

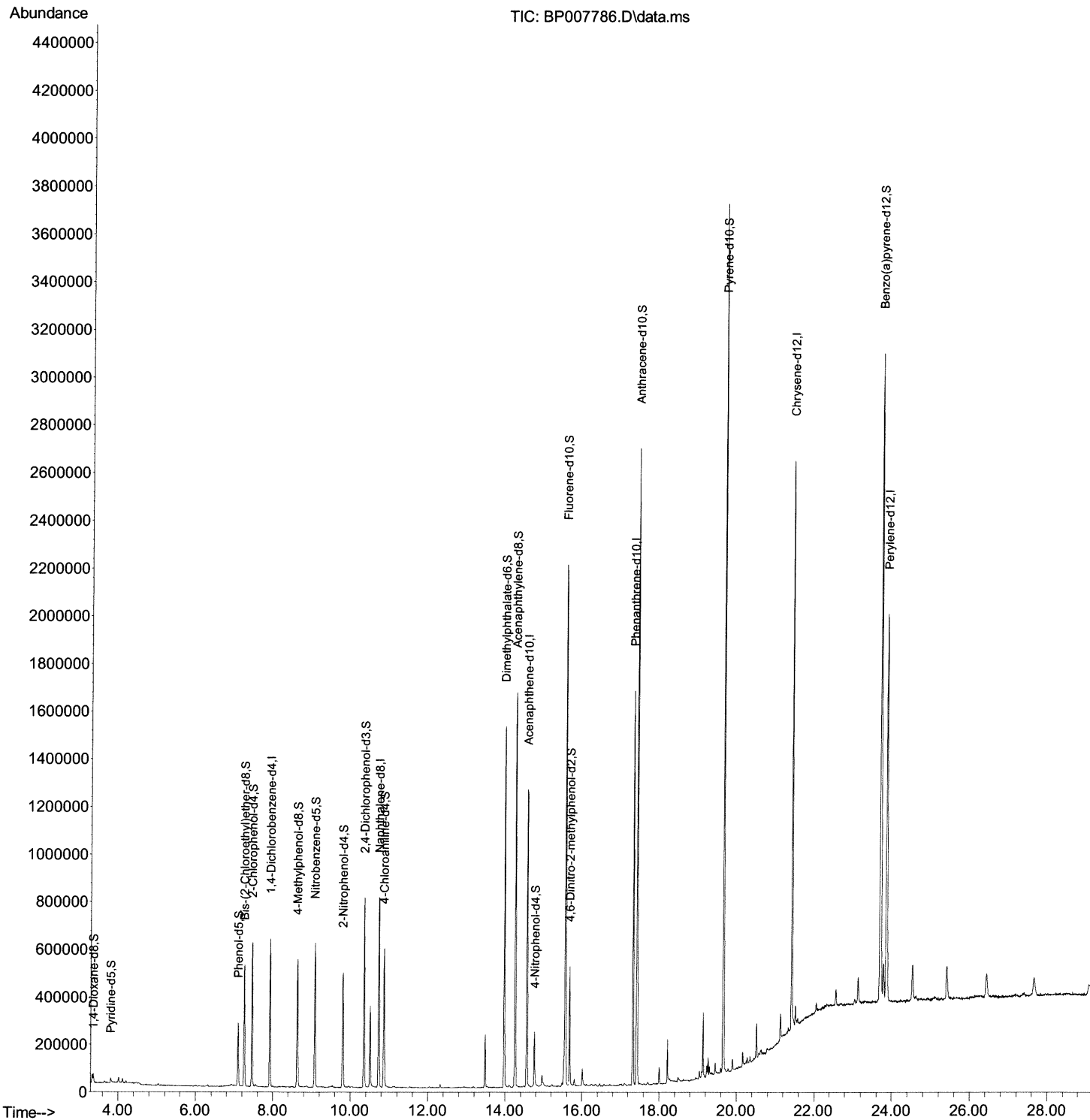
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007786.D
Acq On : 29 Oct 2021 17:02
Operator : CG/JU
Sample : M4282-13
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY83

Manual Integrations
APPROVED

mohammad
11/1/2021 1:36:25 PM

Quant Time: Oct 29 17:28:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 27 15:37:43 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

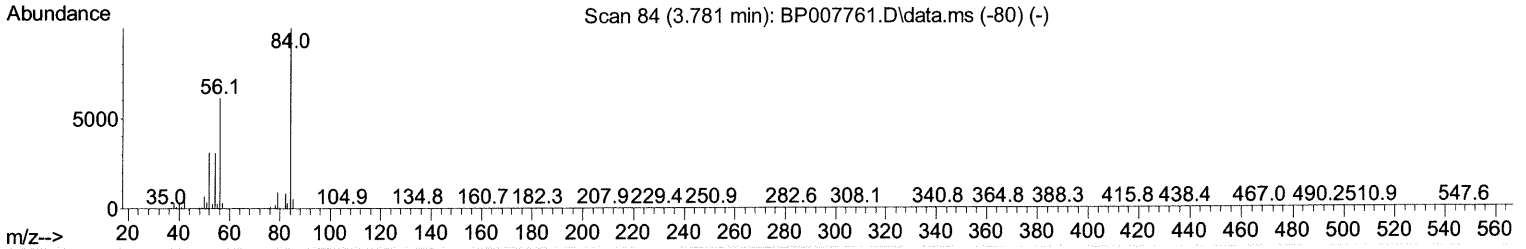
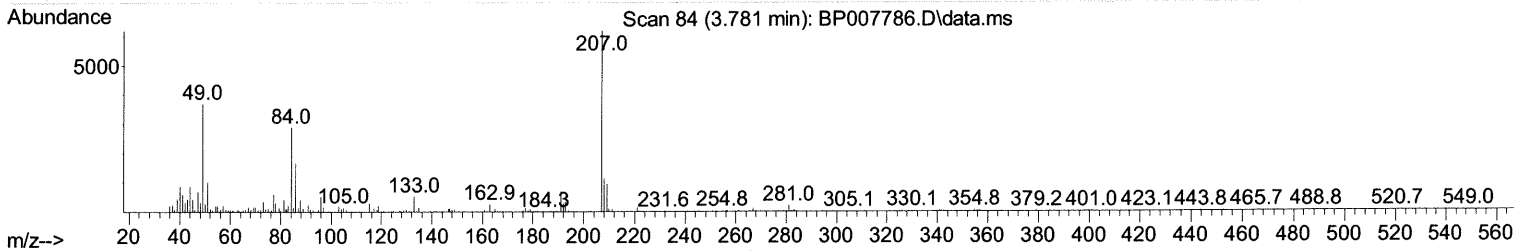
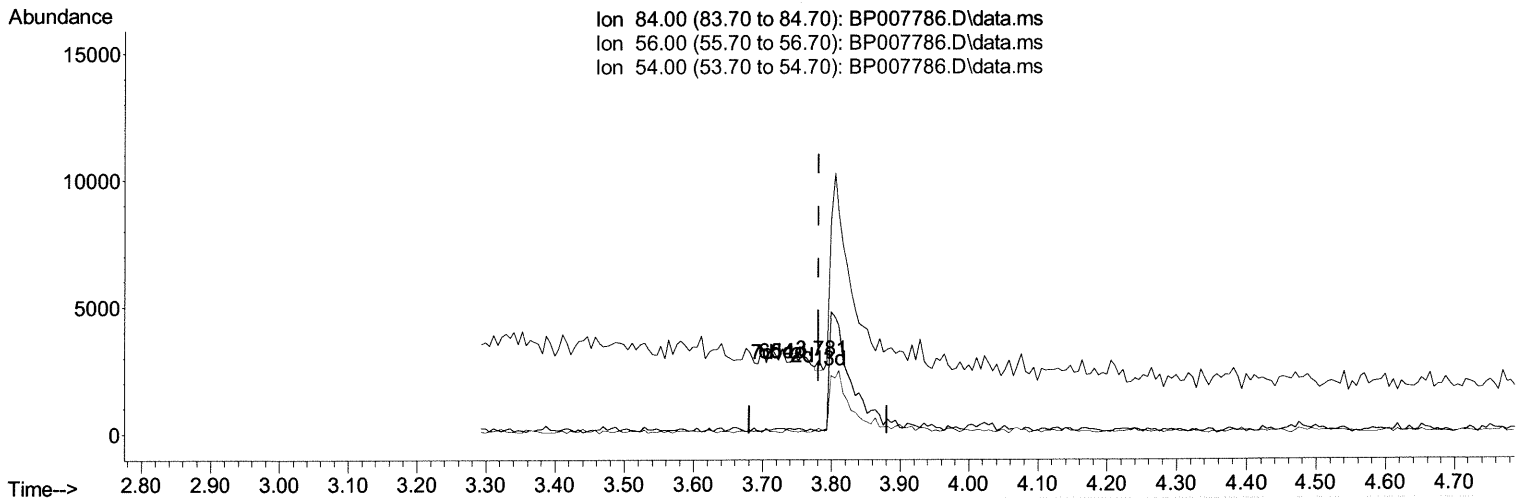
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TIC: BP007786.D\data.ms

(4) Pyridine-d5 (S)

3.781min (+ 0.000) 0.01 ng/ul

response 135

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	3.21#
54.00	31.70	7.40#
0.00	0.00	0.00

Quantitation Report (Qedit)

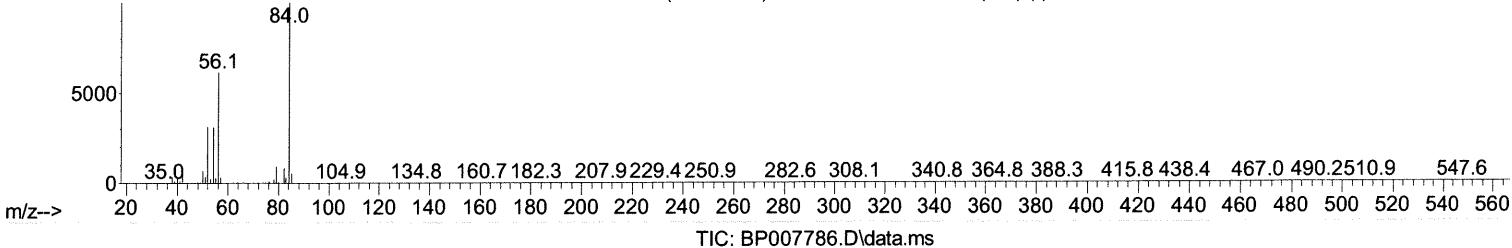
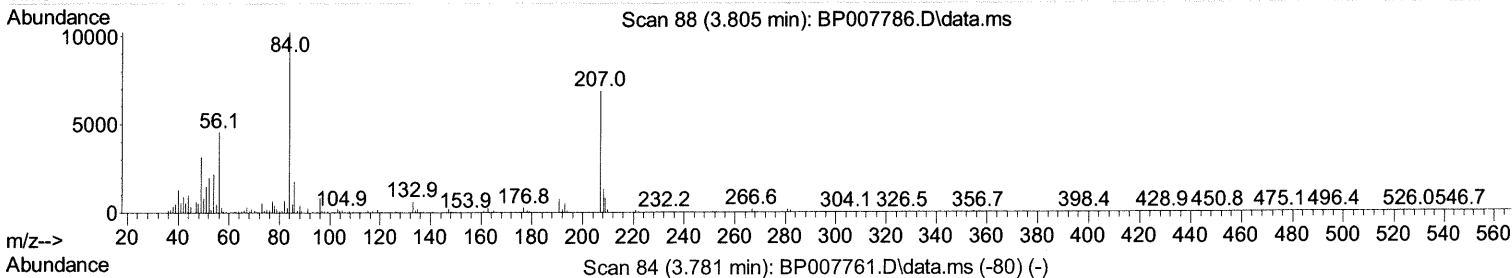
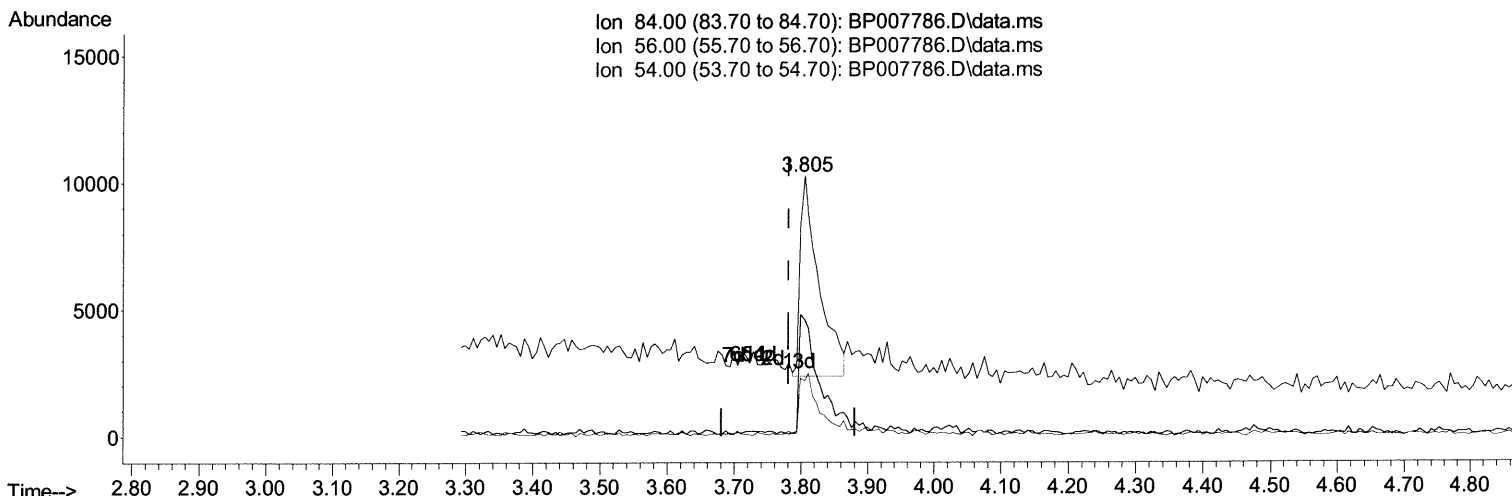
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Manual Integrations
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(4) Pyridine-d5 (S)

3.805min (+ 0.024) 1.36 ng/ul m 11/02/21 JU

response 15247

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	44.72#
54.00	31.70	21.43#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
 Data File : BP007786.D
 Acq On : 29 Oct 2021 17:02
 Operator : CG/JU
 Sample : M4282-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY83

Manual Integrations
 APPROVED

mohammad
 11/1/2021 1:36:25 PM

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	169009	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	678426	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.563	164	446784	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.322	188	991986	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	1187878	20.000	ng/ul	-0.01
88) Perylene-d12	23.857	264	1307681	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	18018	4.843	ng/ul	0.00
4) Pyridine-d5	3.805	84	15247m	1.355	ng/ul	0.02 11/02/21 JU
7) Phenol-d5	7.093	99	150169	11.698	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	218946	30.720	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	272454	27.289	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	200495	19.946	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	150156	32.478	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	157935	31.277	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	296419	28.840	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	307733	22.644	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	996751	32.935	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1155868	31.161	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	58673	10.576	ng/ul	0.00
60) Fluorene-d10	15.557	176	868195	34.029	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	113252	21.344	ng/ul	0.00
73) Anthracene-d10	17.422	188	1506097	36.944	ng/ul	0.00
81) Pyrene-d10	19.657	212	1927795	34.858	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	2168635	33.759	ng/ul	-0.01

Target Compounds Qvalue

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