

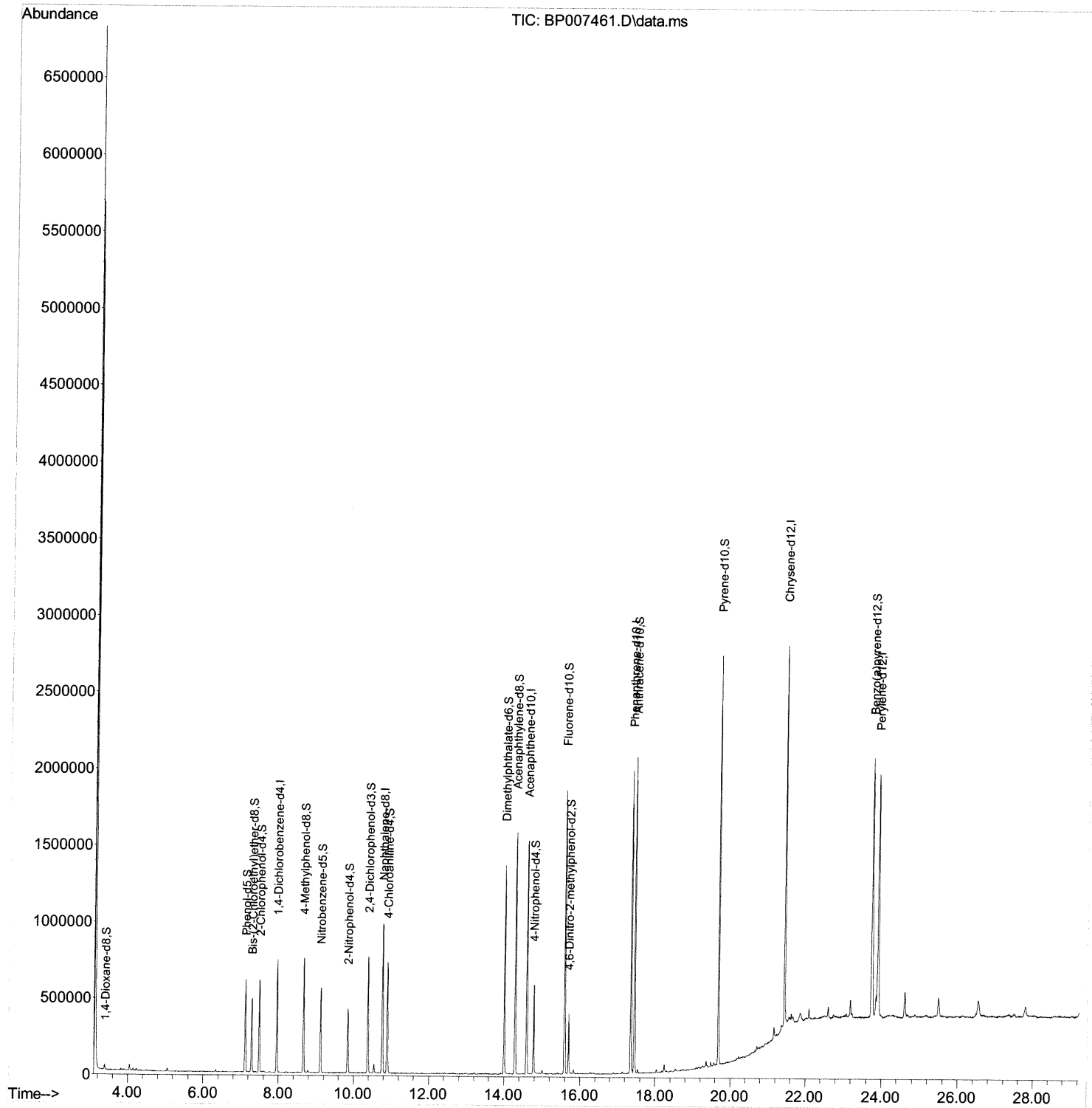
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007461.D
Acq On : 18 Oct 2021 13:32
Operator : CG/JU
Sample : M4181-18
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B16

Manual Integrations
APPROVED

mohammad
10/19/2021 12:36:33 PM

Quant Time: Oct 18 14:19:04 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)

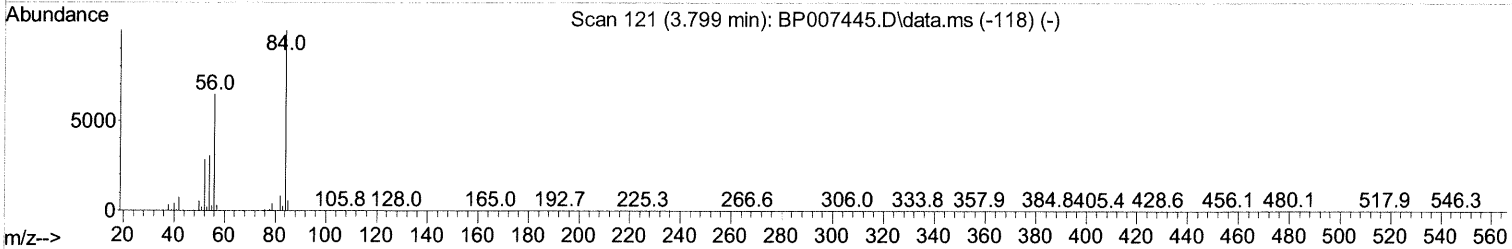
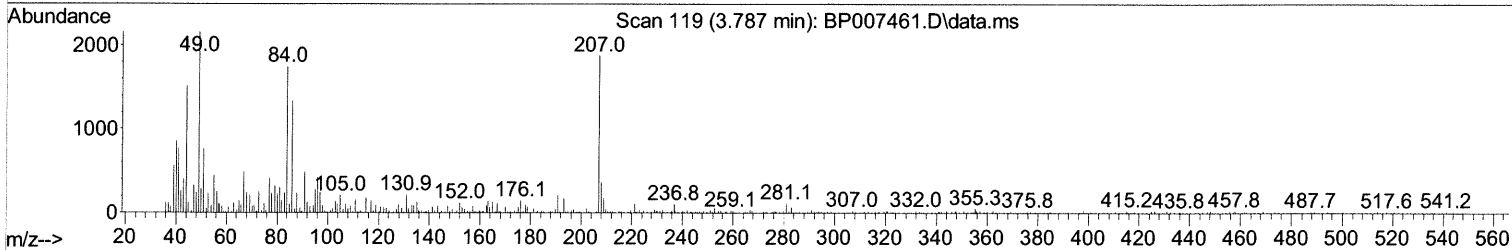
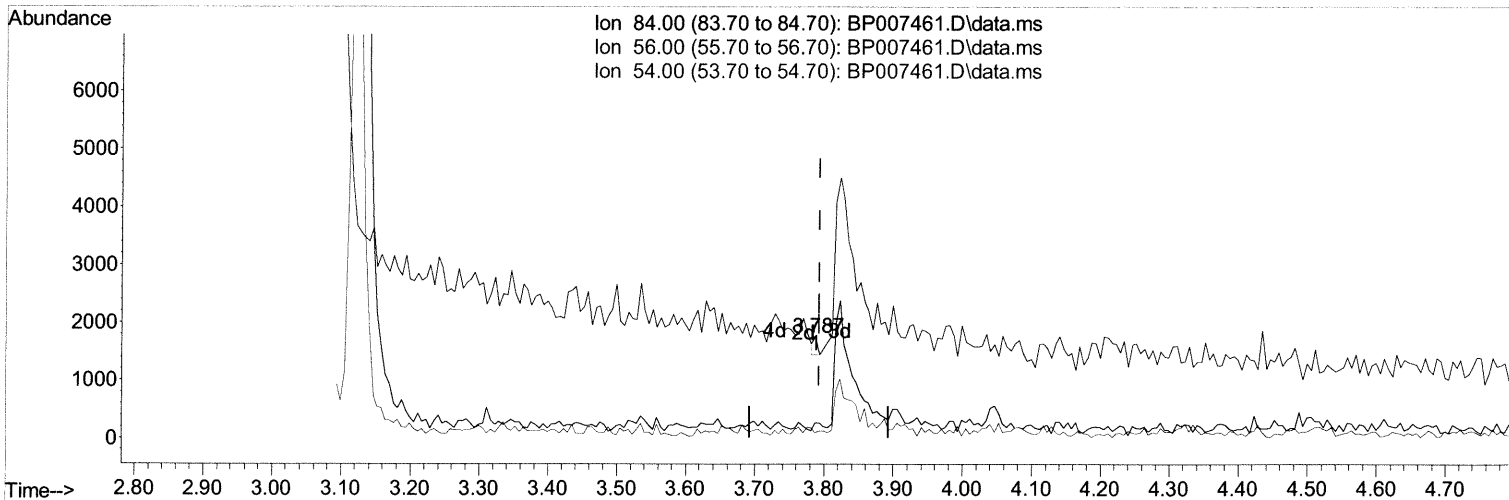
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TIC: BP007461.D\data.ms

(4) Pyridine-d5 (S)

3.787min (-0.006) 0.01 ng/ul

response 106

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	14.74#
54.00	31.70	4.99#
0.00	0.00	0.00

Quantitation Report (Qedit)

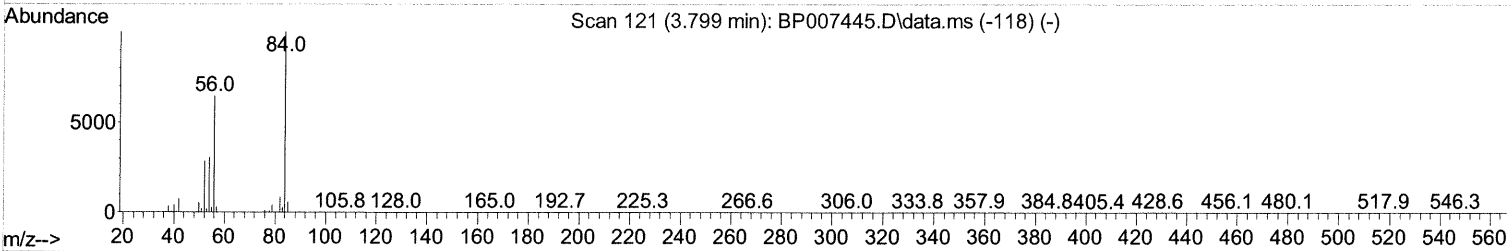
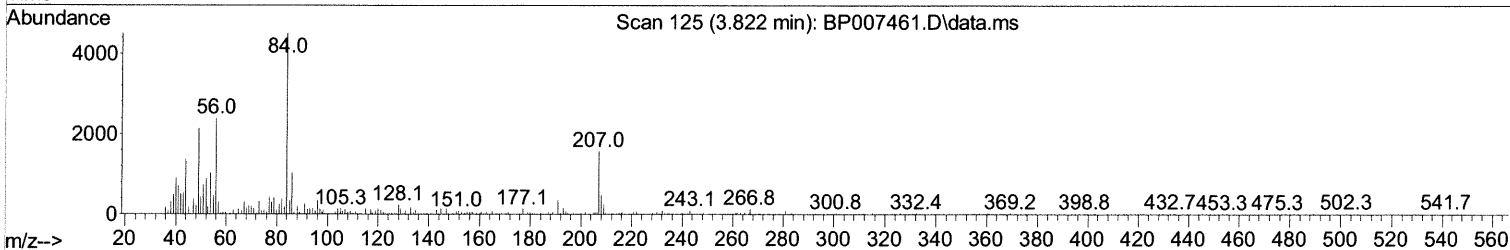
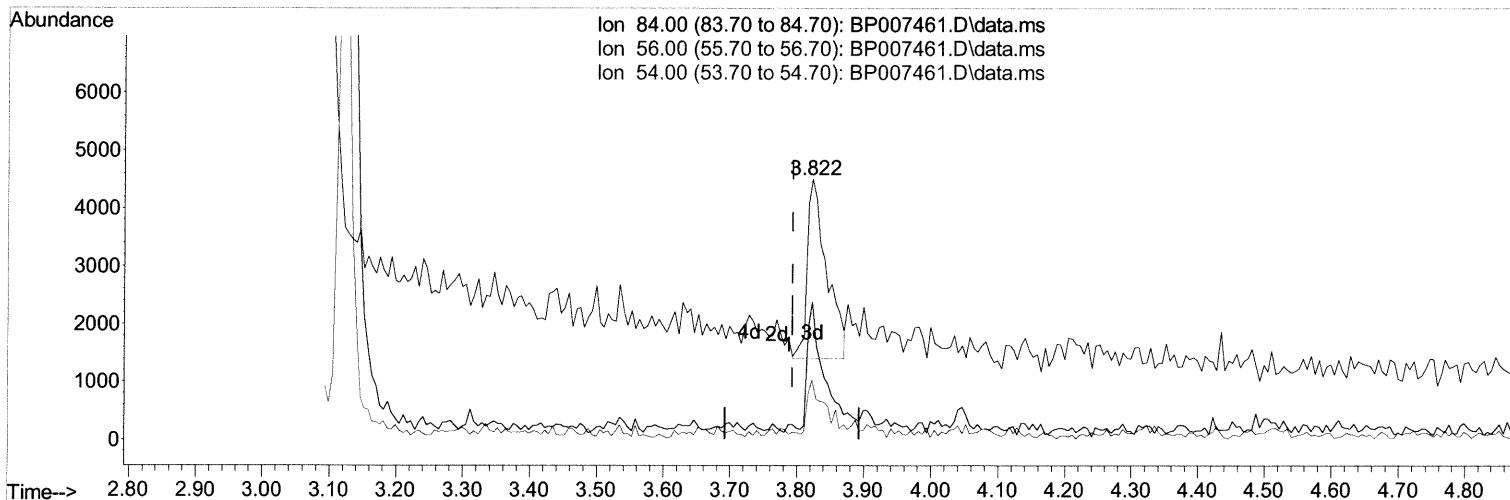
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TIC: BP007461.D\data.ms

(4) Pyridine-d5 (S)

3.822min (+ 0.029) 0.44 ng/ul m 10/20/21 JU

response 6236

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	52.96#
54.00	31.70	22.87#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007461.D
 Acq On : 18 Oct 2021 13:32
 Operator : CG/JU
 Sample : M4181-18
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B16

Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.958	152	205087	20.000 ng/ul	0.00
20) Naphthalene-d8	10.763	136	838935	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.598	164	553542	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.351	188	1176426	20.000 ng/ul	0.00
79) Chrysene-d12	21.439	240	1205953	20.000 ng/ul	0.00
88) Perylene-d12	23.898	264	1258676	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.381	96	16257	3.131 ng/ul	0.00
4) Pyridine-d5	3.822	84	6236m	0.439 ng/ul	0.03
7) Phenol-d5	7.116	99	350633	21.892 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	206130	21.578 ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	275070	21.831 ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	277471	22.491 ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	130696	24.617 ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	137065	27.639 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	280562	23.195 ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	385763	23.584 ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	904554	23.620 ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1073545	22.306 ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	131118	21.338 ng/ul	0.00
60) Fluorene-d10	15.592	176	769356	24.006 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	82916	17.590 ng/ul	0.00
73) Anthracene-d10	17.451	188	1200508	24.029 ng/ul	0.00
81) Pyrene-d10	19.680	212	1343674	22.288 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.739	264	1361327	21.736 ng/ul	0.00

Target Compounds Qvalue

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