

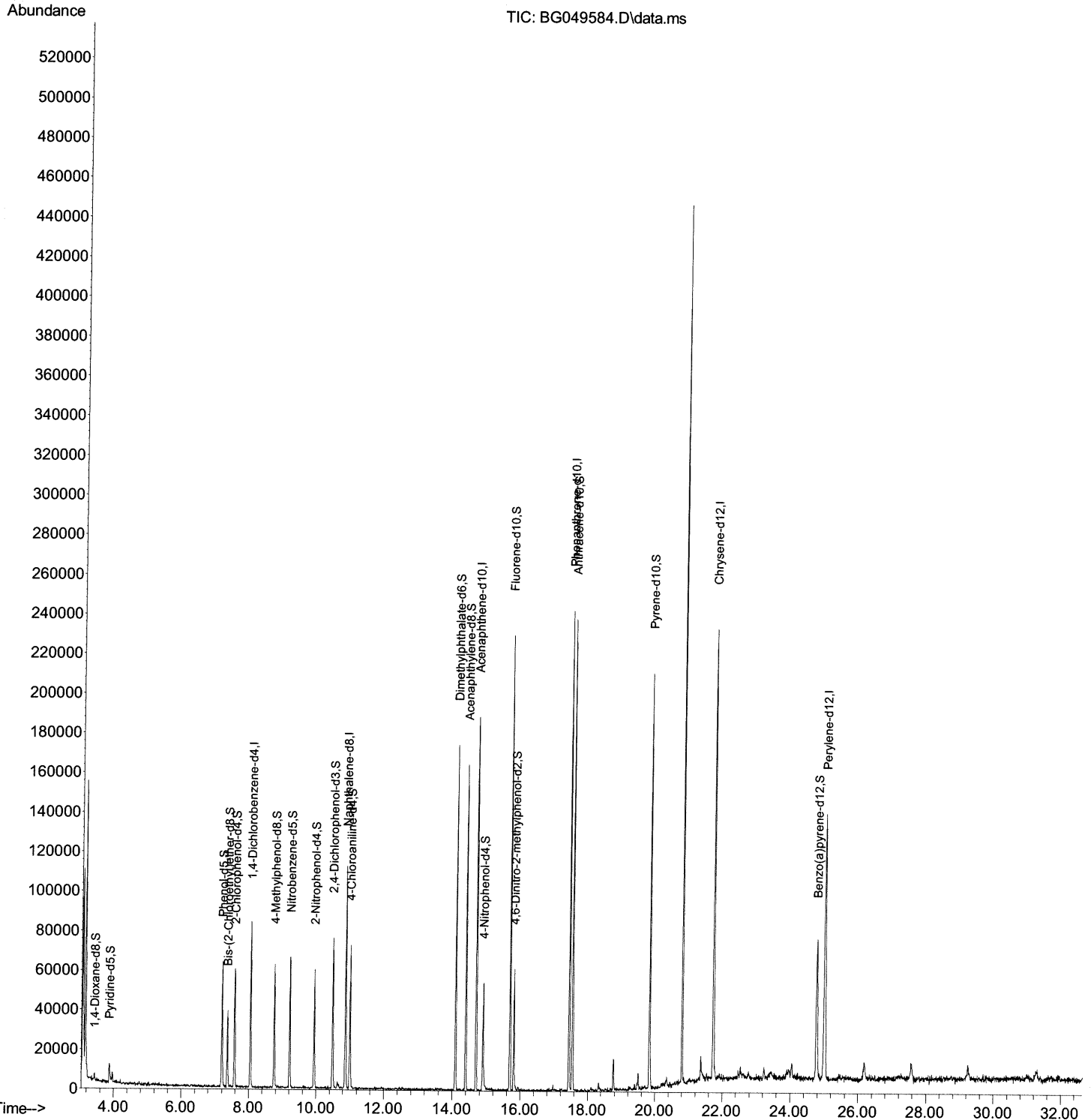
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG073021\
Data File : BG049584.D
Acq On : 30 Jul 2021 19:05
Operator : CG/JU
Sample : M3122-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGEP3

Manual Integrations
APPROVED

mohammad
8/2/2021 3:45:53 PM

Quant Time: Jul 31 02:13:50 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



Quantitation Report (Qedit)

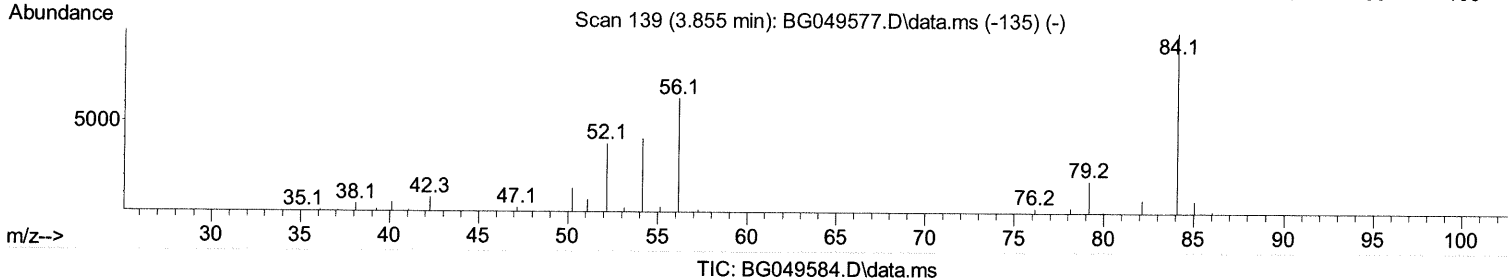
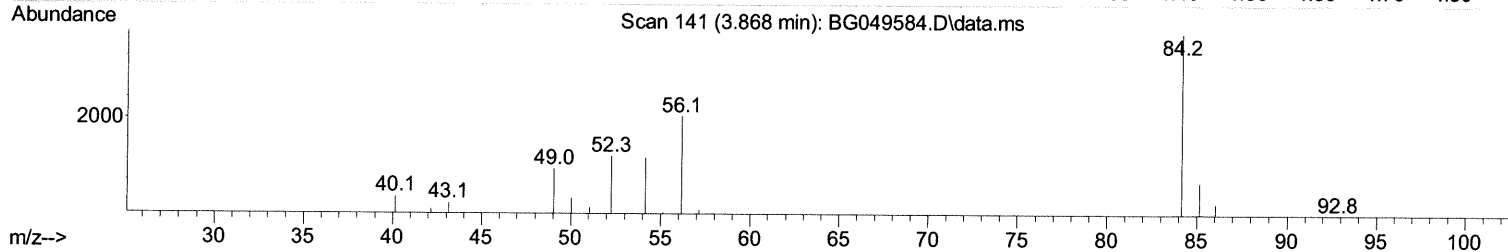
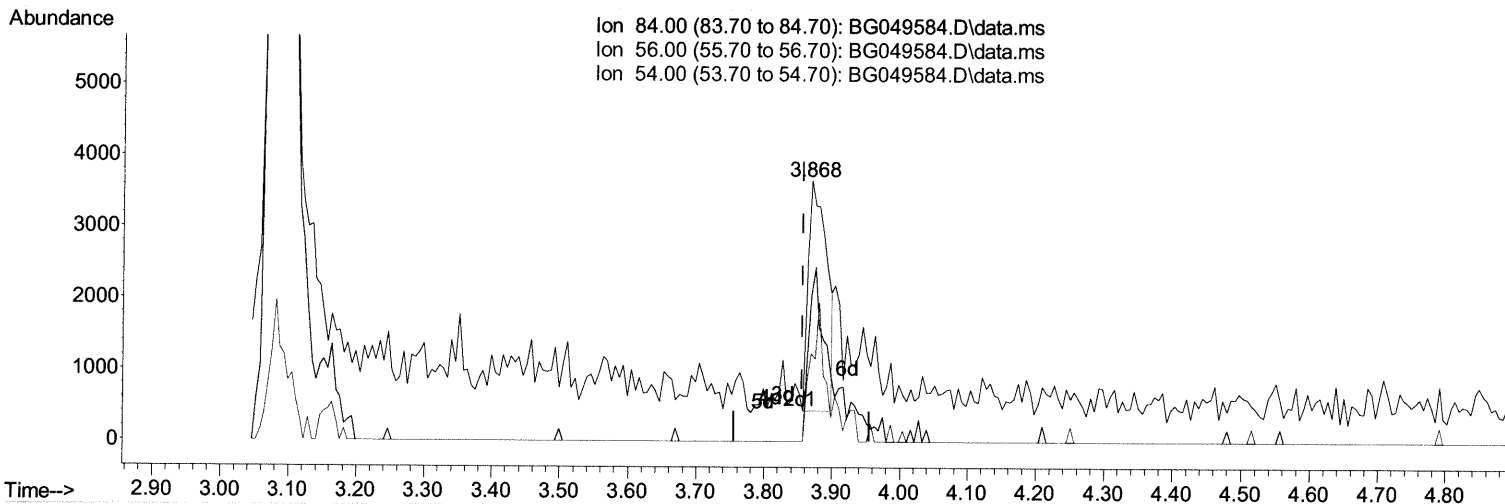
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(4) Pyridine-d5 (S)

3.868min (+ 0.013) 4.42 ng/ul

response 6088

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	55.76
54.00	34.10	33.70
0.00	0.00	0.00

Quantitation Report (Qedit)

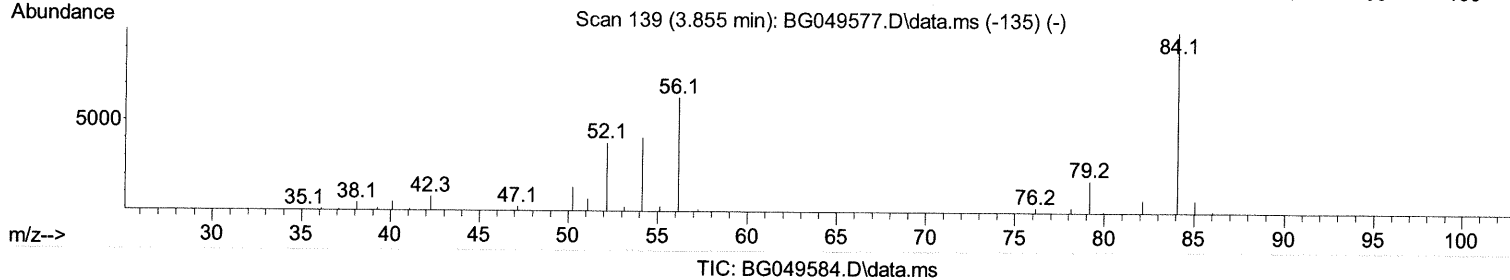
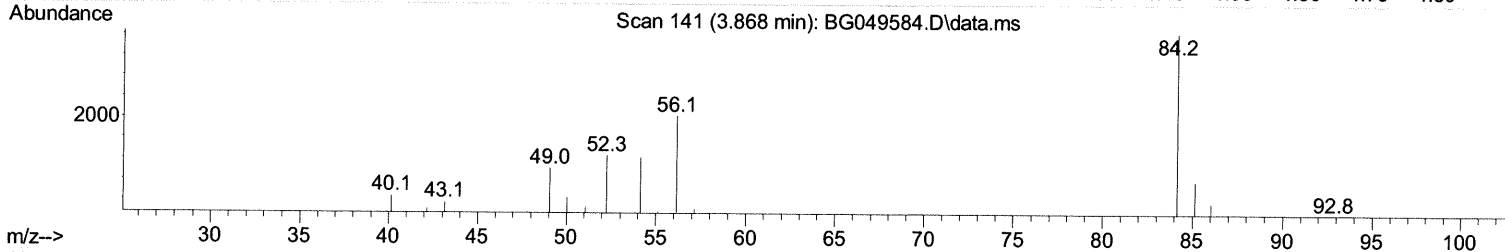
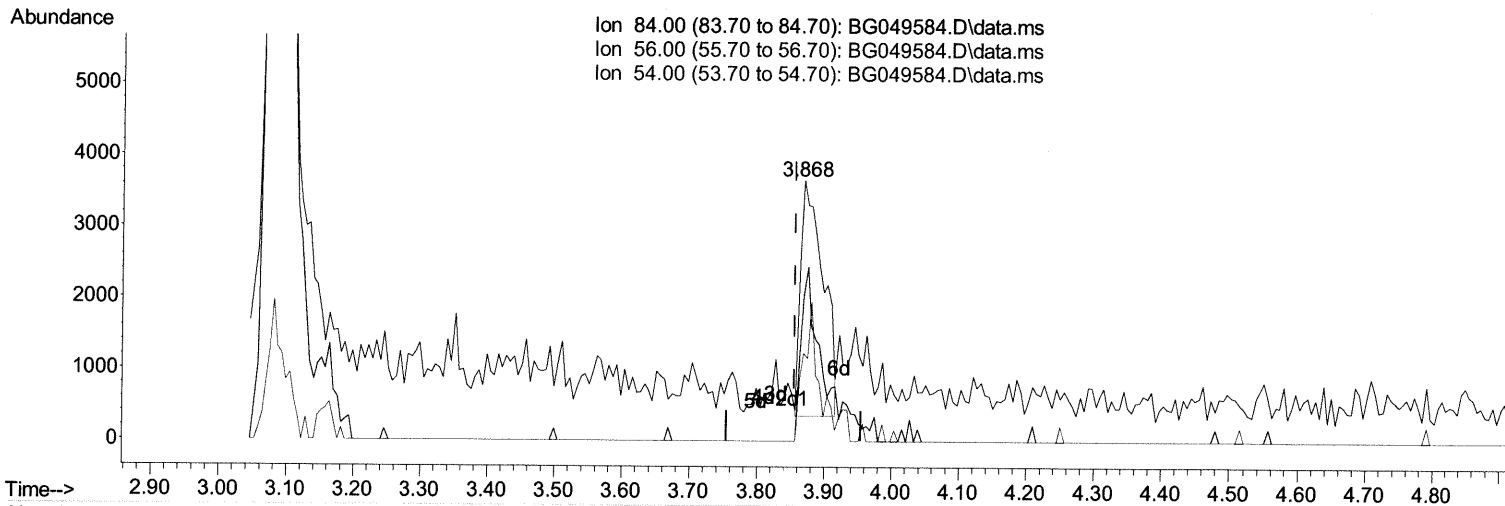
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(4) Pyridine-d5 (S)

3.868min (+ 0.013) 5.58 ng/ul m 08/03/21 JU

response 7678

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	55.76
54.00	34.10	33.70
0.00	0.00	0.00

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Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	21918	20.000	ng/ul	0.00
20) Naphthalene-d8	10.871	136	91970	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.695	164	67115	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.450	188	146296	20.000	ng/ul	0.00
79) Chrysene-d12	21.727	240	141572	20.000	ng/ul	0.00
88) Perylene-d12	25.010	264	146163	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	1581	3.300	ng/uL	0.00
4) Pyridine-d5	3.868	84	7678m	5.580	ng/ul	> 0.01
7) Phenol-d5	7.199	99	33134	17.086	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.364	67	20263	18.450	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	24017	17.278	ng/ul	0.00
15) 4-Methylphenol-d8	8.756	113	23227	15.890	ng/ul	0.00
21) Nitrobenzene-d5	9.214	128	13699	19.374	ng/ul	0.00
24) 2-Nitrophenol-d4	9.943	143	16618	20.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.483	165	26256	17.423	ng/ul	0.00
31) 4-Chloroaniline-d4	11.000	131	34453	16.404	ng/ul	0.00
46) Dimethylphthalate-d6	14.096	166	102663	19.969	ng/ul	0.00
49) Acenaphthylene-d8	14.390	160	115292	19.308	ng/ul	0.00
54) 4-Nitrophenol-d4	14.895	143	13067	15.434	ng/ul	0.00
60) Fluorene-d10	15.688	176	89507	20.209	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.805	200	12995	14.335	ng/ul	0.00
73) Anthracene-d10	17.544	188	141777	20.851	ng/ul	0.00
81) Pyrene-d10	19.835	212	120783	16.685	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.775	264	79904	10.401	ng/ul	0.00

Target Compounds Qvalue

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