

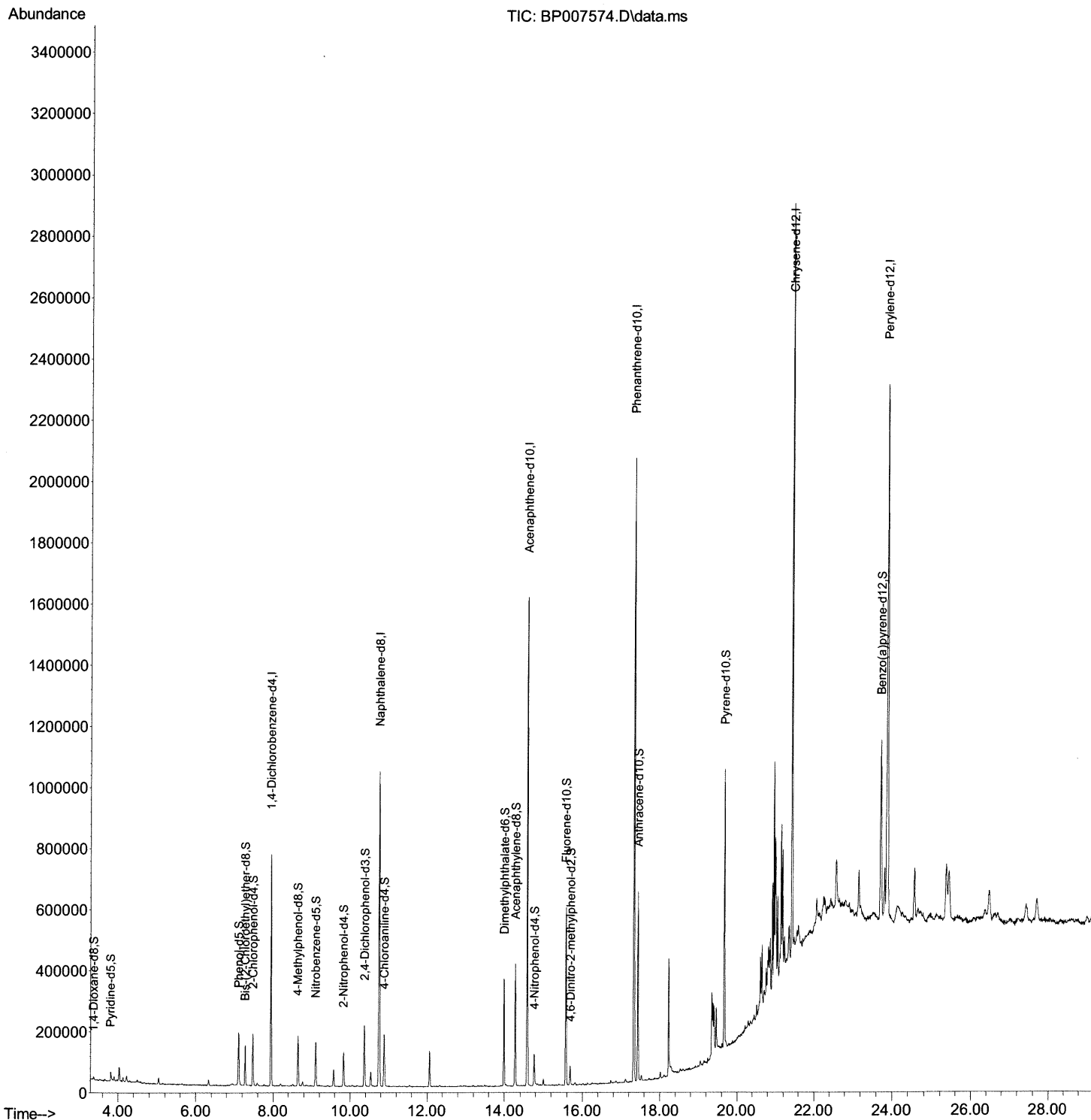
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007574.D
Acq On : 22 Oct 2021 04:16
Operator : CG/JU
Sample : M4217-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGC1

Manual Integrations
APPROVED

mohammad
10/22/2021 2:55:02 PM

Quant Time: Oct 22 04:47:44 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 20 10:49:33 2021
Response via : Initial Calibration



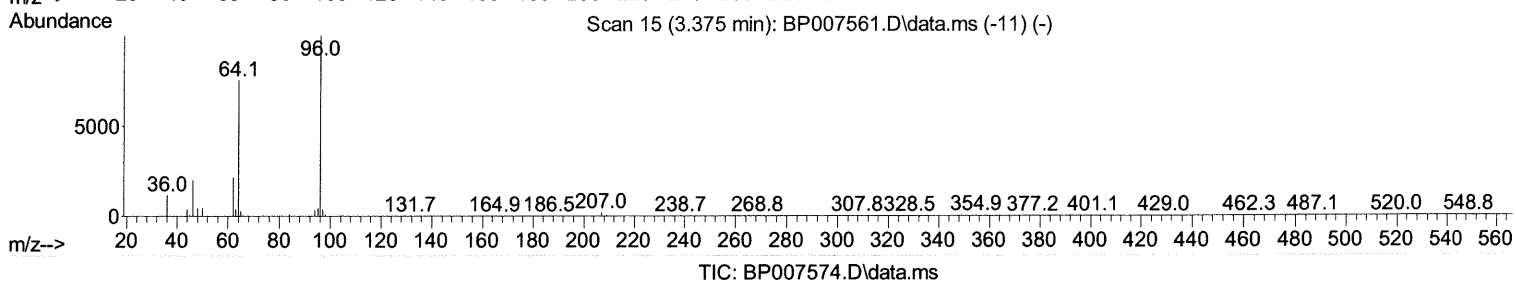
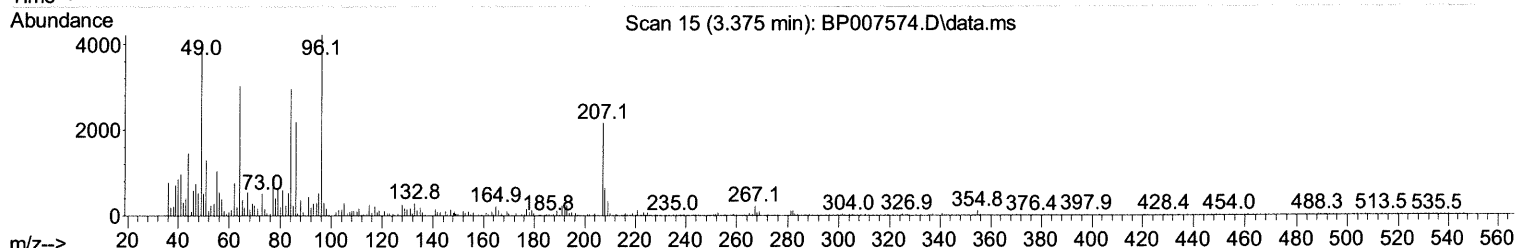
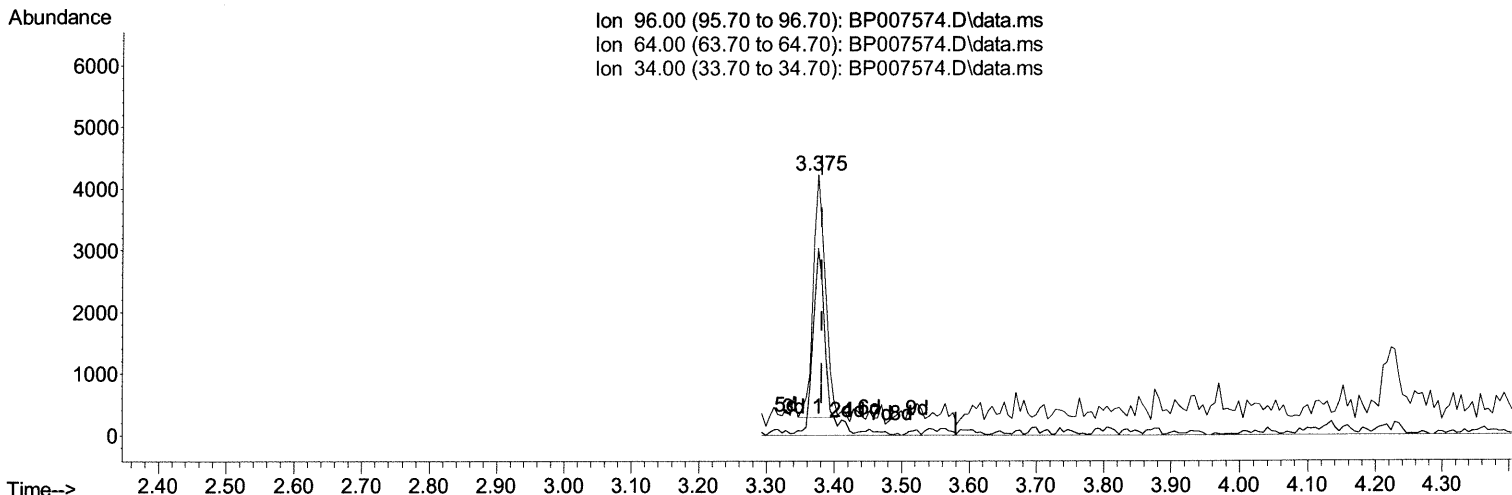
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(3) 1,4-Dioxane-d8 (S)

3.375min (-0.006) 0.93 ng/uL

response 4983

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	71.71
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

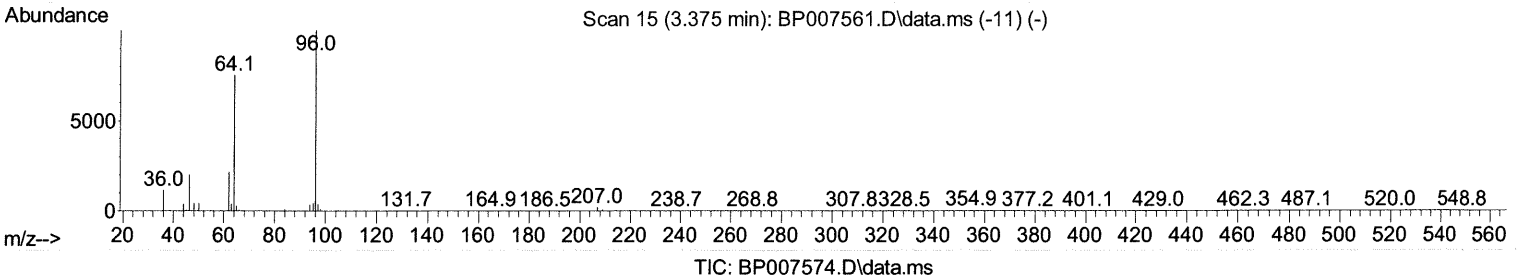
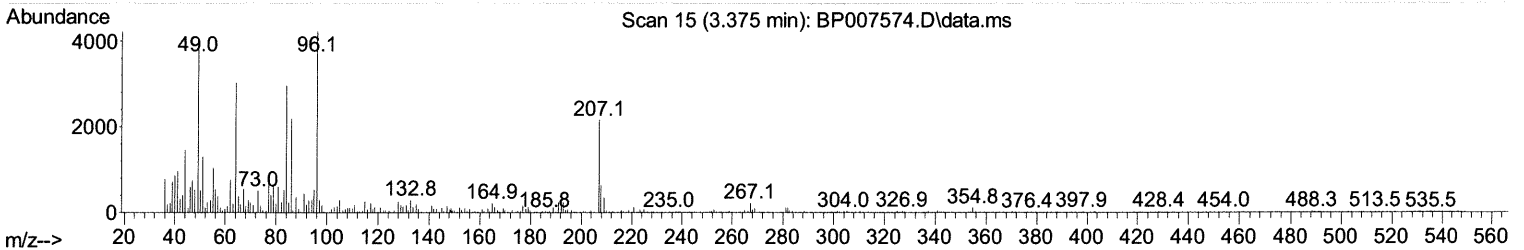
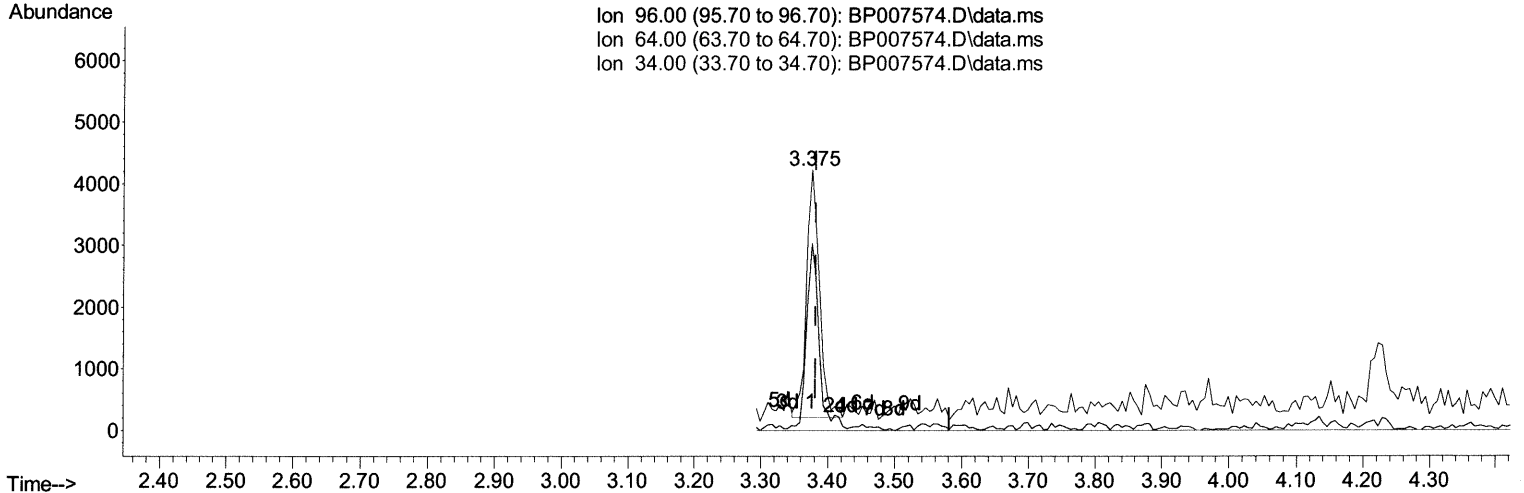
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TIC: BP007574.D\data.ms

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

3.375min (-0.006) 1.01 ng/uL m 10/26/21 JU

response 5385

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	71.71
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

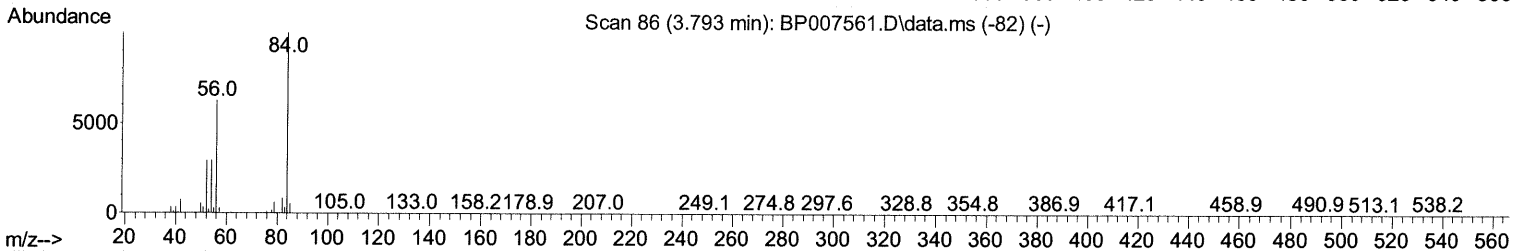
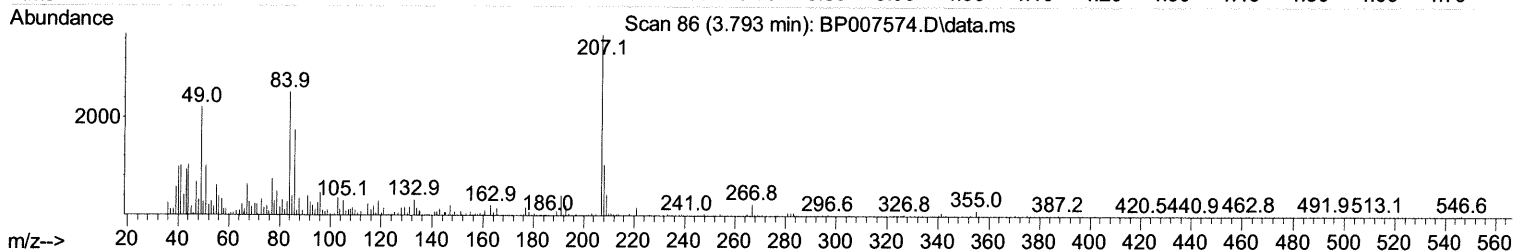
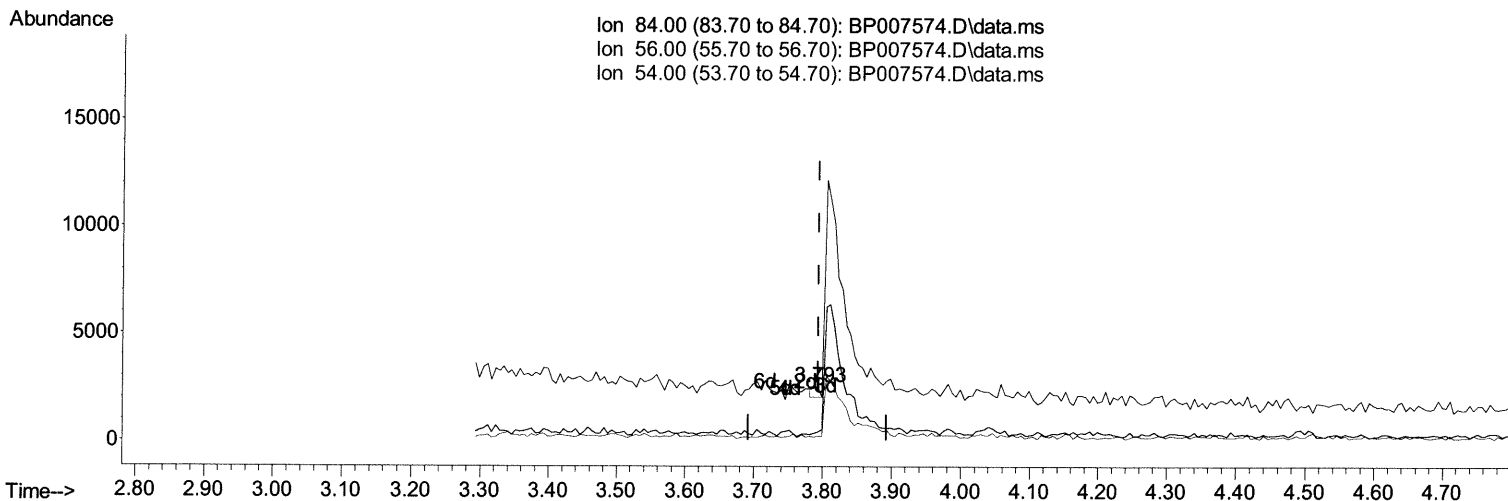
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(4) Pyridine-d5 (S)

3.793min (+ 0.000) 0.02 ng/ul

response 317

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	15.56#
54.00	31.70	7.06#
0.00	0.00	0.00

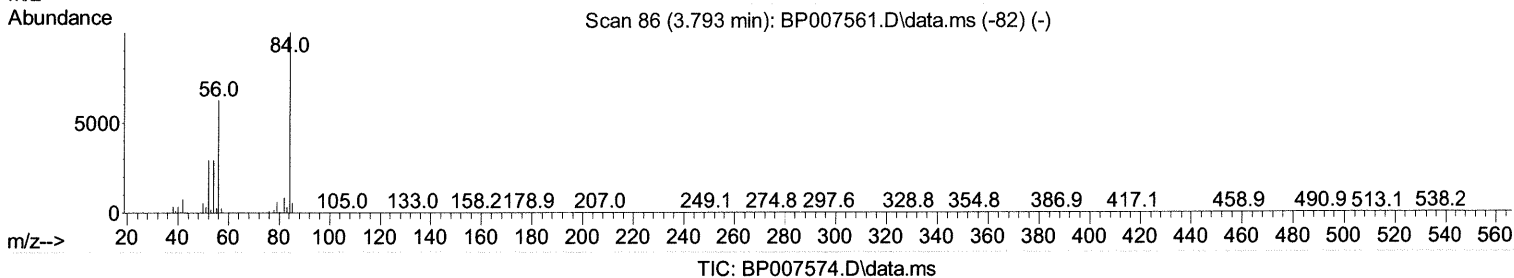
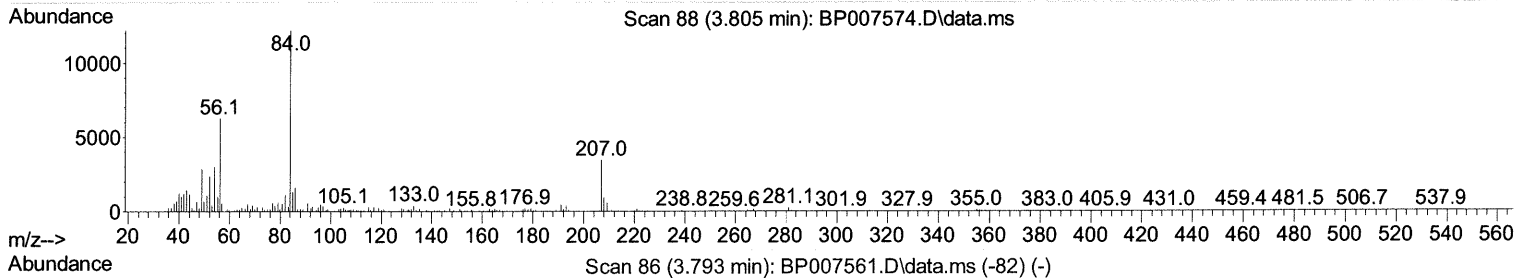
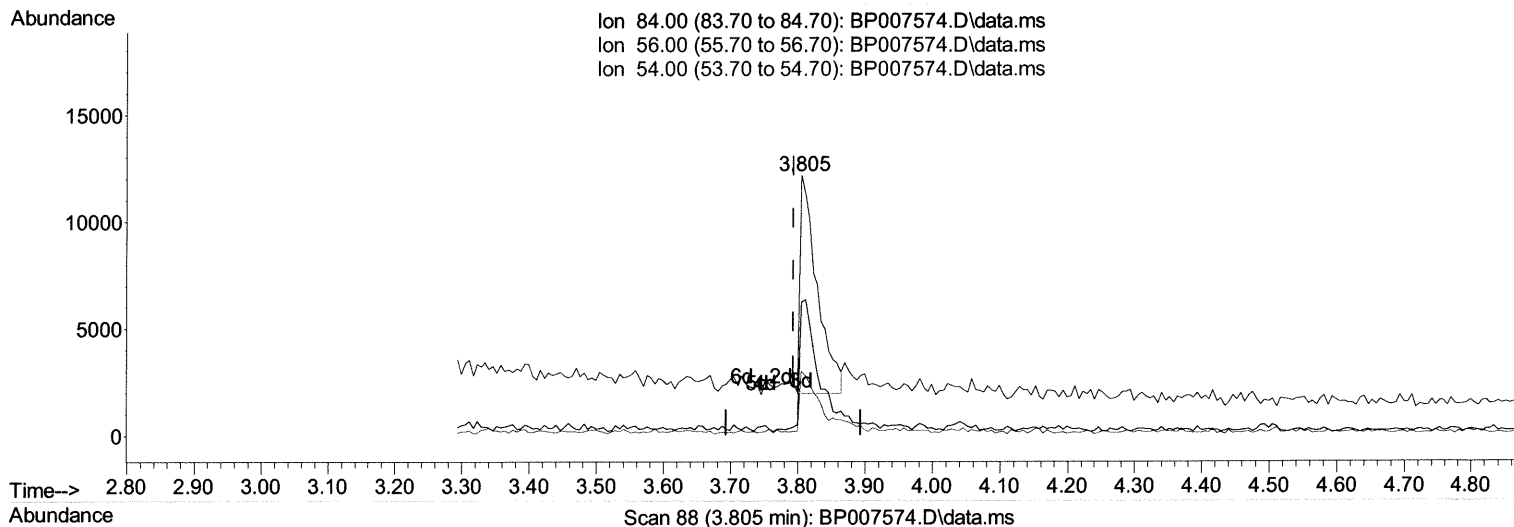
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Instrument :
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(4) Pyridine-d5 (S)

3.805min (+ 0.012) 1.23 ng/ul m 10/26/2020

response 17968

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	51.38#
54.00	31.70	24.67#
0.00	0.00	0.00

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 ALS Vial : 34 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	210682	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	848343	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	555502	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	1209291	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	1279263	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1342232	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	5385m	1.010	ng/ul	0.00
4) Pyridine-d5	3.805	84	17968m	1.231	ng/ul	0.01
7) Phenol-d5	7.105	99	96422	5.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	56201	5.727	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	76881	5.940	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	61561	4.857	ng/ul	-0.01
21) Nitrobenzene-d5	9.099	128	35342	6.583	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.828	143	36028	7.184	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	73147	5.980	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	91442	5.528	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	242596	6.312	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	282992	5.859	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	26081	4.229	ng/ul	-0.01
60) Fluorene-d10	15.575	176	220263	6.849	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	13835	2.855	ng/ul	0.00
73) Anthracene-d10	17.433	188	357777	6.967	ng/ul	0.00
81) Pyrene-d10	19.663	212	466259	7.291	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	448642	6.718	ng/ul	-0.01

Target Compounds Qvalue

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