

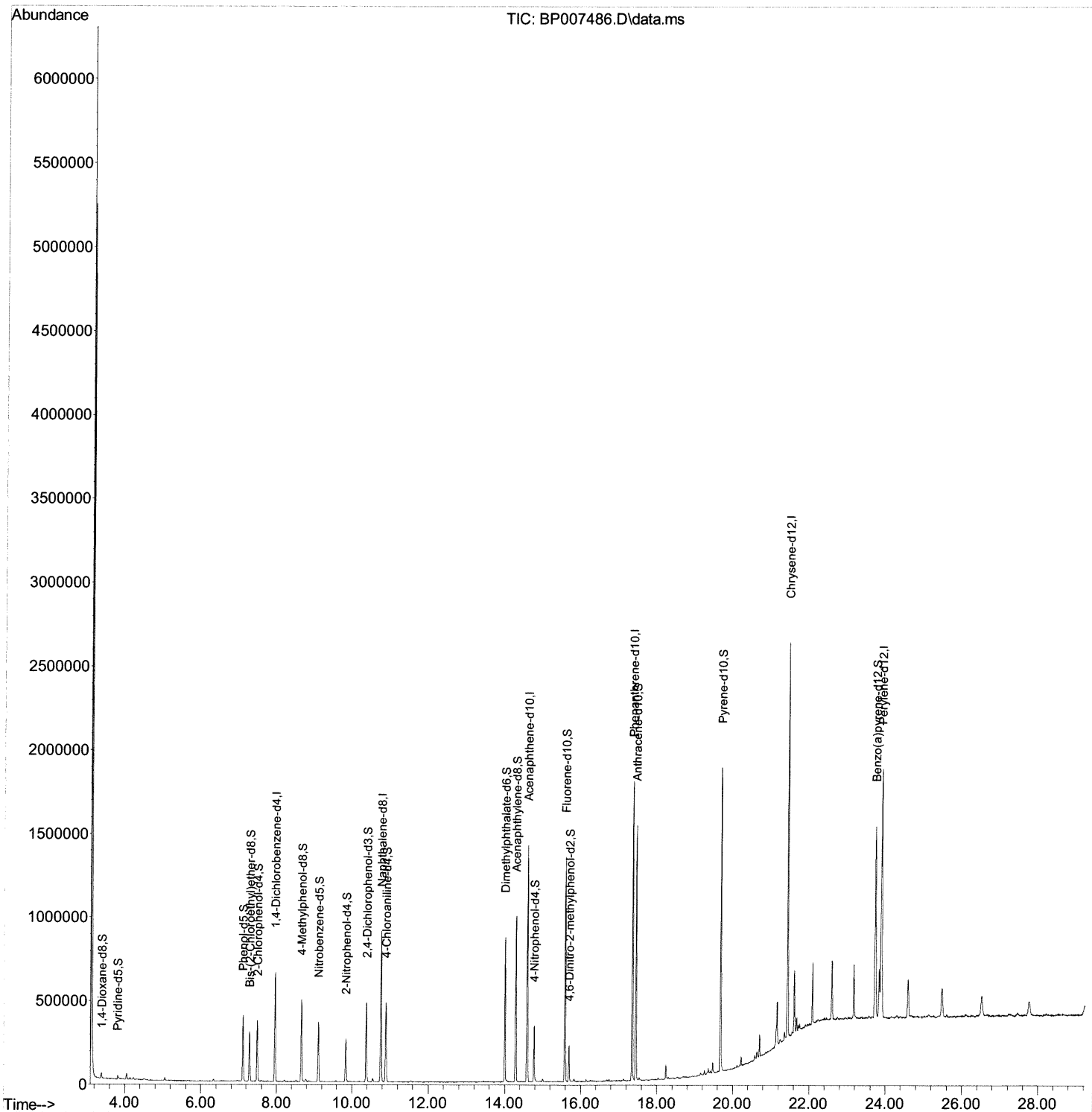
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007486.D
Acq On : 19 Oct 2021 10:15
Operator : CG/JU
Sample : M4196-18
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGD6

Manual Integrations
APPROVED

mohammad
10/19/2021 12:37:04 PM

Quant Time: Oct 19 10:51:13 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 18 11:26:58 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

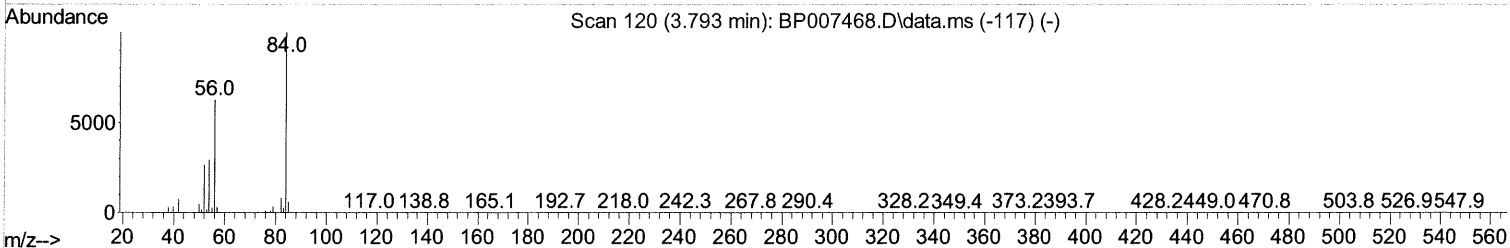
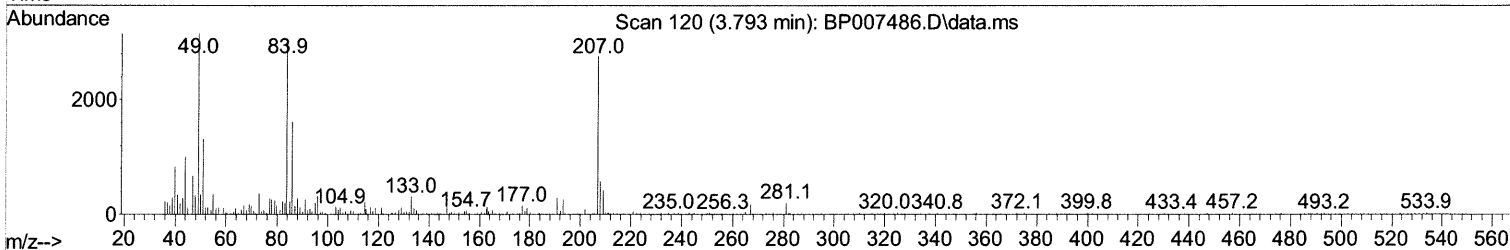
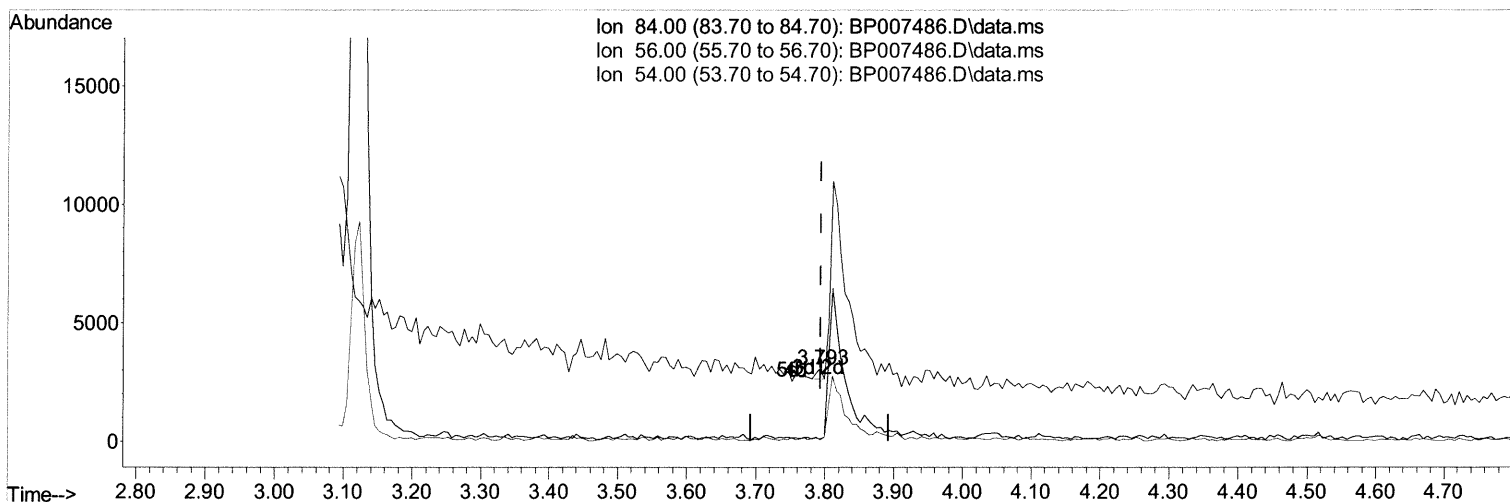
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TIC: BP007486.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.000) 0.02 ng/ul

response 247

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	3.42#
54.00	31.70	2.47#
0.00	0.00	0.00

Quantitation Report (Qedit)

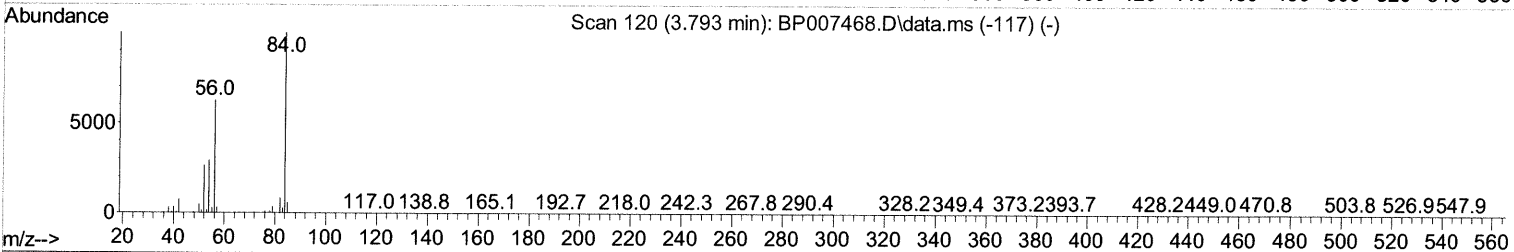
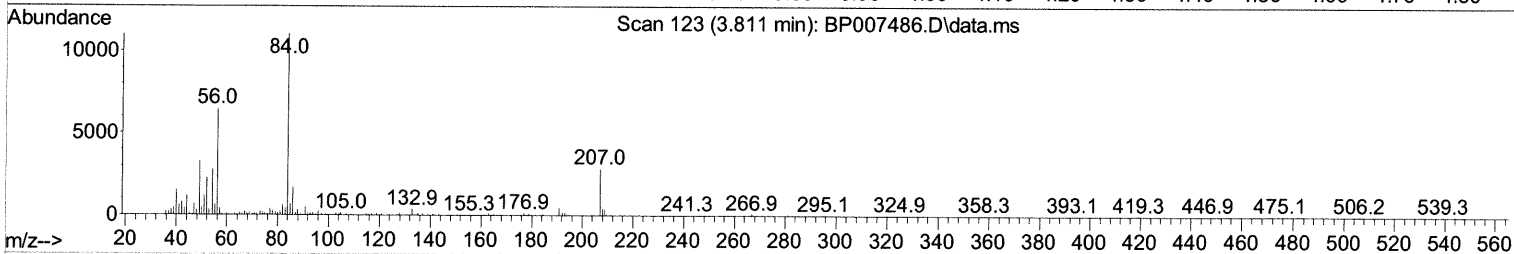
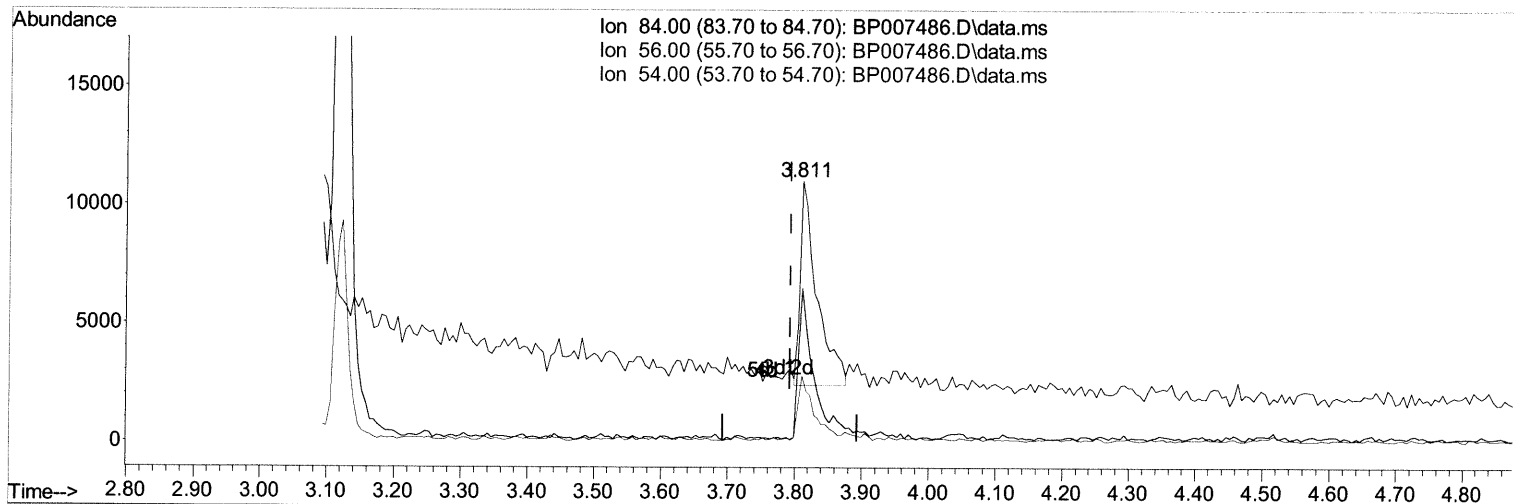
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TIC: BP007486.D\data.ms

(4) Pyridine-d5 (S)

3.811min (+ 0.018) 1.20 ng/ul m 10/20/21 JU

response 14847

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	58.73
54.00	31.70	25.08#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007486.D
 Acq On : 19 Oct 2021 10:15
 Operator : CG/JU
 Sample : M4196-18
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGD6

Manual Integrations
 APPROVED

mohammad
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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.958	152	179056	20.000 ng/ul	0.00
20) Naphthalene-d8	10.763	136	733964	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.593	164	484110	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1061274	20.000 ng/ul	0.00
79) Chrysene-d12	21.433	240	1146167	20.000 ng/ul	0.00
88) Perylene-d12	23.886	264	1176430	20.000 ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.381	96	17610	3.884 ng/uL	0.00
4) Pyridine-d5	3.811	84	14847m >	1.196 ng/ul >	0.02 10/20/21
7) Phenol-d5	7.111	99	227511	16.270 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	129427	15.518 ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	173395	15.762 ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	180123	16.723 ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	83114	17.894 ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	83911	19.341 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	175538	16.588 ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	258080	18.035 ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	582686	17.397 ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	713143	16.943 ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	78771	14.657 ng/ul	0.00
60) Fluorene-d10	15.587	176	518503	18.499 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	48222	11.340 ng/ul	0.00
73) Anthracene-d10	17.445	188	842351	18.690 ng/ul	0.00
81) Pyrene-d10	19.675	212	957853	16.717 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	868082	14.830 ng/ul	0.00

Target Compounds

Qvalue

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