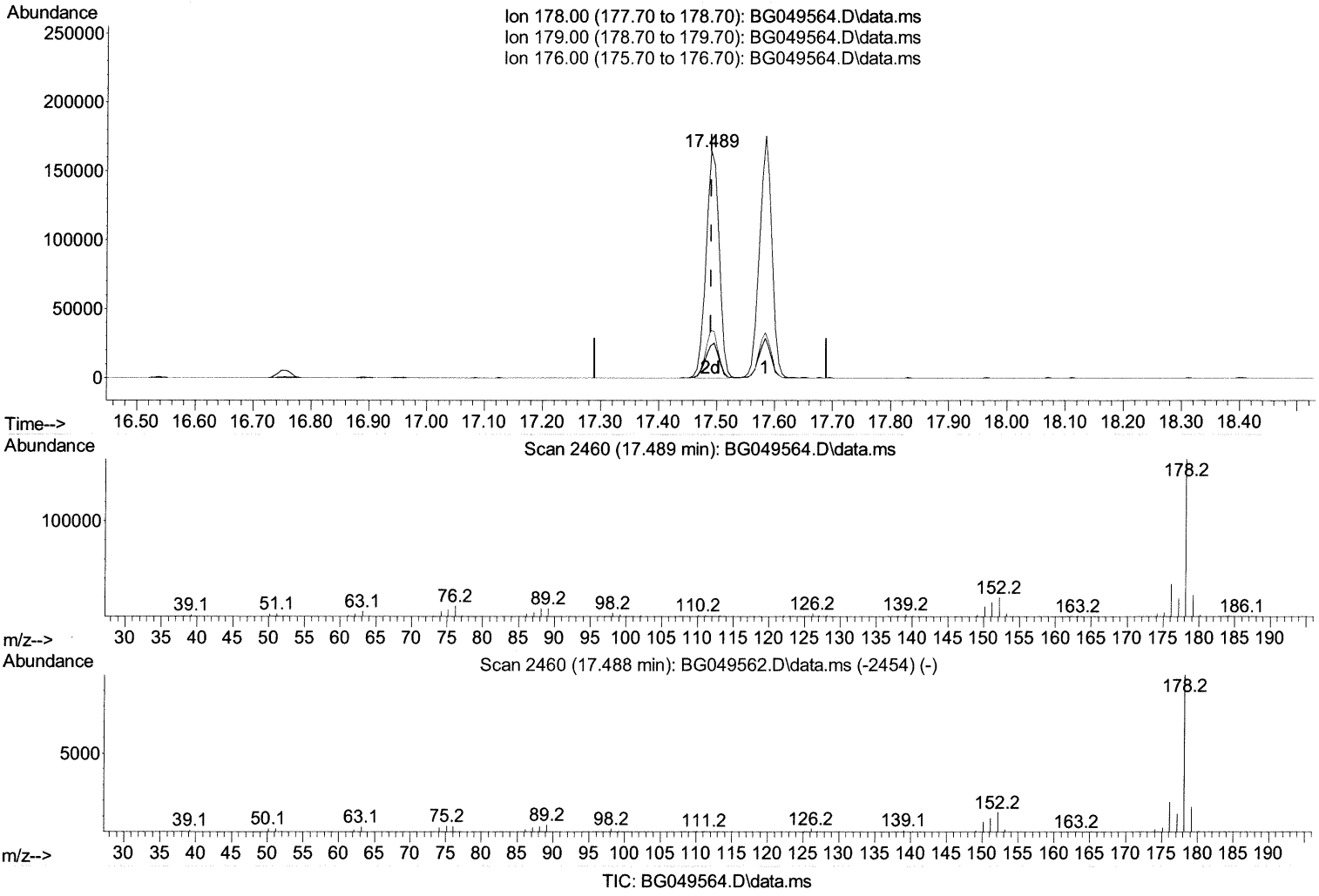
Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM



TIC: BG049564.D\data.ms

(72) Phenanthrene

17.489min (-0.000) 35.32 ng/ul m 08/02/21 JU

response 260216

Ion	Exp%	Act%
178.00	100.00	100.00
179.00	16.50	14.14
176.00	18.80	20.82
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
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Instrument :
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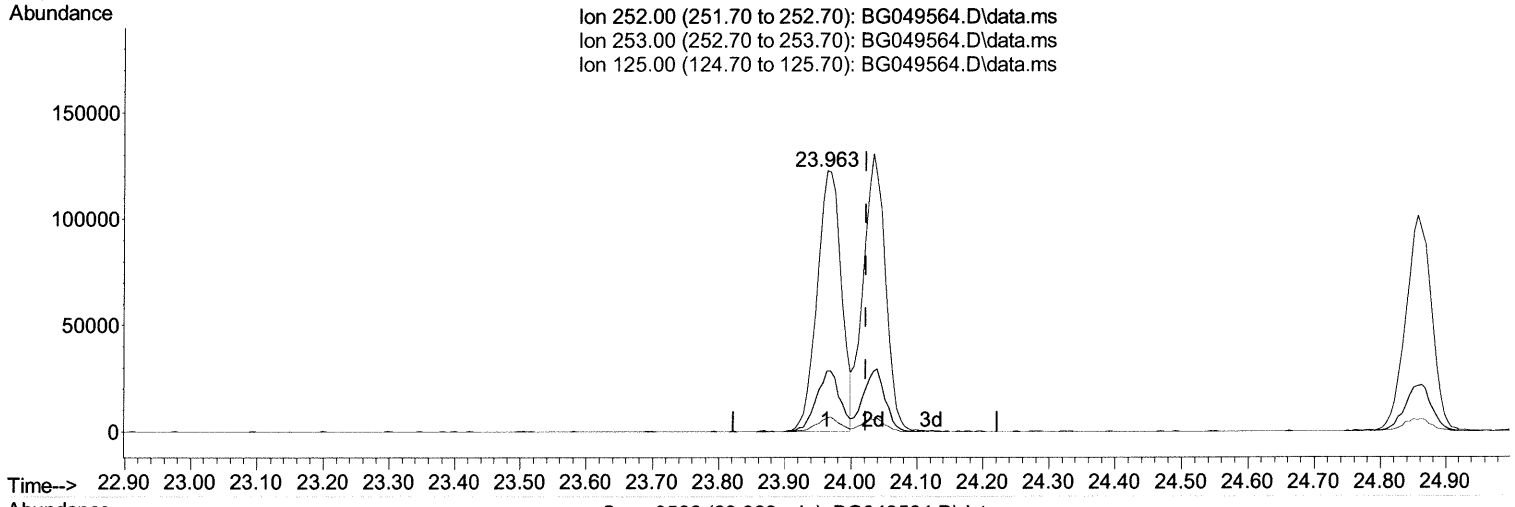
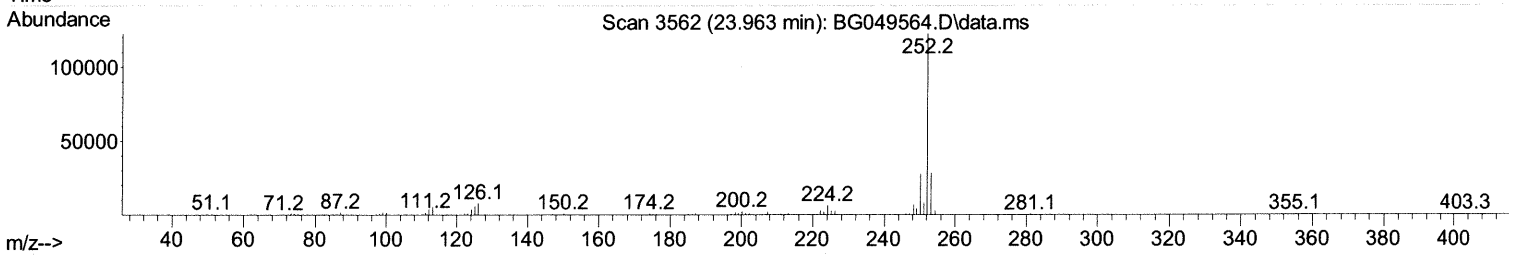
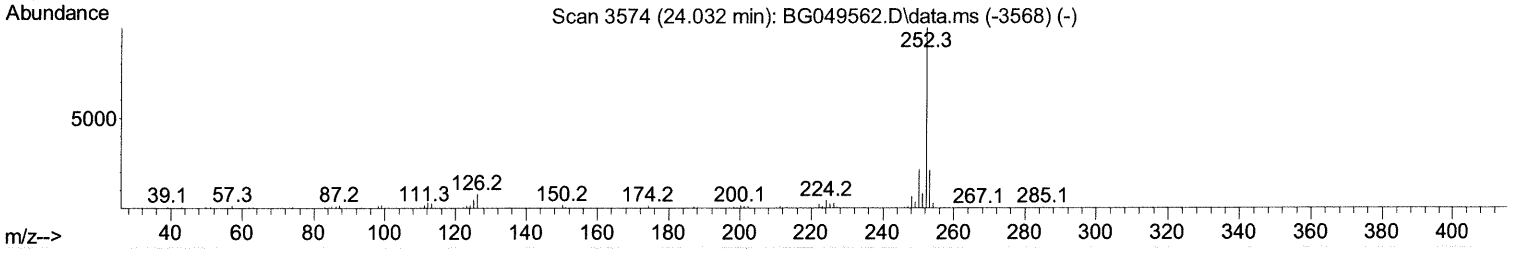
Manual Integrations APPROVED
mohammad
7/30/2021 12:00:12 PM







TIC: BG049564.D\data.ms

(91) Benzo(k)fluoranthene

23.963min (-0.059) 36.08 ng/ul

response 311181

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.60	23.29
125.00	5.80	5.03
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 30 05:25:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SLCS967

Manual Integrations APPROVED
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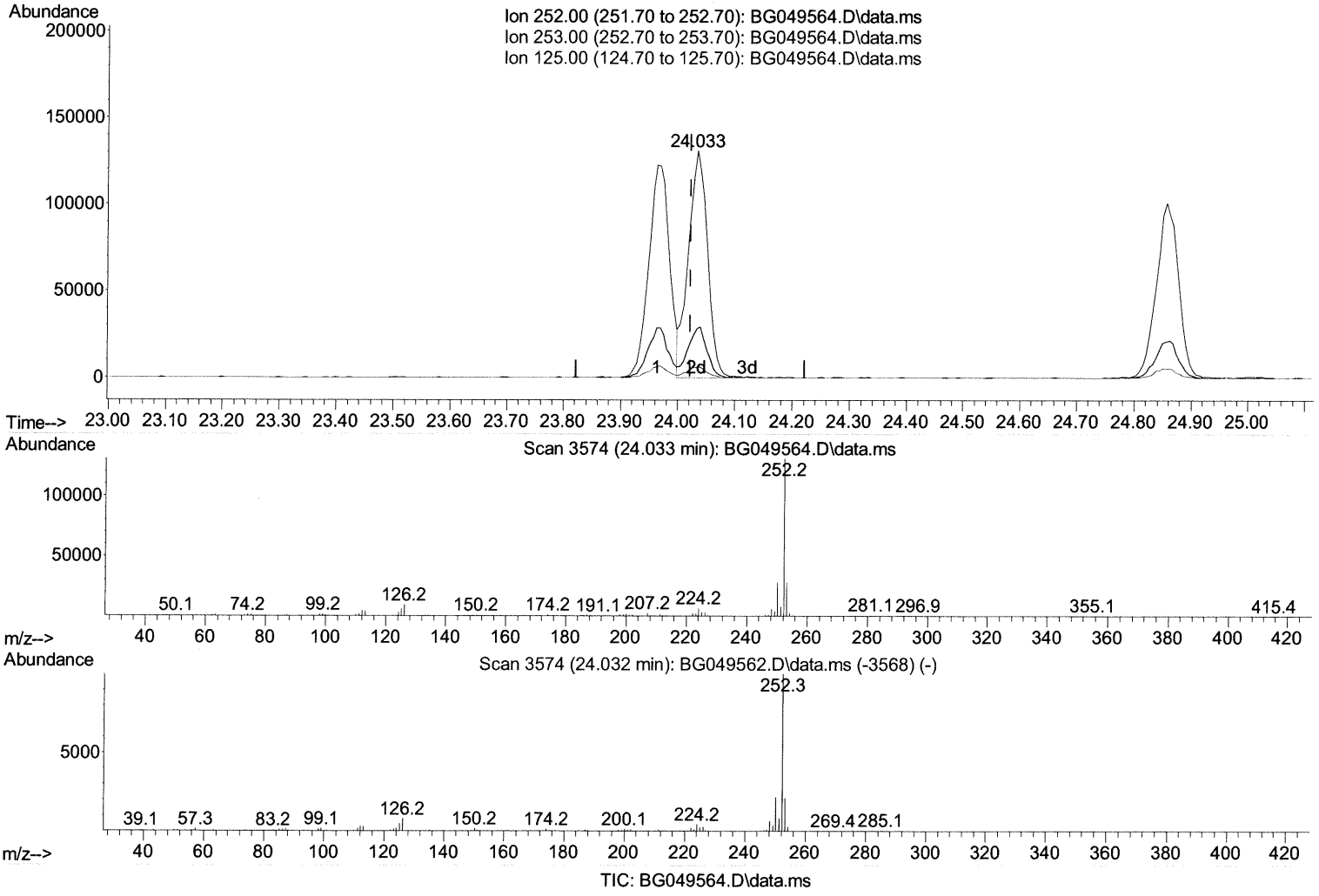
7/30/2021 12:00:12 PM



TIC: BG049564.D\data.ms

(91) Benzo(k)fluoranthene

24.033min (+ 0.012) 35.27 ng/ul m 08/02/21 JU

response 304192

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.60	21.53
125.00	5.80	4.40#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049564.D
Acq On : 29 Jul 2021 18:01
Operator : CG/JU
Sample : PB137967BS
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jul 30 05:25:08 2021
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Quant Title : SVOA CALIBRATION
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ClientSampleId : SLCS967

Manual Integrations APPROVED
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Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)	
Internal Standards								
1) 1,4-Dichlorobenzene-d4	8.049	152	17661	20.000	ng/ul		0.00	
20) Naphthalene-d8	10.869	136	79910	20.000	ng/ul	#	0.00	
38) Acenaphthene-d10	14.699	164	60051	20.000	ng/ul		0.00	
64) Phenanthrene-d10	17.448	188	141503m	20.000	ng/ul	>	0.00	08/02/21 JU
79) Chrysene-d12	21.731	240	135158	20.000	ng/ul		0.00	
88) Perylene-d12	25.009	264	140619	20.000	ng/ul		0.00	
System Monitoring Compounds								
3) 1,4-Dioxane-d8	3.432	96	2716	7.036	ng/uL		0.00	
4) Pyridine-d5	3.849	84	37809	34.100	ng/ul		0.00	
7) Phenol-d5	7.203	99	58058	37.156	ng/ul		0.00	
9) Bis-(2-Chloroethyl)eth...	7.368	67	30611	34.590	ng/ul		0.00	
11) 2-Chlorophenol-d4	7.579	132	42214	37.689	ng/ul		0.00	
15) 4-Methylphenol-d8	8.754	113	44417	37.712	ng/ul		0.00	
21) Nitrobenzene-d5	9.212	128	21186	34.484	ng/ul		0.00	
24) 2-Nitrophenol-d4	9.941	143	25559	35.501	ng/ul		0.00	
28) 2,4-Dichlorophenol-d3	10.487	165	44933	34.316	ng/ul		0.00	
31) 4-Chloroaniline-d4	10.998	131	56059	30.720	ng/ul		0.00	
46) Dimethylphthalate-d6	14.094	166	157963	34.340	ng/ul		0.00	
49) Acenaphthylene-d8	14.394	160	191756	35.890	ng/ul		0.00	
54) 4-Nitrophenol-d4	14.893	143	27520	36.329	ng/ul		0.00	
60) Fluorene-d10	15.692	176	143393	36.184	ng/ul		0.00	
65) 4,6-Dinitro-2-methylph...	15.809	200	31015m	35.371	ng/ul	>	0.00	08/02/21 JU
73) Anthracene-d10	17.548	188	235086	35.746	ng/ul		0.00	
81) Pyrene-d10	19.833	212	269455	38.989	ng/ul		0.00	
92) Benzo(a)pyrene-d12	24.785	264	255818	34.613	ng/ul		0.02	
Target Compounds							Qvalue	
2) 1,4-Dioxane	3.467	88	5230	9.912	ng/uL#		70	
5) Pyridine	3.873	79	40867	34.457	ng/ul		98	
6) Benzaldehyde	7.180	77	37982	40.736	ng/ul		98	
8) Phenol	7.227	94	57409	37.611	ng/ul		98	
10) Bis(2-Chloroethyl)ether	7.462	93	37353	35.520	ng/ul		99	
12) 2-Chlorophenol	7.609	128	42369	39.292	ng/ul		97	
13) 2-Methylphenol	8.490	108	45404	39.028	ng/ul		97	
14) 2,2'-oxybis(1-Chloropr...	8.578	45	43614	35.391	ng/ul#		96	
16) Acetophenone	8.872	105	71034	37.473	ng/ul#		91	
17) N-Nitroso-di-n-propyla...	8.854	70	37275	38.133	ng/ul#		88	
18) 4-Methylphenol	8.819	108	47540	38.844	ng/ul		86	
19) Hexachloroethane	9.136	117	18535	38.155	ng/ul#		84	
22) Nitrobenzene	9.259	77	62412	34.023	ng/ul		97	
23) Isophorone	9.782	82	105015	34.164	ng/ul		99	
25) 2-Nitrophenol	9.976	139	25419	37.517	ng/ul#		85	
26) 2,4-Dimethylphenol	10.029	107	57302	35.487	ng/ul		100	
27) Bis(2-Chloroethoxy)met...	10.264	93	52531	34.265	ng/ul		98	
29) 2,4-Dichlorophenol	10.516	162	45107	35.240	ng/ul		99	
30) Naphthalene	10.922	128	141829	34.457	ng/ul		97	
32) 4-Chloroaniline	11.028	127	56092	32.313	ng/ul		92	
33) Hexachlorobutadiene	11.204	225	34141	34.887	ng/ul		90	
34) Caprolactam	11.779	113	16677m	38.309	ng/ul	>		08/02/21 JU
35) 4-Chloro-3-methylphenol	12.150	107	56995	39.574	ng/ul		95	

Quantitation Report (QT Reviewed)

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration

Instrument : BNA_G
ClientSampleId : SLCS967

Manual Integrations APPROVED
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7/30/2021 12:00:12 PM

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
36) 2-Methylnaphthalene	12.526	142	107534	35.695 ng/ul	95
37) 1-Methylnaphthalene	12.743	142	106711	35.054 ng/ul	89
39) 1,2,4,5-Tetrachloroben...	12.890	216	60307	33.789 ng/ul	89
40) Hexachlorocyclopentadiene	12.872	237	20489	19.052 ng/ul	99
41) 2,4,6-Trichlorophenol	13.131	196	40613	37.028 ng/ul#	79
42) 2,4,5-Trichlorophenol	13.207	196	43551	37.181 ng/ul	99
43) 1,1'-Biphenyl	13.530	154	144820	33.866 ng/ul	95
44) 2-Chloronaphthalene	13.577	162	115530	34.719 ng/ul	97
45) 2-Nitroaniline	13.777	65	46440	39.326 ng/ul	96
47) Dimethylphthalate	14.141	163	149230	33.360 ng/ul	99
48) 2,6-Dinitrotoluene	14.270	165	33447	37.465 ng/ul#	90
50) Acenaphthylene	14.417	152	179139	35.110 ng/ul	96
51) 3-Nitroaniline	14.599	138	33105	37.180 ng/ul	97
52) Acenaphthene	14.764	153	127860	34.940 ng/ul	99
53) 2,4-Dinitrophenol	14.811	184	16533	37.170 ng/ul	87
55) 4-Nitrophenol	14.911	109	37645	36.648 ng/ul	97
56) Dibenzofuran	15.098	168	181963	36.825 ng/ul	100
57) 2,4-Dinitrotoluene	15.057	165	49504	39.119 ng/ul#	97
58) 2,3,4,6-Tetrachlorophenol	15.322	232	42034	39.544 ng/ul#	98
59) Diethylphthalate	15.510	149	158086	34.146 ng/ul	96
61) Fluorene	15.745	166	150946	35.984 ng/ul	99
62) 4-Chlorophenyl-phenyle...	15.733	204	79819	36.553 ng/ul	89
63) 4-Nitroaniline	15.762	138	34685	39.283 ng/ul	89
66) 4,6-Dinitro-2-methylph...	15.827	198	27397m	> 33.942 ng/ul	> 08/02/21 JU
67) N-Nitrosodiphenylamine	15.950	169	132426	35.710 ng/ul	99
68) 4-Bromophenyl-phenylether	16.632	248	54825	34.797 ng/ul	91
69) Hexachlorobenzene	16.755	284	63860	35.823 ng/ul	95
70) Atrazine	16.896	200	58307	34.425 ng/ul	99
71) Pentachlorophenol	17.102	266	31075	37.648 ng/ul	91
72) Phenanthrene	17.489	178	260216m	> 35.324 ng/ul	> 08/02/21 JU
74) Anthracene	17.583	178	257412	34.483 ng/ul	98
75) 1,2,3,4-Tetrachloroben...	13.501	216	63123	32.829 ng/uL	98
76) Pentachlorobenzene	15.016	250	72452	35.614 ng/uL	98
77) Carbazole	17.854	167	230660	34.562 ng/ul	97
78) Di-n-butylphthalate	18.400	149	285469	34.397 ng/ul	100
80) Fluoranthene	19.498	202	315421	37.588 ng/ul	98
82) Pyrene	19.863	202	312241	37.477 ng/ul	99
83) Butylbenzylphthalate	20.738	149	123044	37.933 ng/ul	95
84) 3,3'-Dichlorobenzidine	21.625	252	107470	34.606 ng/ul	99
85) Benzo(a)anthracene	21.713	228	299952	36.299 ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.607	149	175751	36.509 ng/ul	99
87) Chrysene	21.778	228	282358	35.699 ng/ul	98
89) Di-n-octyl phthalate	22.835	149	289606	35.036 ng/ul	100
90) Benzo(b)fluoranthene	23.963	252	311181	35.307 ng/ul	96
91) Benzo(k)fluoranthene	24.033	252	304192m	> 35.266 ng/ul	> 08/02/21 JU
93) Benzo(a)pyrene	24.856	252	273110	35.719 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	28.762	276	357620	36.656 ng/ul	98
95) Dibenzo(a,h)anthracene	28.833	278	294162	36.436 ng/ul	99
96) Benzo(g,h,i)perylene	29.937	276	293995	36.245 ng/ul#	92

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