

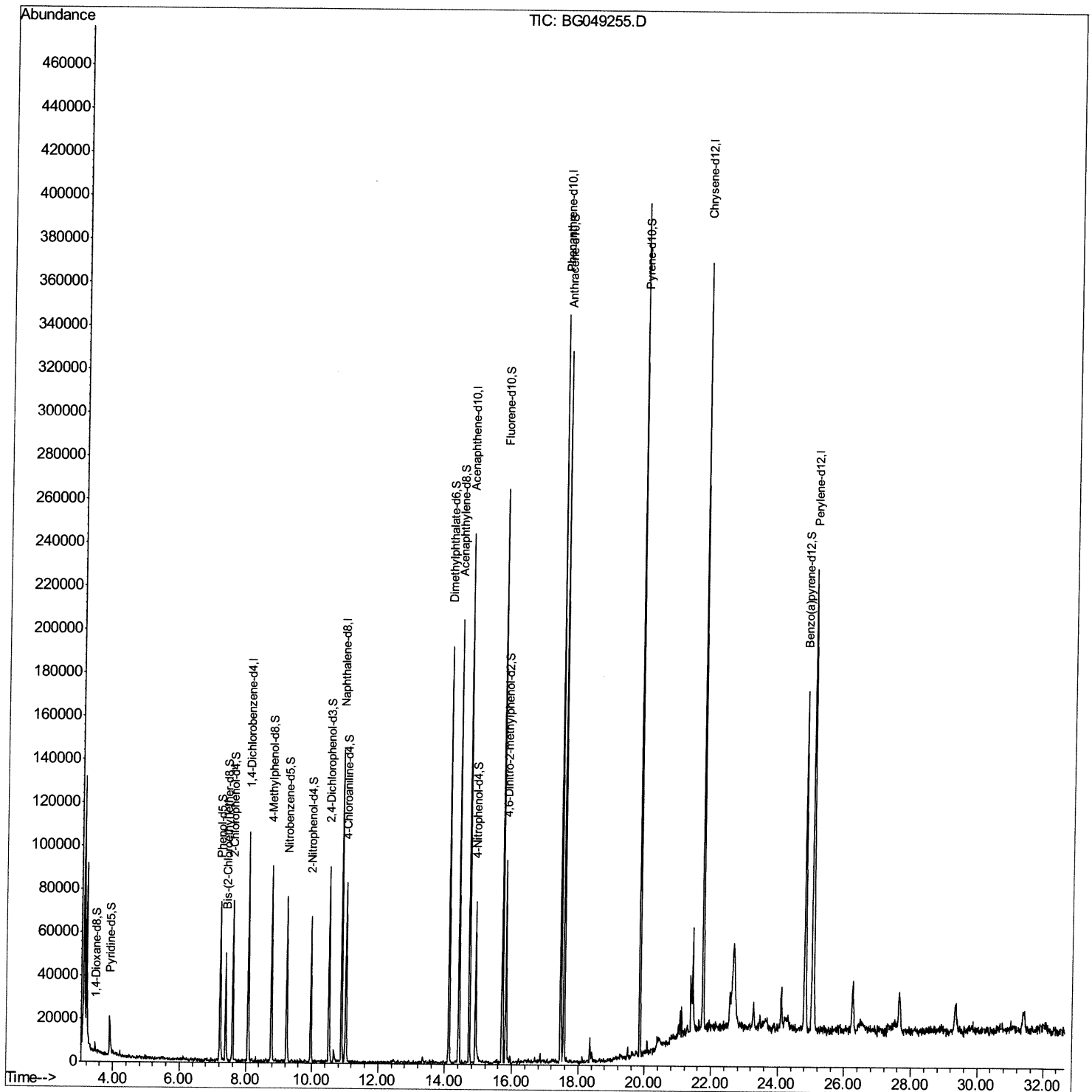
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
Data File : BG049255.D
Acq On : 10 Jul 2021 2:54
Operator : CG/JU
Sample : M2878-17
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGE09

Manual Integrations
APPROVED

mohammad
7/12/2021 12:34:05 PM

Quant Time: Jul 10 03:37:31 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 09 22:53:59 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

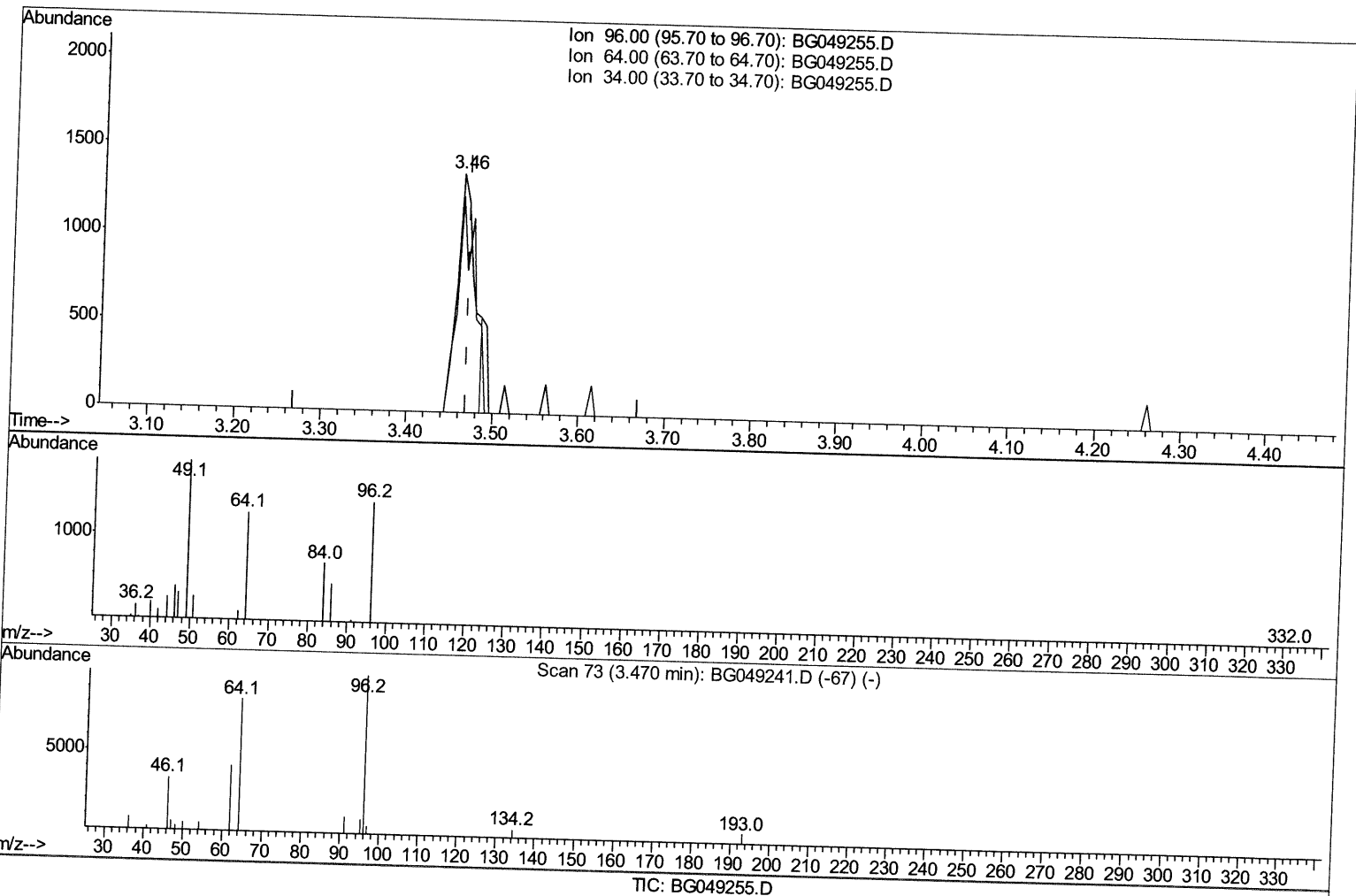
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(3) 1,4-Dioxane-d8 (S)

3.461min (-0.008) 3.08ng/uL

response 1924

Ion	Exp%	Act%
96.00	100	100
64.00	62.20	90.35#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

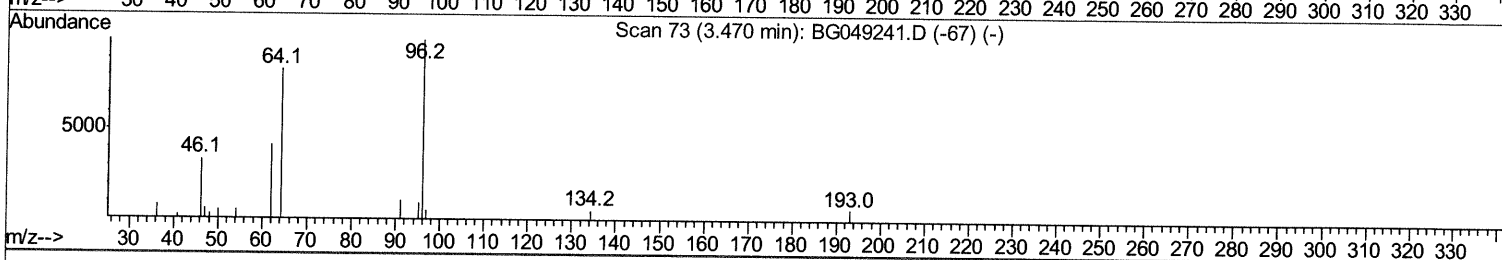
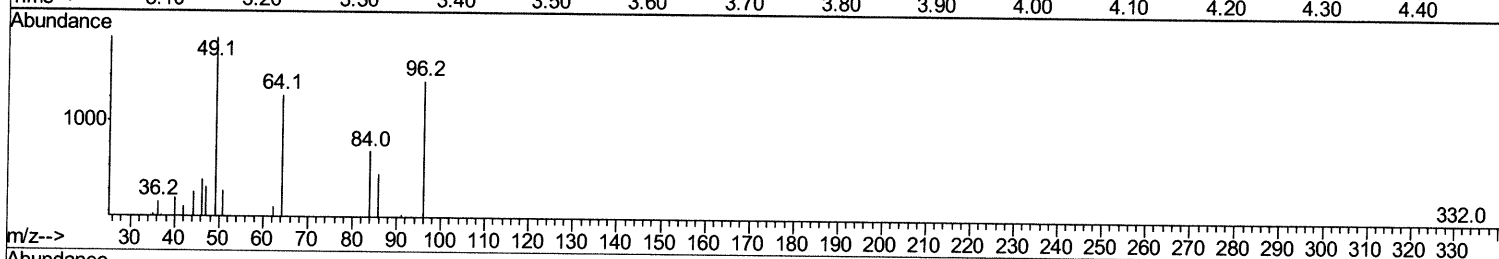
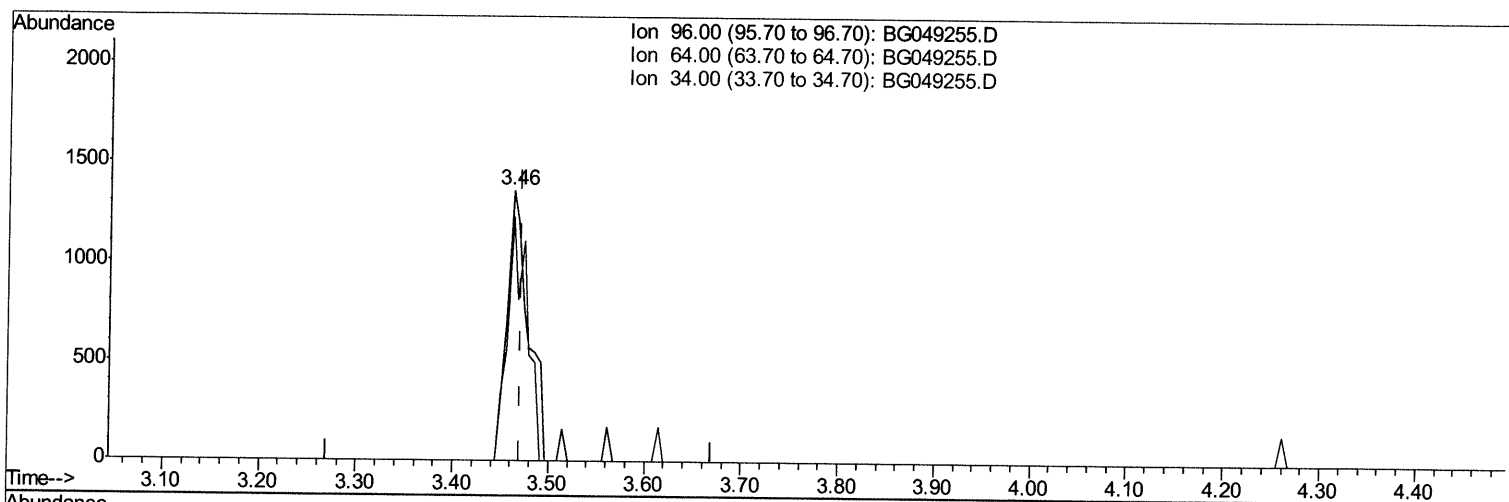
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TIC: BG049255.D

(3) 1,4-Dioxane-d8 (S)

3.461min (-0.008) 3.36ng/uL m 07/12/21 JU

response 2097

Ion	Exp%	Act%
96.00	100	100
64.00	62.20	90.35#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

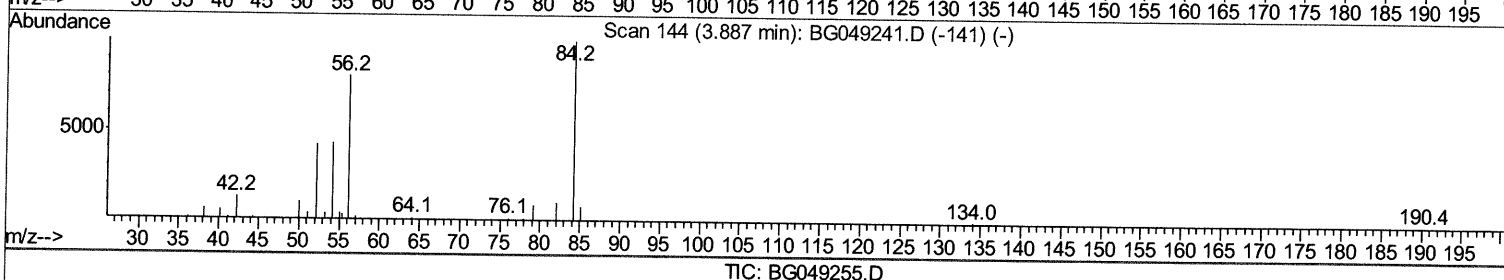
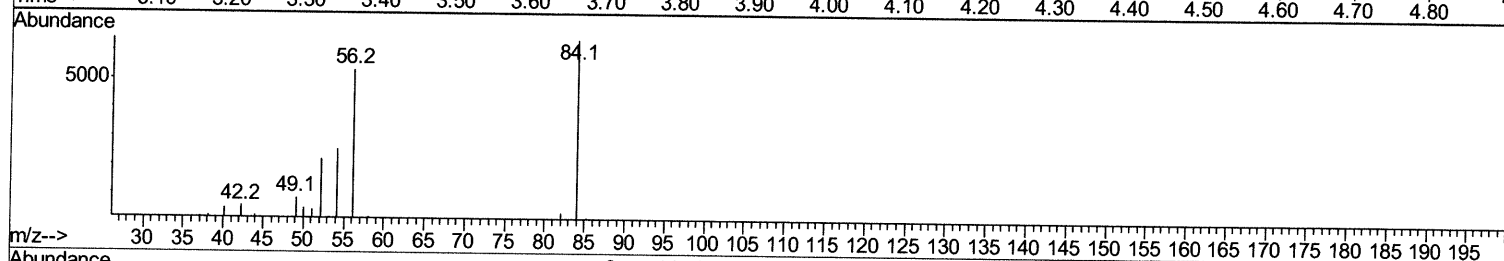
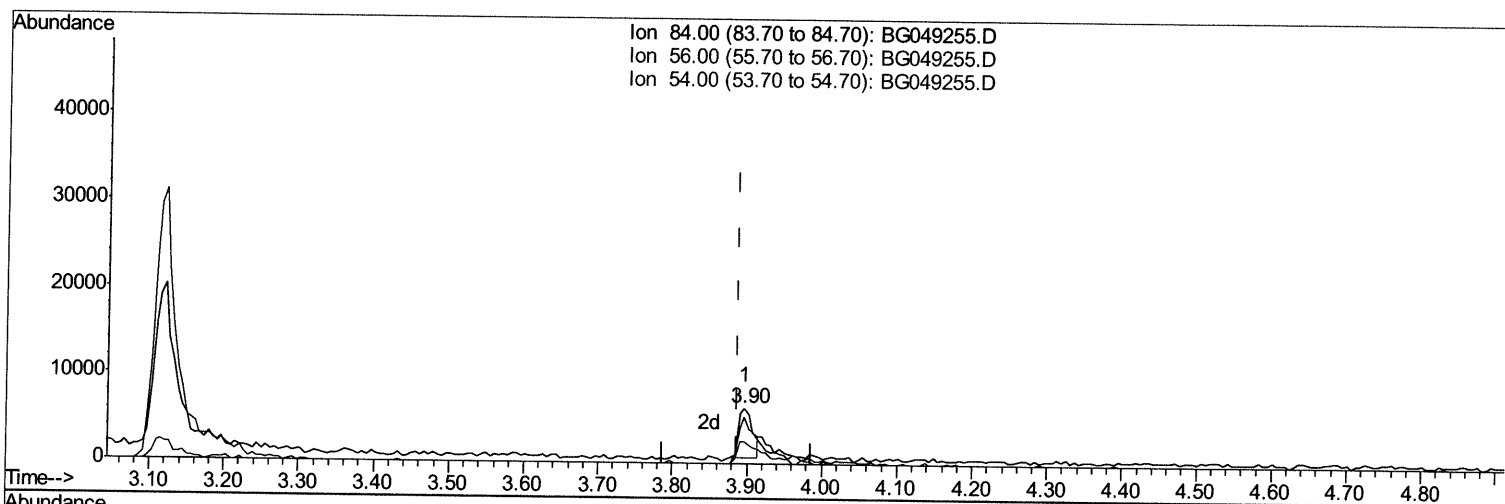
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Manual Integrations
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(4) Pyridine-d5 (S)

3.896min (+0.009) 4.07ng/ul

response 7457

Ion	Exp%	Act%
84.00	100	100
56.00	76.20	83.31
54.00	42.60	40.14
0.00	0.00	0.00

Quantitation Report (Qedit)

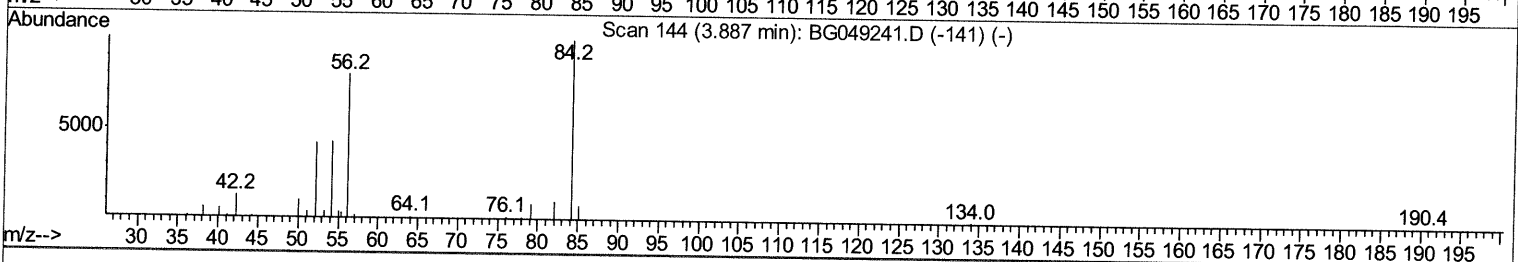
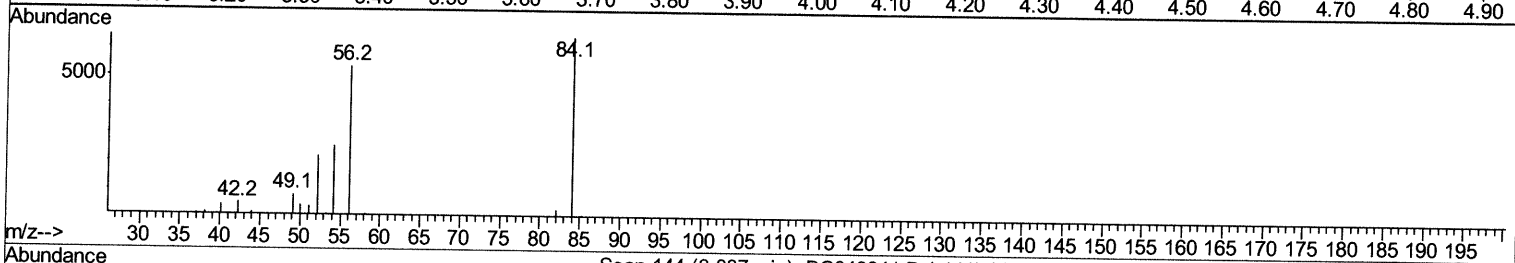
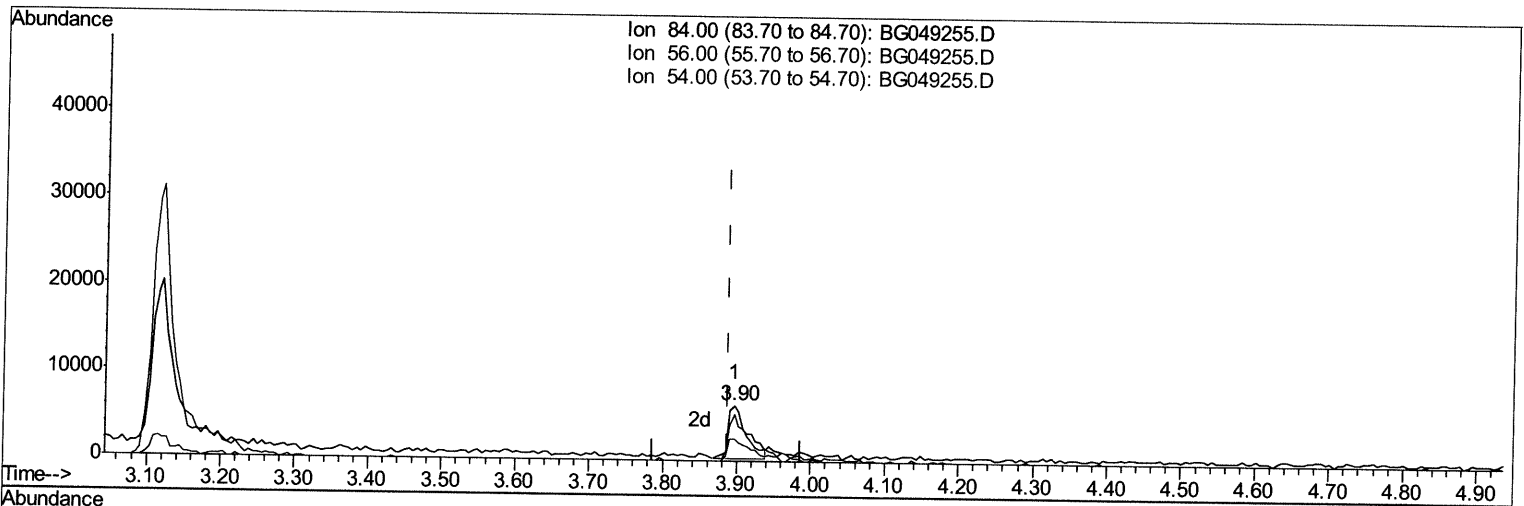
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 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
 APPROVED

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TIC: BG049255.D

(4) Pyridine-d5 (S)

3.896min (+0.009) 6.25ng/ul m 07/13/21 JU.

response 11459

Ion	Exp%	Act%
84.00	100	100
56.00	76.20	83.31
54.00	42.60	40.14
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	29339	20.00	ng/ul	0.00
20) Naphthalene-d8	10.89	136	118618	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.72	164	87763	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.47	188	212249	20.00	ng/ul	0.00
79) Chrysene-d12	21.75	240	228856	20.00	ng/ul	0.00
88) Perylene-d12	25.05	264	231480	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	2097m	3.36	ng/uL	0.00
4) Pyridine-d5	3.90	84	11459m	6.25	ng/ul	0.00
7) Phenol-d5	7.22	99	39602	15.52	ng/ul	0.00
9) Bis-(2-Chloroethyl) ether-d	7.39	67	25148	16.99	ng/ul	0.00
11) 2-Chlorophenol-d4	7.60	132	31908	16.92	ng/ul	0.00
15) 4-Methylphenol-d8	8.77	113	33466	16.37	ng/ul	0.00
21) Nitrobenzene-d5	9.24	128	16324	17.93	ng/ul	0.00
24) 2-Nitrophenol-d4	9.97	143	18401	17.61	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.51	165	33228	16.30	ng/ul	0.00
31) 4-Chloroaniline-d4	11.02	131	42655	16.05	ng/ul	0.00
46) Dimethylphthalate-d6	14.12	166	128267	19.00	ng/ul	0.00
49) Acenaphthylene-d8	14.41	160	141260	17.84	ng/ul	0.00
54) 4-Nitrophenol-d4	14.91	143	17249	15.44	ng/ul	0.00
60) Fluorene-d10	15.71	176	108085	18.91	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.82	200	21594	16.01	ng/ul	0.00
73) Anthracene-d10	17.57	188	193091	19.98	ng/ul	0.00
81) Pyrene-d10	19.85	212	215223	18.35	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.81	264	181268	15.06	ng/ul	0.00

Target Compounds

Ovalue

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