

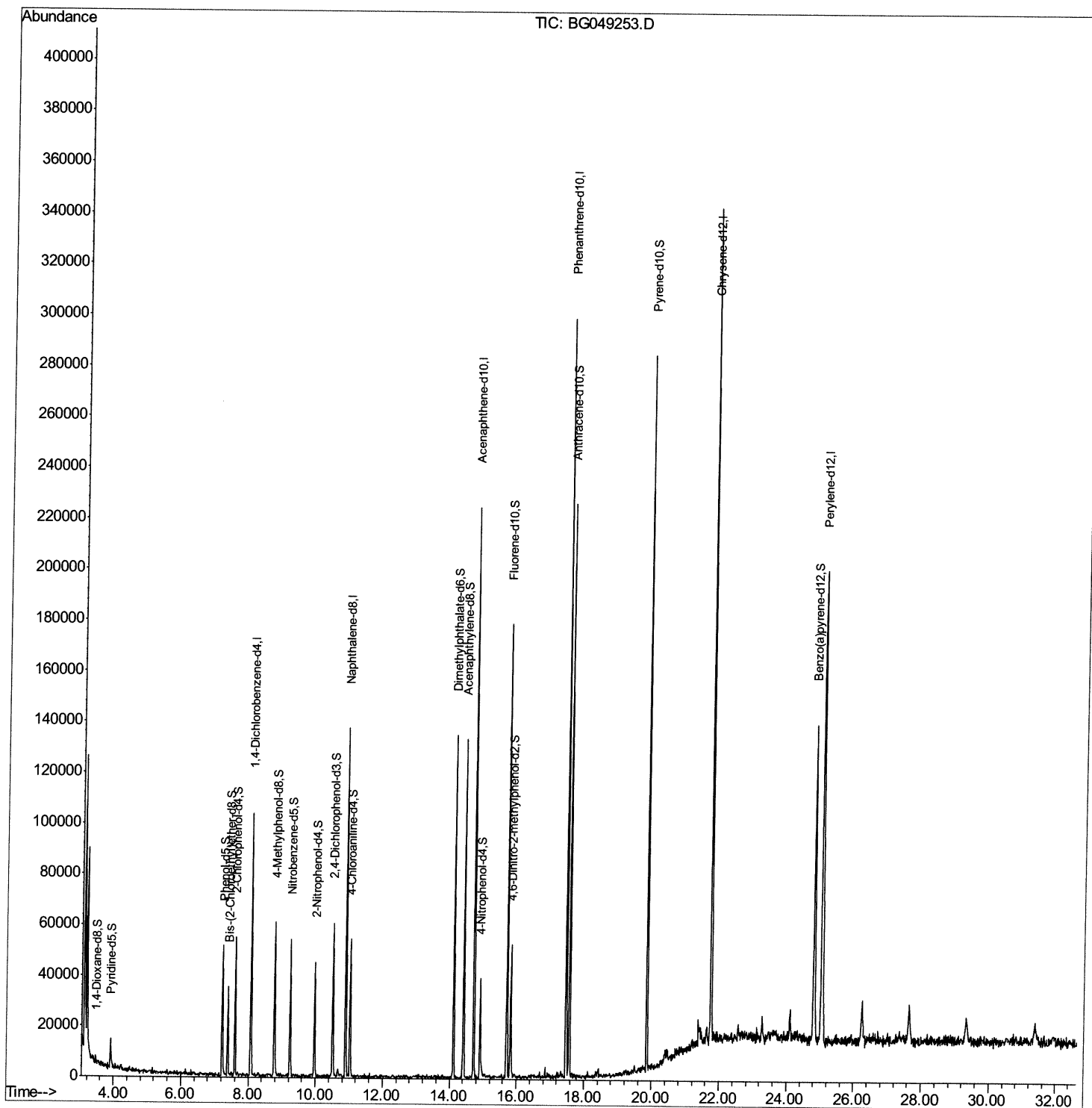
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
Data File : BG049253.D
Acq On : 10 Jul 2021 1:31
Operator : CG/JU
Sample : M2878-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGE08

Manual Integrations
APPROVED

mohammad
7/12/2021 12:33:52 PM

Quant Time: Jul 10 02:15:36 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)

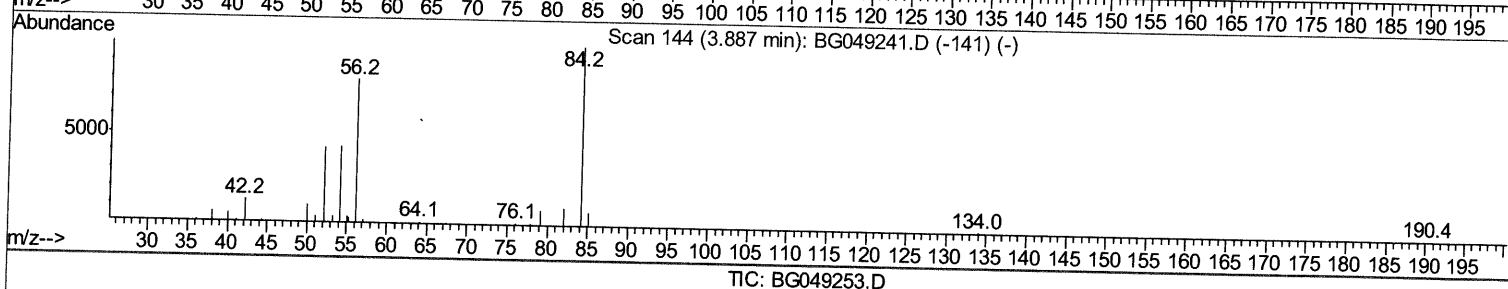
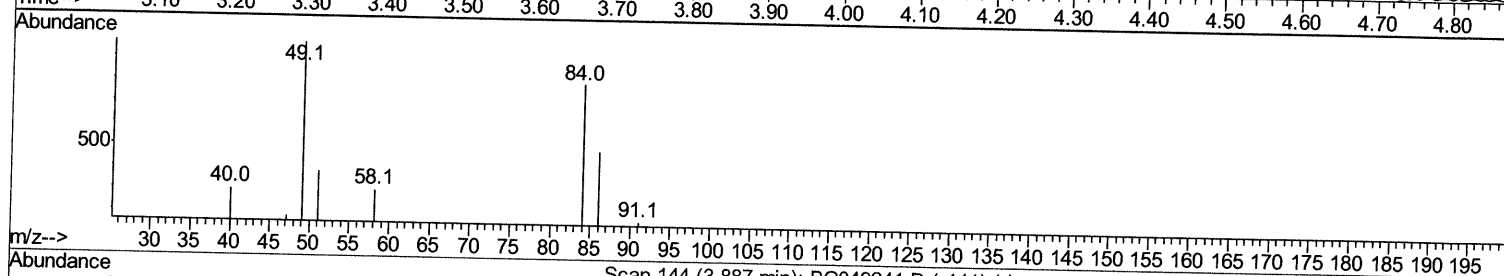
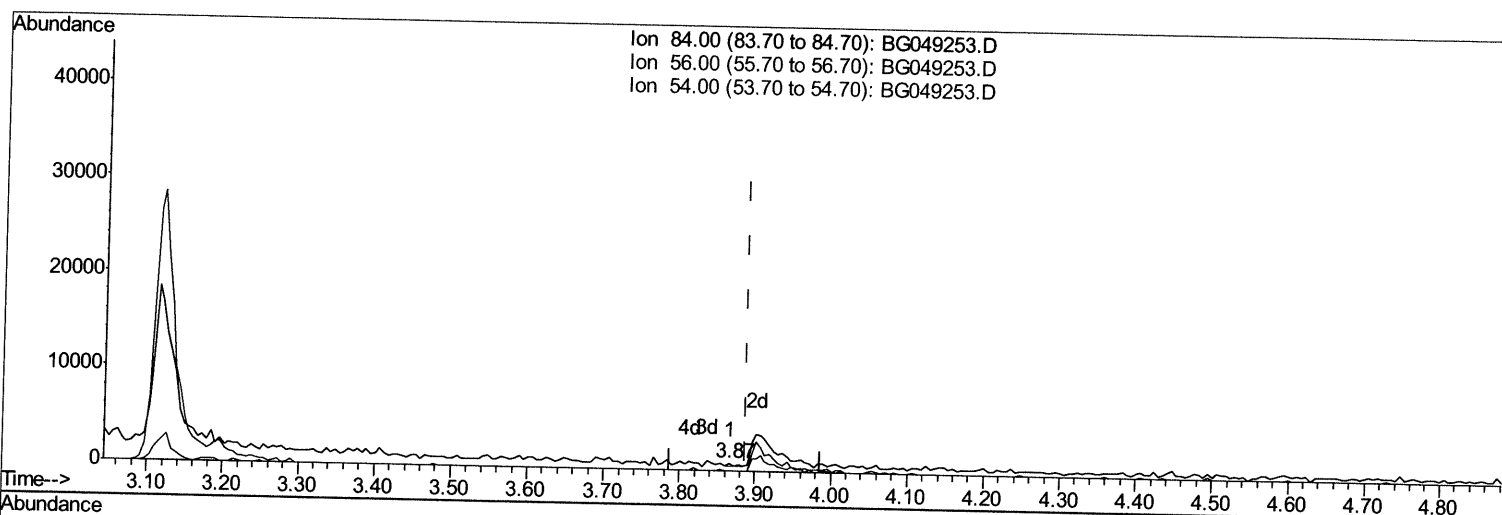
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(4) Pyridine-d5 (S)

3.866min (-0.021) 0.11ng/ul

response 186

Ion	Exp%	Act%
84.00	100	100
56.00	76.20	0.00#
54.00	42.60	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

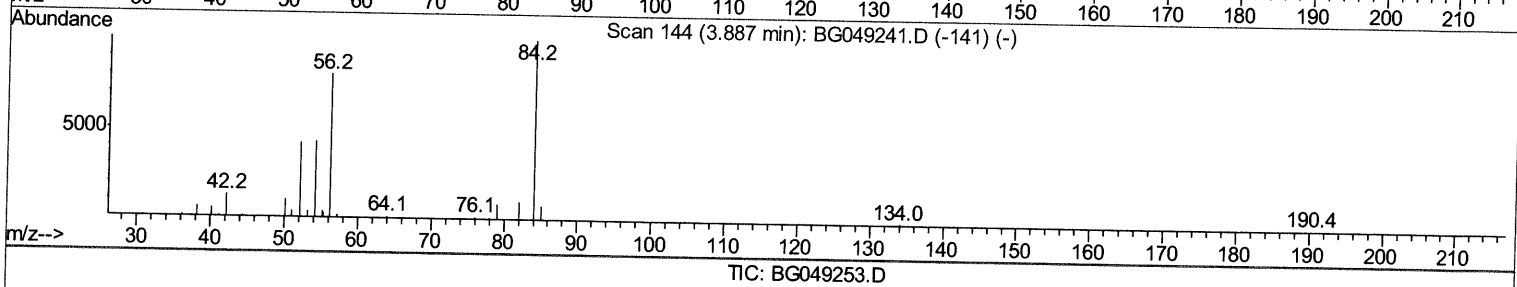
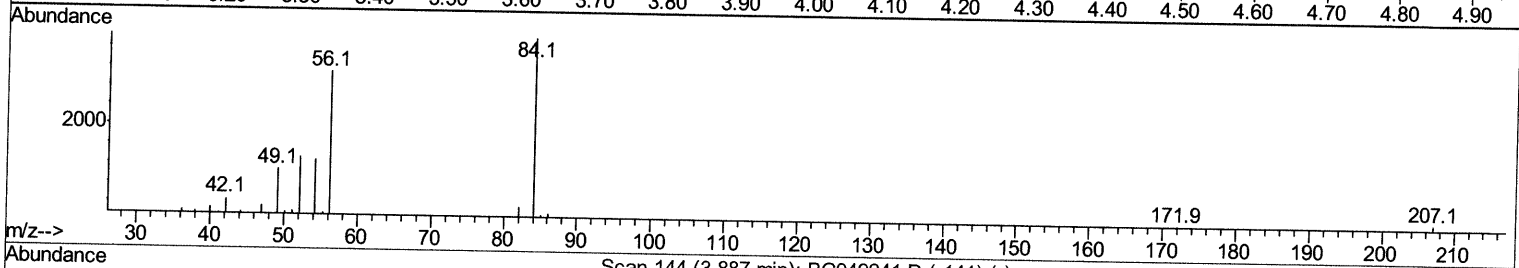
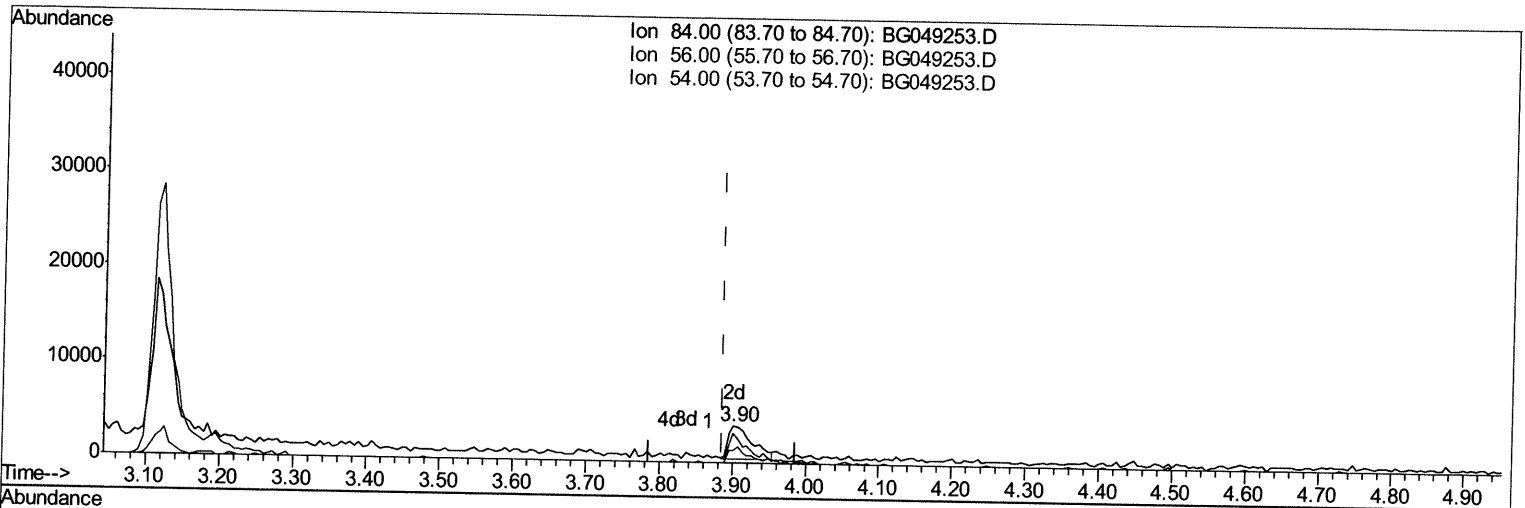
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Manual Integrations
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 Response via : Initial Calibration



TIC: BG049253.D

(4) Pyridine-d5 (S)

3.901min (+0.014) 4.64ng/ul m 07/13/21 JU

response 7682

Ion	Exp%	Act%
84.00	100	100
56.00	76.20	80.72
54.00	42.60	33.77#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\
 Data File : BG049253.D
 Acq On : 10 Jul 2021 1:31
 Operator : CG/JU
 Sample : M2878-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGE08

Manual Integrations
 APPROVED

mohammad
 7/12/2021 12:33:52 PM

Quant Time: Jul 10 02:15:36 2021
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	26502	20.00	ng/ul	0.00
20) Naphthalene-d8	10.89	136	108383	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.72	164	81335	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.47	188	192892	20.00	ng/ul	0.00
79) Chrysene-d12	21.75	240	208608	20.00	ng/ul	0.00
88) Perylene-d12	25.05	264	212643	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	1267	2.25	ng/ul	0.00
4) Pyridine-d5	3.90	84	7682m	4.64	ng/ul	0.01
7) Phenol-d5	7.23	99	26463	11.48	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.39	67	16928	12.66	ng/ul	0.00
11) 2-Chlorophenol-d4	7.60	132	21705	12.74	ng/ul	0.00
15) 4-Methylphenol-d8	8.78	113	22504	12.18	ng/ul	0.00
21) Nitrobenzene-d5	9.24	128	11499	13.83	ng/ul	0.00
24) 2-Nitrophenol-d4	9.96	143	12797	13.40	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.51	165	22239	11.94	ng/ul	0.00
31) 4-Chloroaniline-d4	11.03	131	28296	11.66	ng/ul	0.00
46) Dimethylphthalate-d6	14.11	166	85318	13.64	ng/ul	0.00
49) Acenaphthylene-d8	14.41	160	97439	13.28	ng/ul	0.00
54) 4-Nitrophenol-d4	14.91	143	10258	9.91	ng/ul	0.00
60) Fluorene-d10	15.71	176	70184	13.25	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.82	200	12737	10.39	ng/ul	0.00
73) Anthracene-d10	17.57	188	130912	14.91	ng/ul	0.00
81) Pyrene-d10	19.85	212	153096	14.32	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.82	264	140243	12.69	ng/ul	0.00

Target Compounds

Ovalue

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