

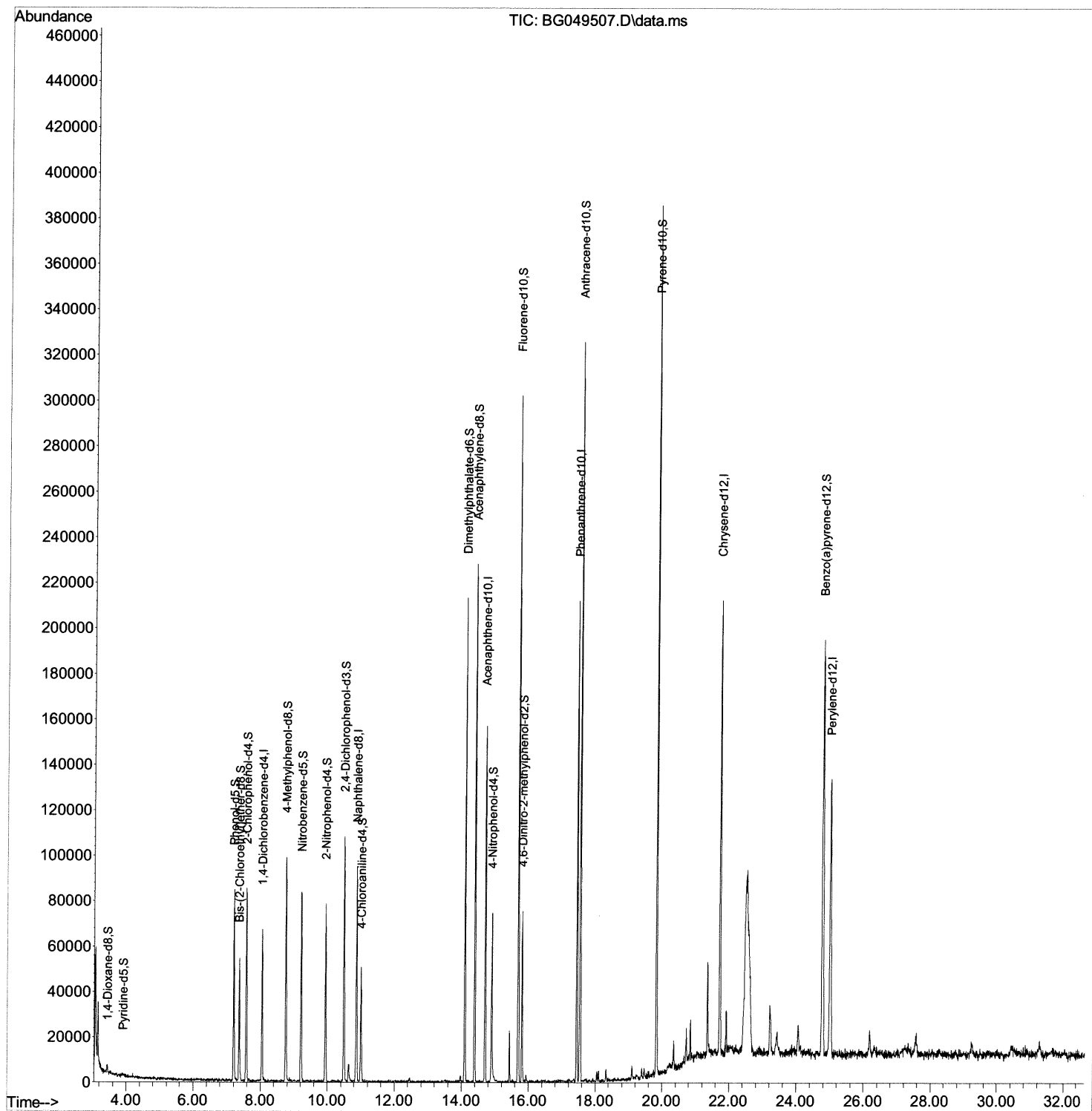
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049507.D
Acq On : 27 Jul 2021 14:30
Operator : CG/JU
Sample : M3084-16
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0034

Manual Integrations
APPROVED

mohammad
7/28/2021 11:29:39 AM

Quant Time: Jul 27 15:03:56 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)

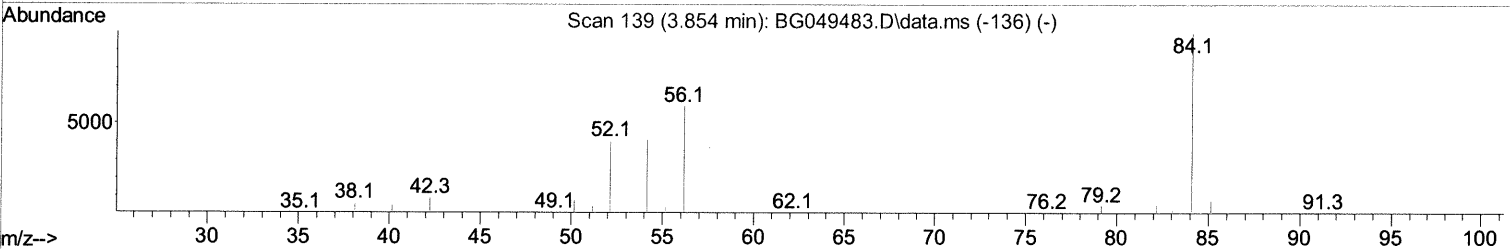
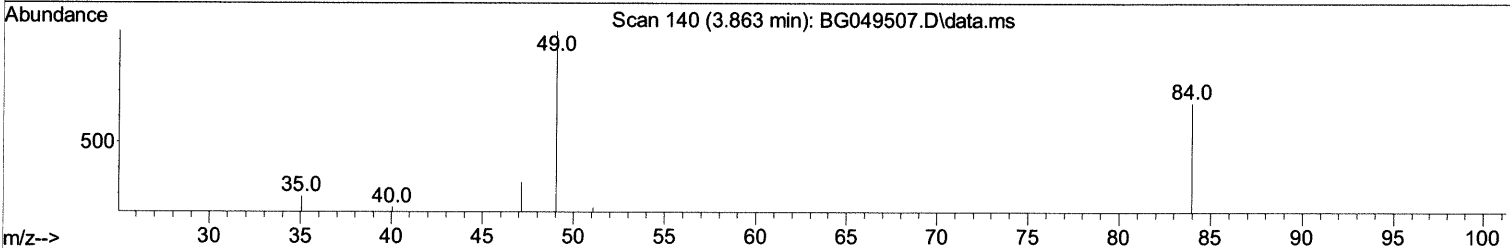
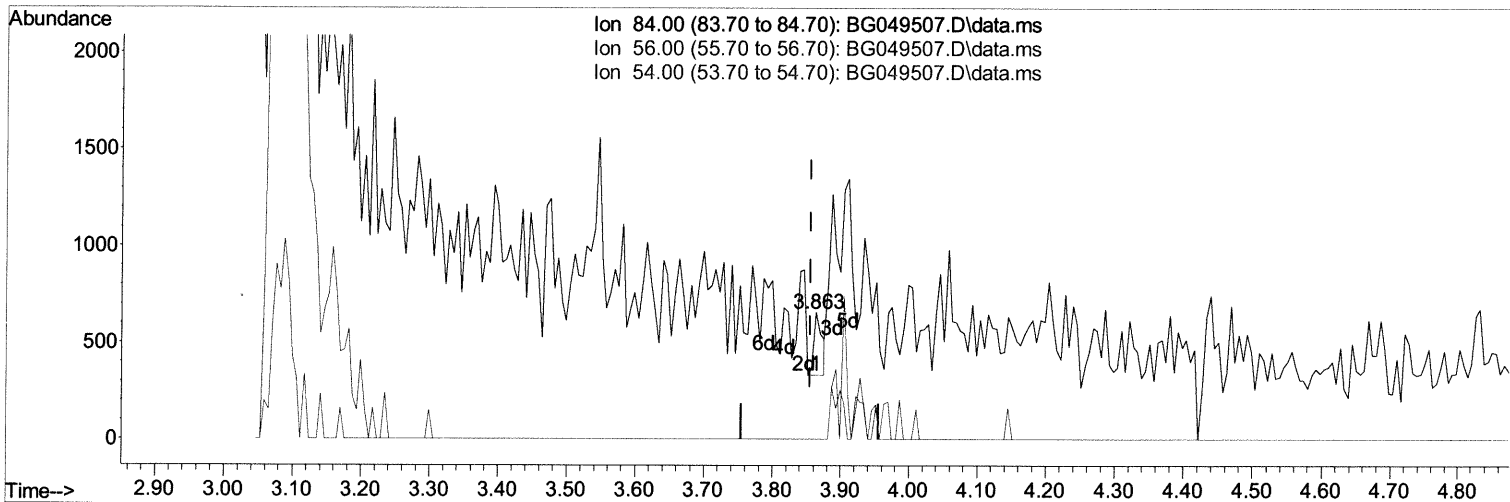
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TIC: BG049507.D\data.ms

(4) Pyridine-d5 (S)

3.863min (+ 0.008) 0.24 ng/ul

response 266

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	0.00#
54.00	34.10	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

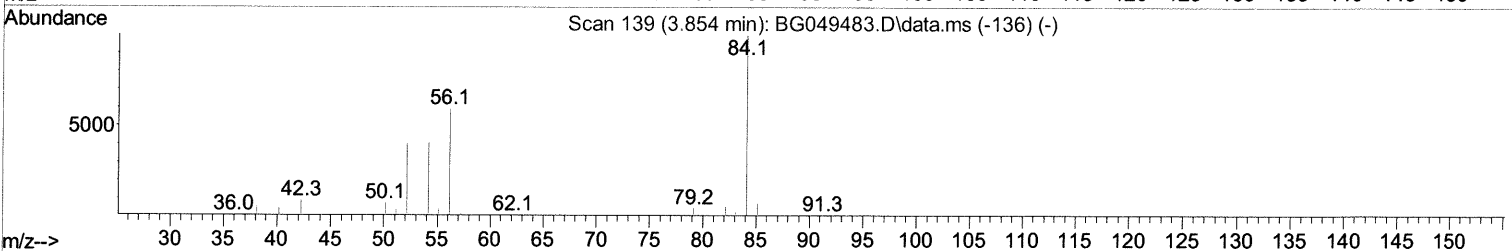
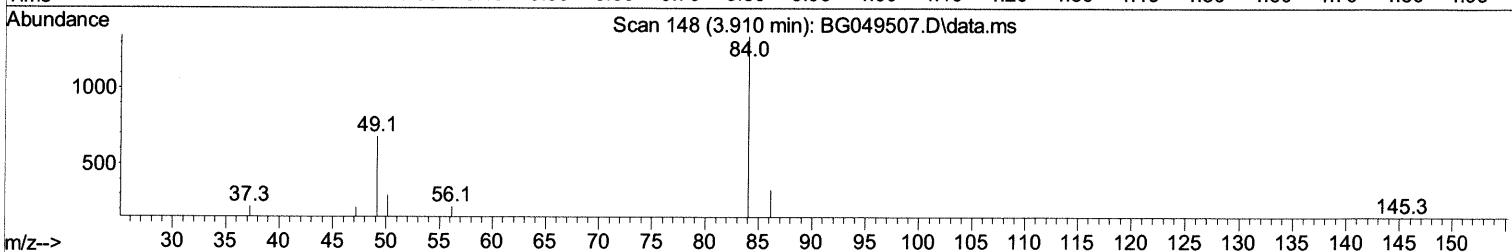
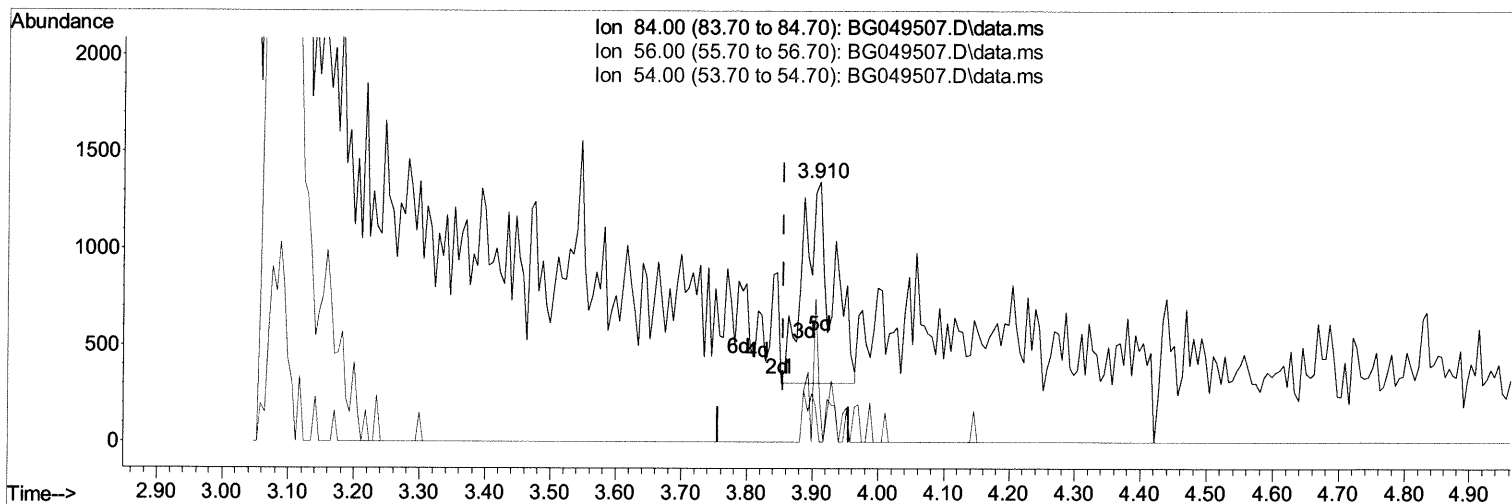
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TIC: BG049507.D\data.ms

(4) Pyridine-d5 (S)

3.910min (+ 0.055) 2.85 ng/ul m 07/29/21 Ju

response 3205

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	16.36#
54.00	34.10	0.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
 Data File : BG049507.D
 Acq On : 27 Jul 2021 14:30
 Operator : CG/JU
 Sample : M3084-16
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0034

Manual Integrations
 APPROVED

mohammad
 7/28/2021 11:29:39 AM

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.052	152	17897	20.000	ng/ul	0.00
20) Naphthalene-d8	10.871	136	77047	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.695	164	58419	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188	132201	20.000	ng/ul	0.00
79) Chrysene-d12	21.727	240	120362	20.000	ng/ul	0.00
88) Perylene-d12	25.005	264	126614	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.440	96	1601	4.093	ng/uL	0.00
4) Pyridine-d5	3.910	84	3205m >	2.852	ng/ul >	0.06
7) Phenol-d5	7.206	99	45673	28.844	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.364	67	26856	29.946	ng/ul	0.00
11) 2-Chlorophenol-d4	7.576	132	34649	30.527	ng/ul	0.00
15) 4-Methylphenol-d8	8.756	113	37755	31.633	ng/ul	0.00
21) Nitrobenzene-d5	9.221	128	18223	30.763	ng/ul	0.00
24) 2-Nitrophenol-d4	9.943	143	20744	29.884	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.489	165	37641	29.815	ng/ul	0.00
31) 4-Chloroaniline-d4	11.006	131	25856	14.695	ng/ul	0.00
46) Dimethylphthalate-d6	14.096	166	130717	29.211	ng/ul	0.00
49) Acenaphthylene-d8	14.390	160	165849	31.909	ng/ul	0.00
54) 4-Nitrophenol-d4	14.895	143	16558	22.469	ng/ul	0.00
60) Fluorene-d10	15.688	176	123268	31.974	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.806	200	15676	19.136	ng/ul	0.00
73) Anthracene-d10	17.545	188	195417	31.805	ng/ul	0.00
81) Pyrene-d10	19.830	212	219315	35.635	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.776	264	198679	29.855	ng/ul	0.00

Target Compounds Qvalue

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