

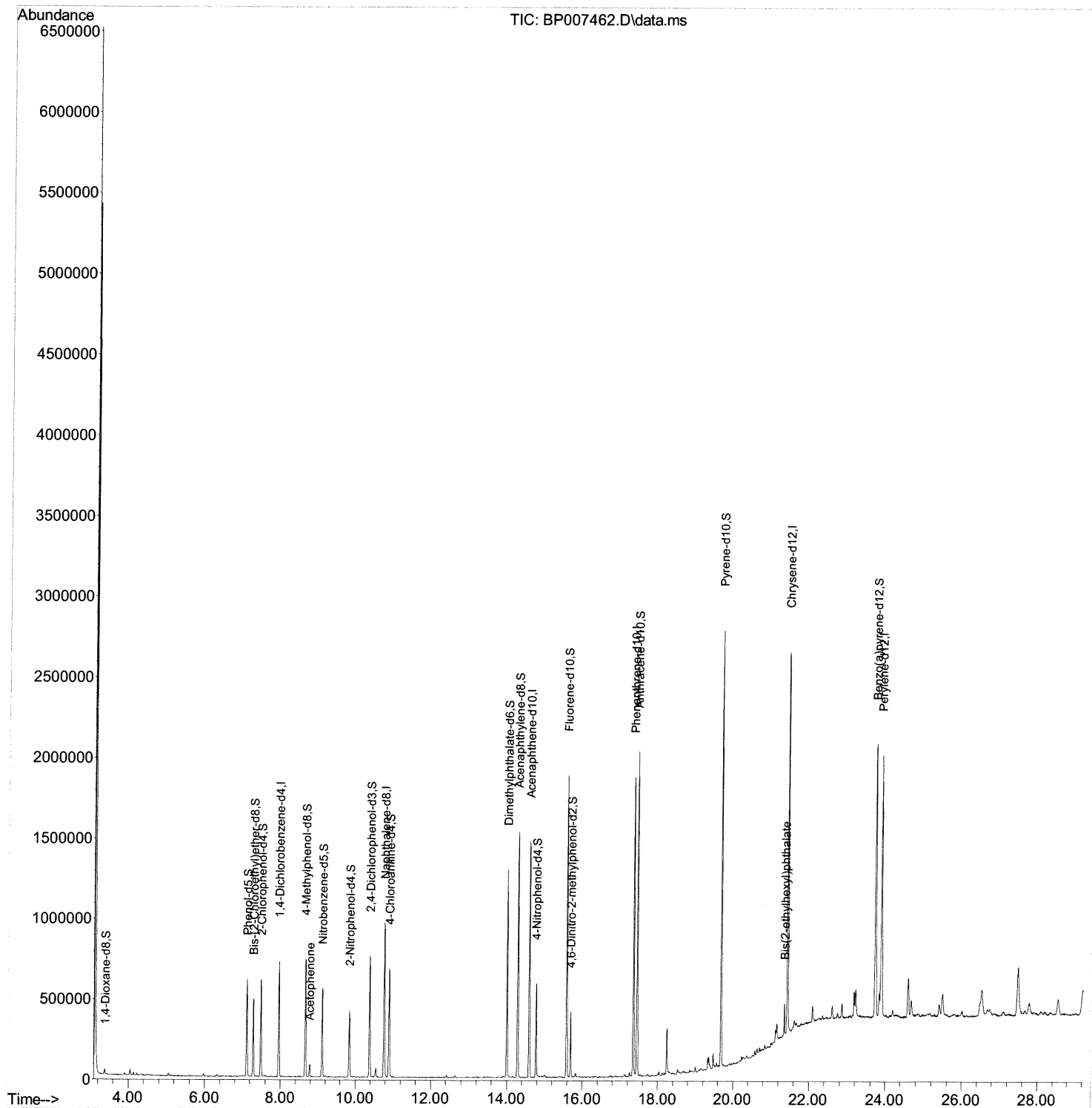
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007462.D
Acq On : 18 Oct 2021 14:06
Operator : CG/JU
Sample : M4181-17
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B15

Manual Integrations
APPROVED

mohammad
10/19/2021 12:36:36 PM

Quant Time: Oct 18 14:32:05 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)

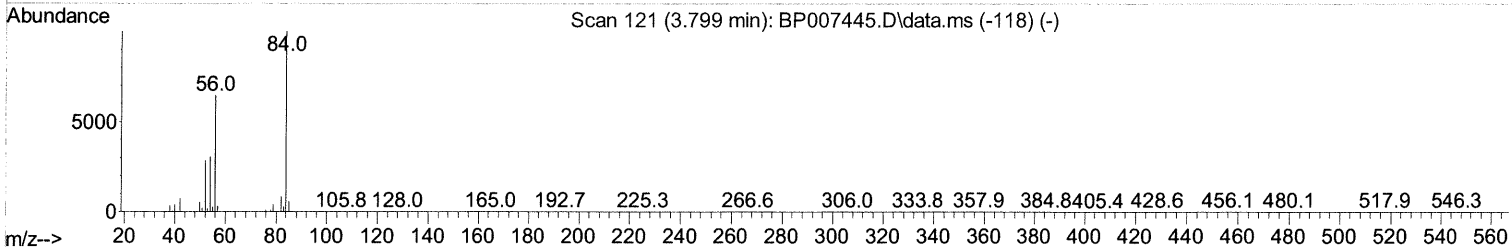
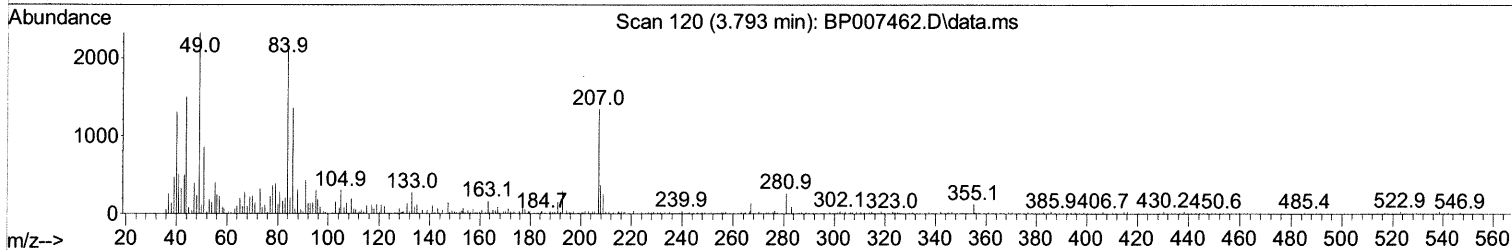
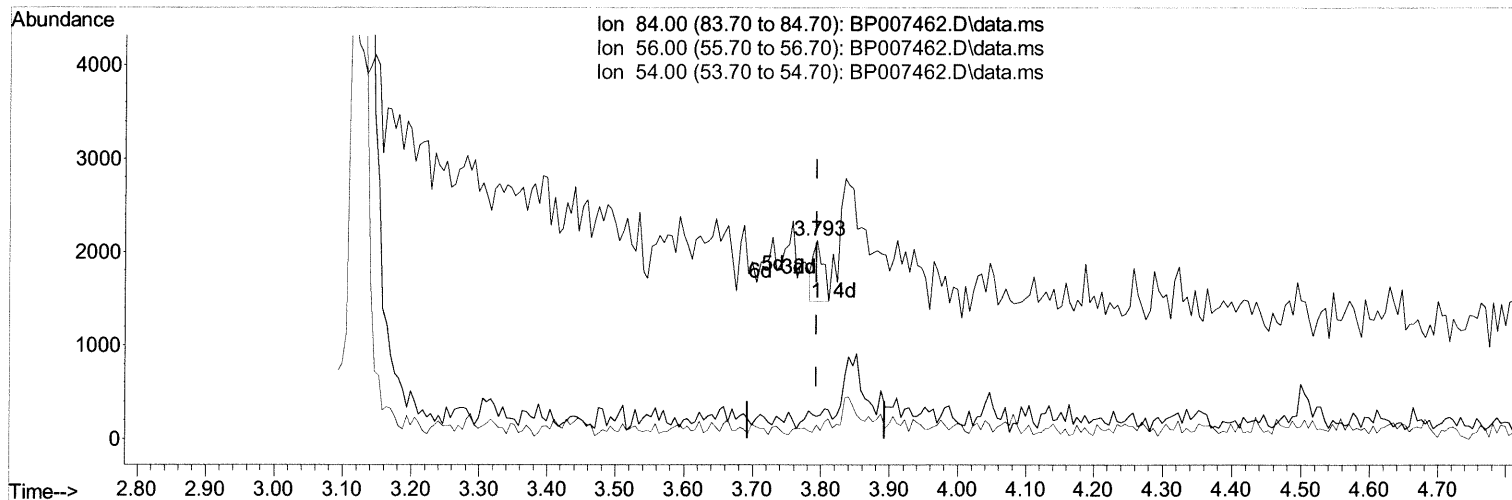
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TIC: BP007462.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.000) 0.05 ng/ul

response 694

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	11.42#
54.00	31.70	7.10#
0.00	0.00	0.00

Quantitation Report (Qedit)

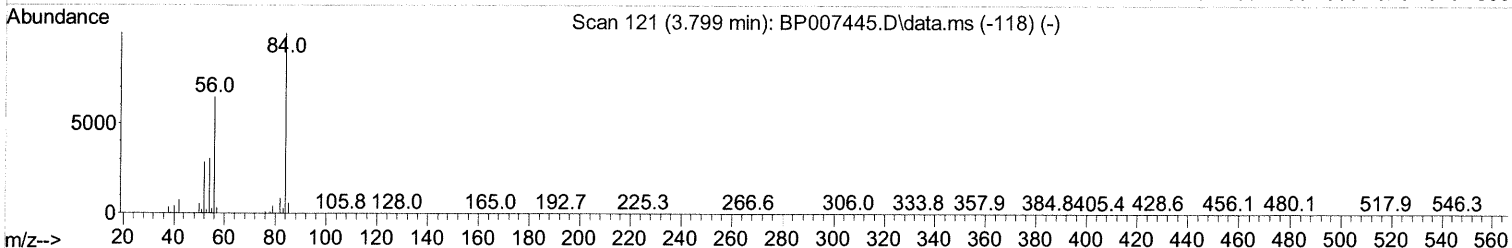
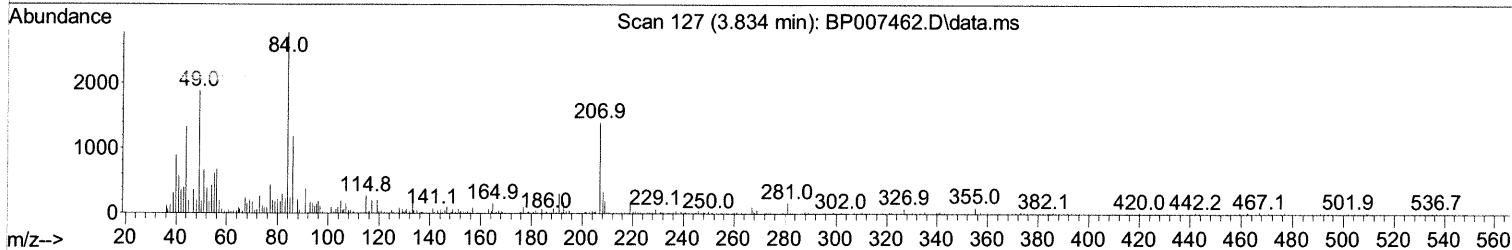
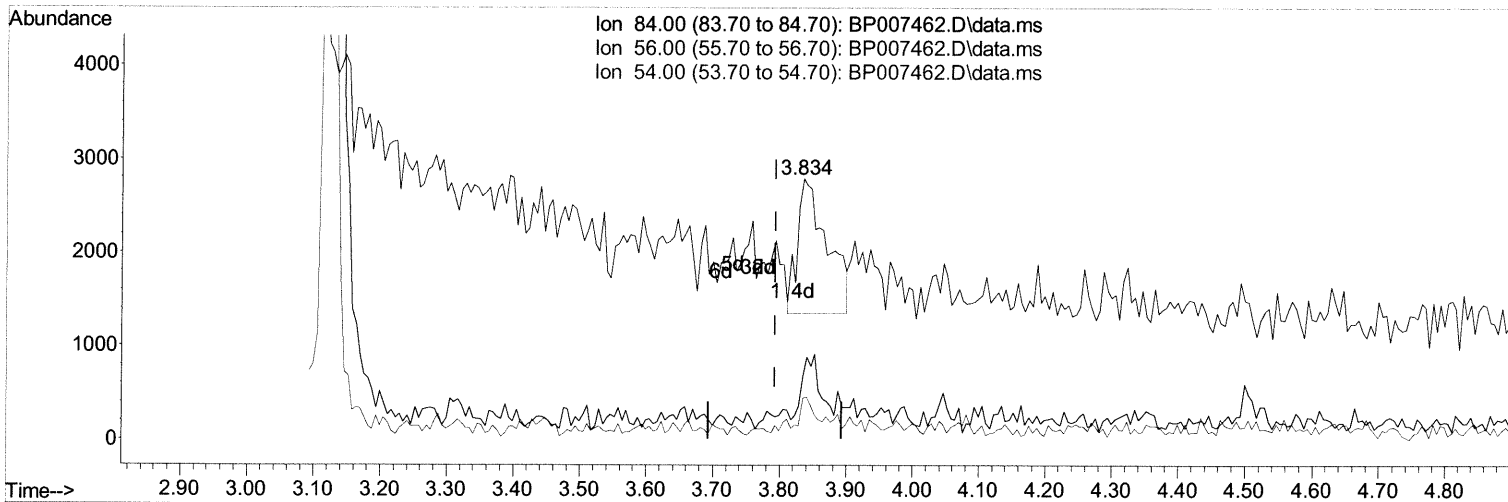
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Manual Integrations
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TIC: BP007462.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.041) 0.33 ng/ul m 10/20/21JU

response 4449

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	24.41#
54.00	31.70	15.65#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007462.D
 Acq On : 18 Oct 2021 14:06
 Operator : CG/JU
 Sample : M4181-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B15

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	195115	20.000 ng/ul	0.00
20) Naphthalene-d8	10.763	136	796291	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.593	164	520086	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1137418	20.000 ng/ul	0.00
79) Chrysene-d12	21.433	240	1212284	20.000 ng/ul	0.00
88) Perylene-d12	23.892	264	1257784	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.381	96	17630	3.569 ng/uL	0.00
4) Pyridine-d5	3.834	84	4449m	0.329 ng/ul	> 0.04
7) Phenol-d5	7.116	99	354674	23.276 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	207998	22.887 ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	276916	23.101 ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	273226	23.279 ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	132051	26.204 ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	130965	27.823 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	280965	24.472 ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	356799	22.982 ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	871981	24.234 ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1032432	22.832 ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	132277	22.911 ng/ul	0.00
60) Fluorene-d10	15.587	176	749399	24.888 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	87476	19.194 ng/ul	0.00
73) Anthracene-d10	17.445	188	1191091	24.658 ng/ul	0.00
81) Pyrene-d10	19.680	212	1343324	22.166 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1386920	22.161 ng/ul	0.00
Target Compounds					
16) Acetophenone	8.781	105	42590	2.372 ng/ul	93
86) Bis(2-ethylhexyl)phtha...	21.357	149	68413	1.607 ng/ul	99

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