

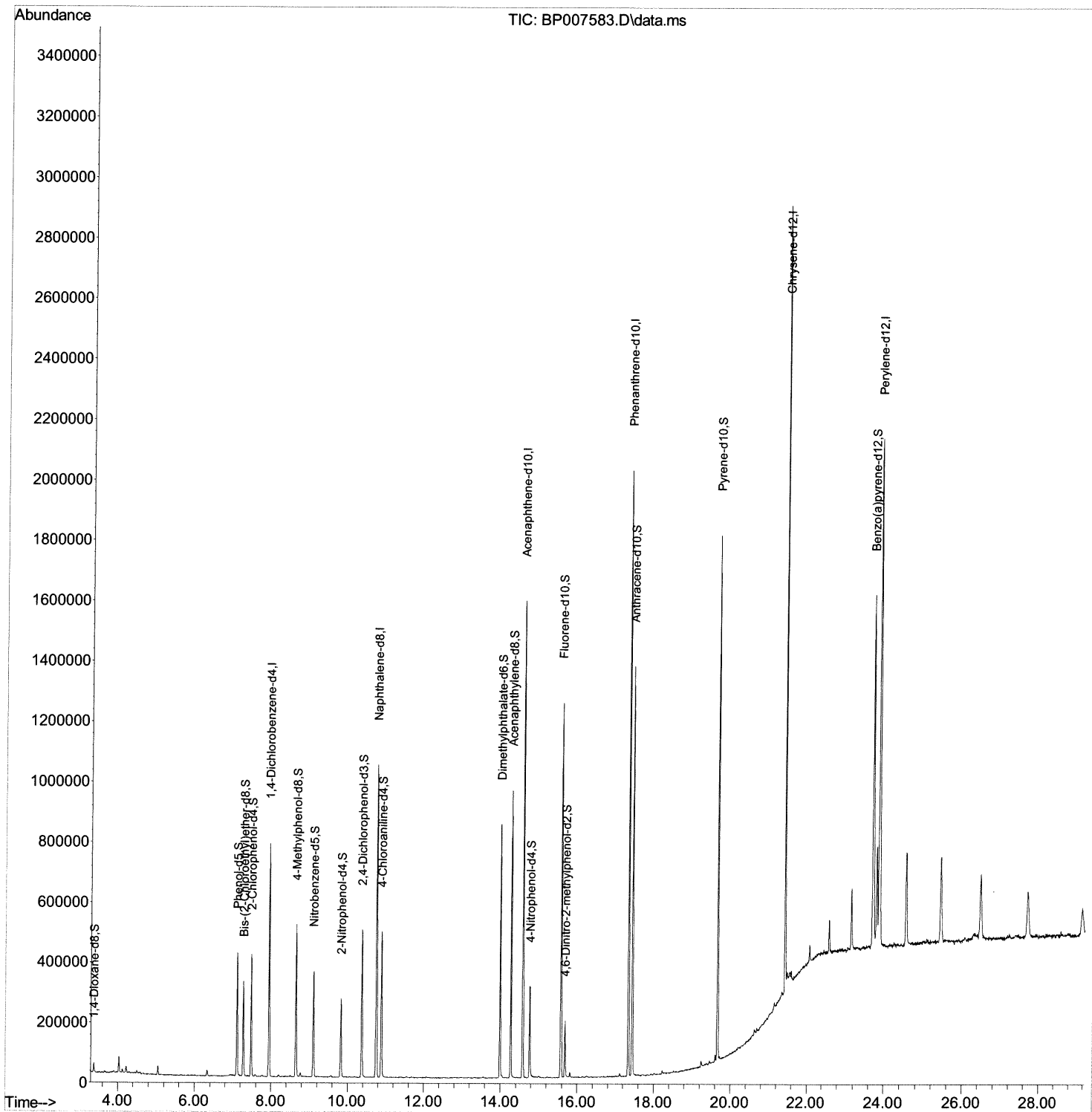
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007583.D
Acq On : 22 Oct 2021 11:09
Operator : CG/JU
Sample : M4217-14
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGF8

Manual Integrations
APPROVED

mohammad
10/22/2021 2:55:16 PM

Quant Time: Oct 22 13:07:39 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



Quantitation Report (Qedit)

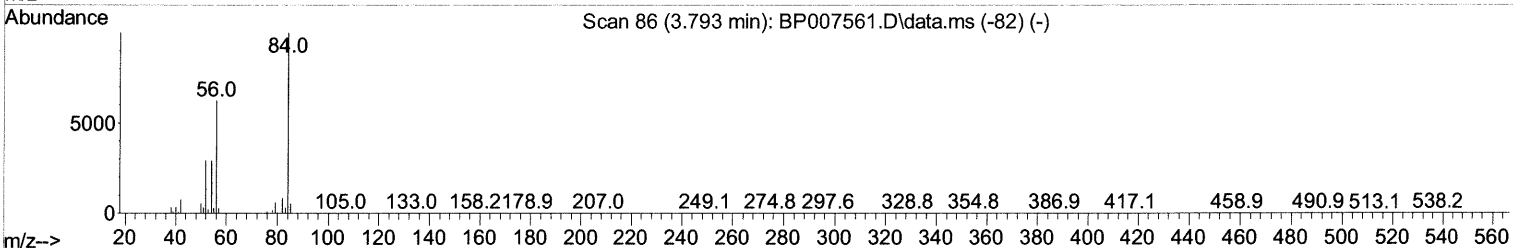
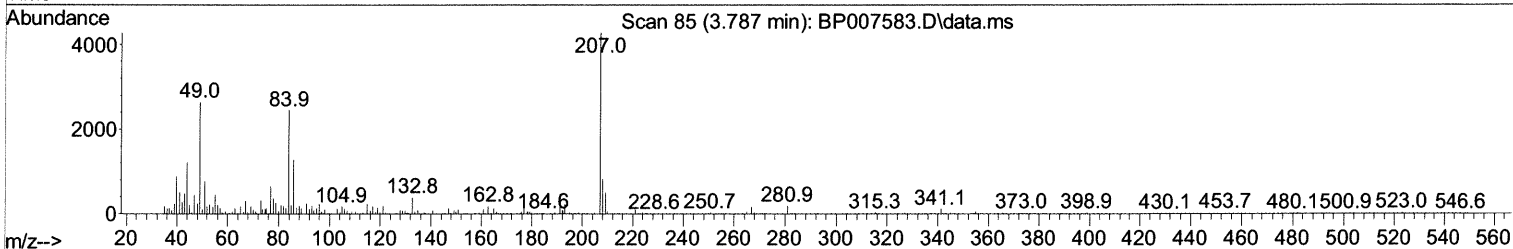
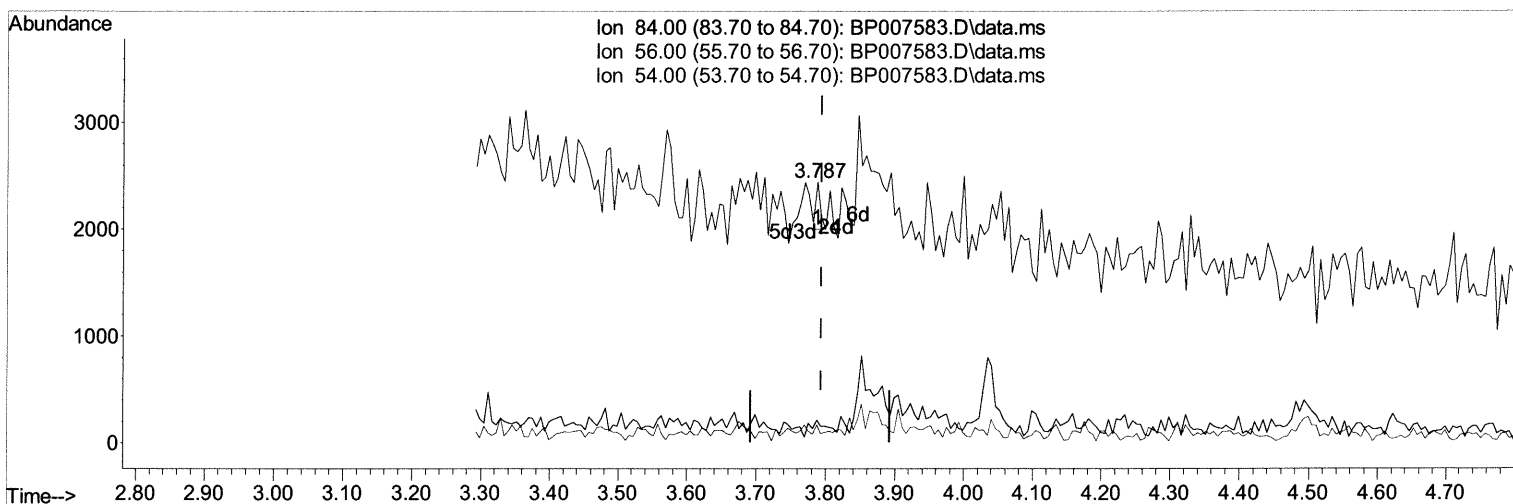
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TIC: BP007583.D\data.ms

(4) Pyridine-d5 (S)

3.787min (-0.006) 0.01 ng/ul

response 195

Ion	Exp%	Act%
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84.00	100.00	100.00
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56.00	66.30	8.60#
-------	-------	-------

54.00	31.70	6.72#
-------	-------	-------

0.00	0.00	0.00
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Quantitation Report (Qedit)

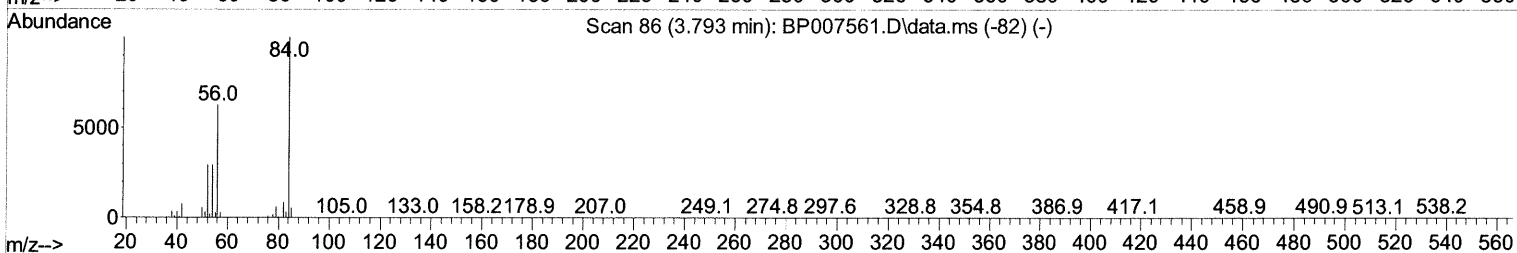
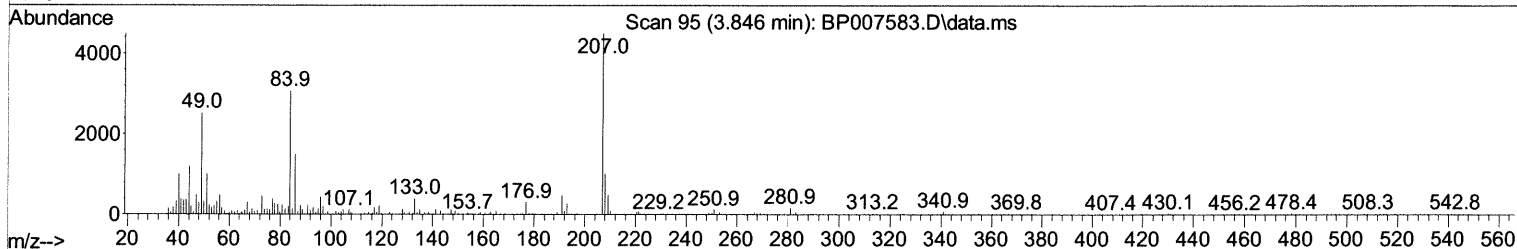
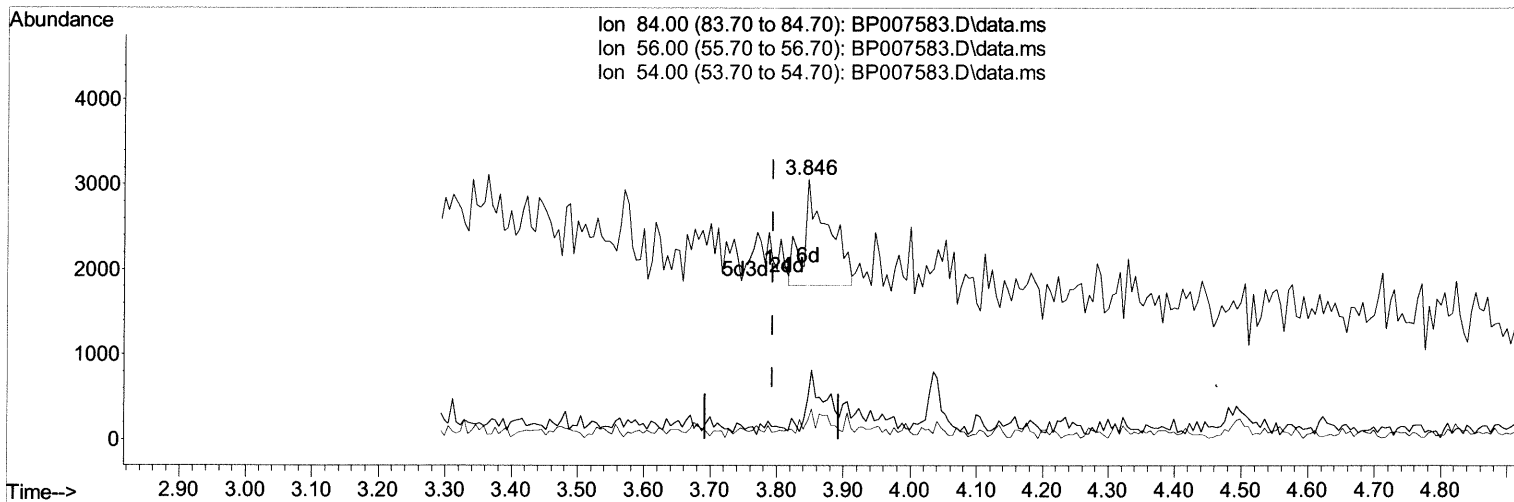
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TIC: BP007583.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.053) 0.22 ng/ul m 10/26/21JU

response 3259

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	15.79#
54.00	31.70	7.18#
0.00	0.00	0.00

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 Data File : BP007583.D
 Acq On : 22 Oct 2021 11:09
 Operator : CG/JU
 Sample : M4217-14
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGF8

Manual Integrations
 APPROVED

mohammad
 10/22/2021 2:55:16 PM

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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.946	152	211885	20.000 ng/ul	0.00
20) Naphthalene-d8	10.751	136	866486	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.581	164	564300	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1222678	20.000 ng/ul	0.00
79) Chrysene-d12	21.421	240	1269669	20.000 ng/ul	0.00
88) Perylene-d12	23.874	264	1304828	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.375	96	15249	2.842 ng/uL	0.00
4) Pyridine-d5	3.846	84	3259m>	0.222 ng/ul>	0.05
7) Phenol-d5	7.105	99	235659	14.242 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	131990	13.374 ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	185350	14.239 ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	184651	14.487 ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	84298	15.373 ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	85290	16.652 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	180988	14.487 ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	252093	14.922 ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	545134	13.963 ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	673706	13.731 ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	74408	11.878 ng/ul	0.00
60) Fluorene-d10	15.575	176	492904	15.087 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	40227	8.211 ng/ul	0.00
73) Anthracene-d10	17.433	188	772409	14.875 ng/ul	0.00
81) Pyrene-d10	19.669	212	913391	14.390 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	869038	13.385 ng/ul	0.00

Target Compounds Qvalue

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