

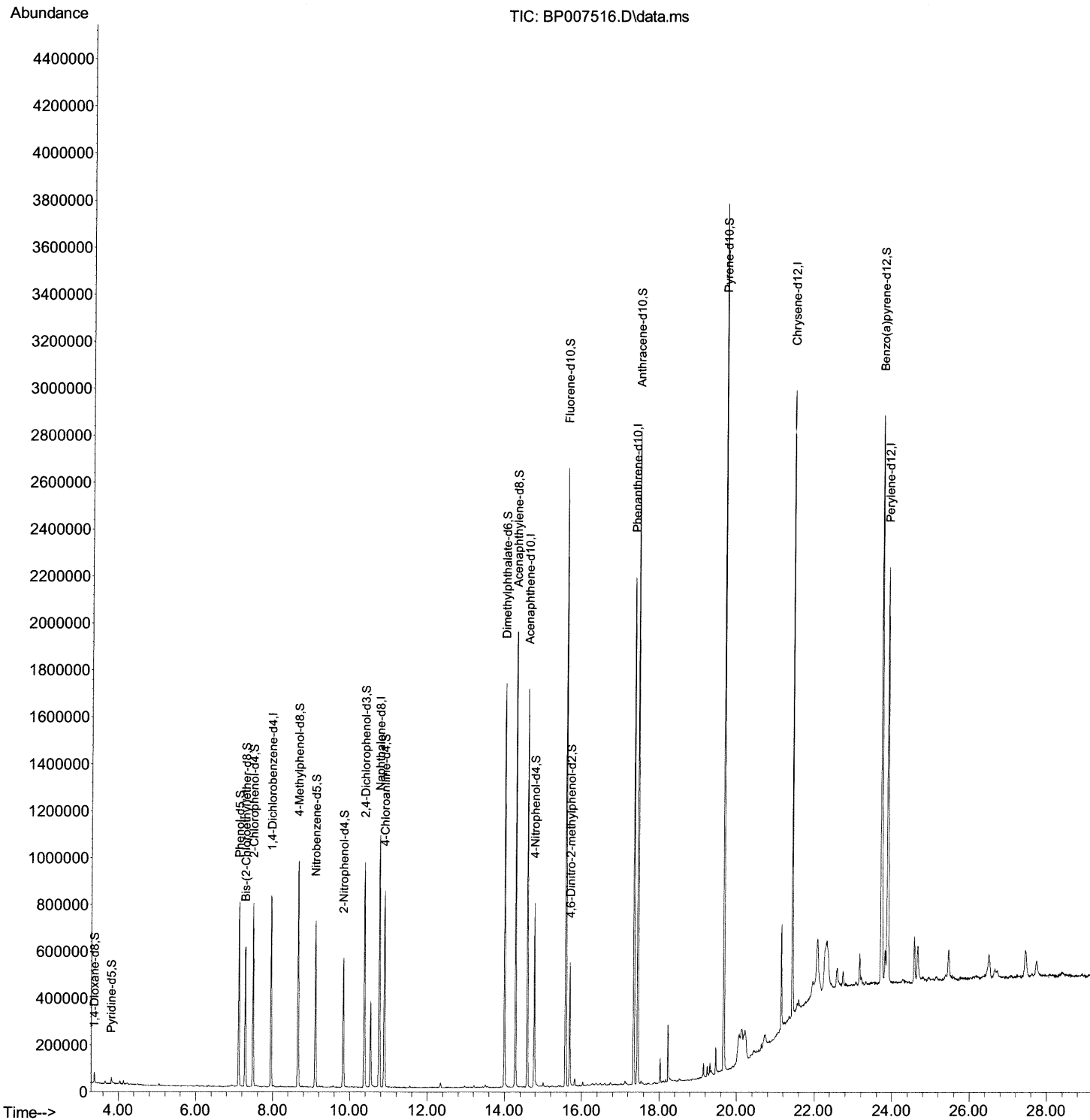
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007516.D
Acq On : 20 Oct 2021 07:46
Operator : CG/JU
Sample : M4218-04
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGF1

Manual Integrations
APPROVED

mohammad
10/20/2021 1:57:18 PM

Quant Time: Oct 20 08:18:44 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)

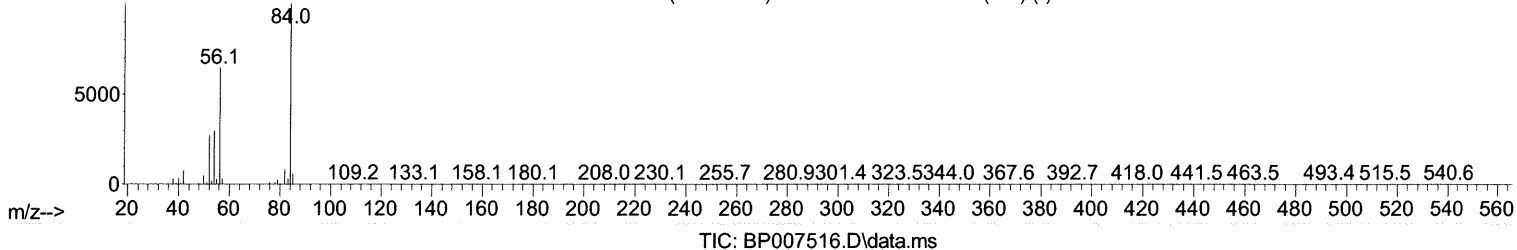
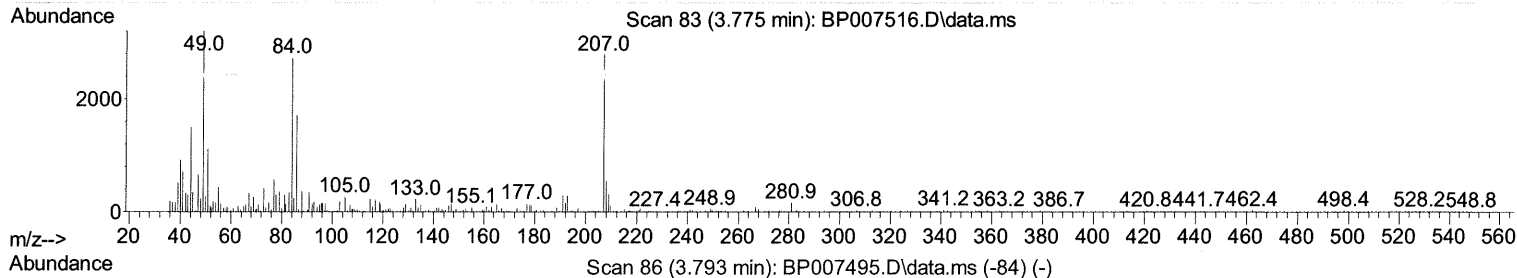
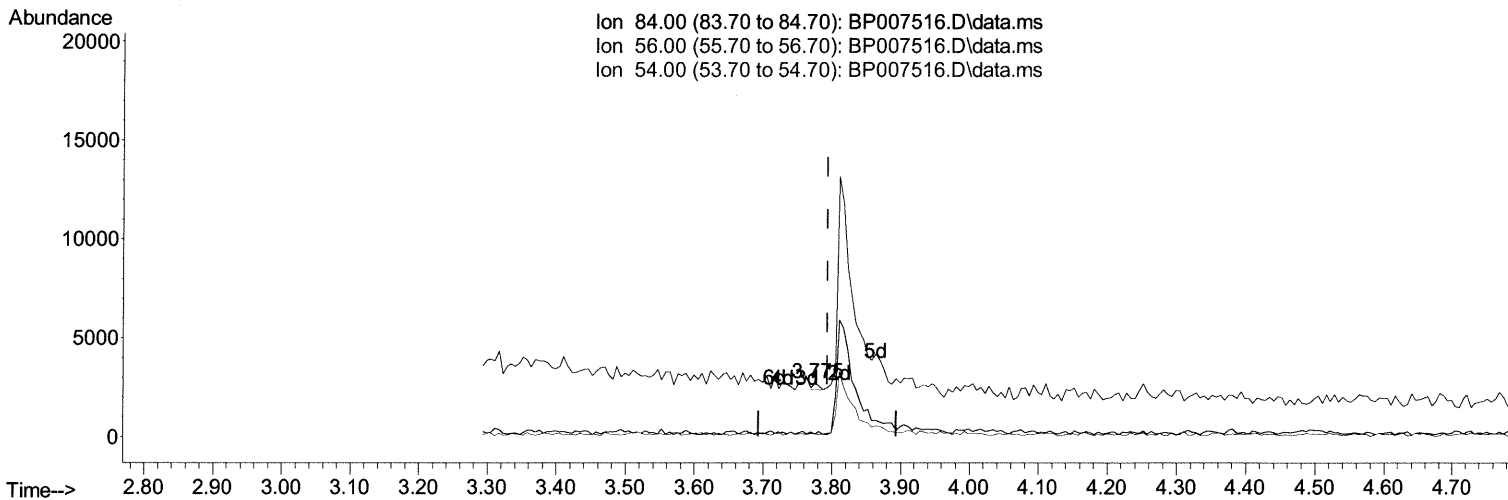
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(4) Pyridine-d5 (S)

3.775min (-0.018) 0.01 ng/ul

response 185

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	5.35#
54.00	31.70	5.79#
0.00	0.00	0.00

Quantitation Report (Qedit)

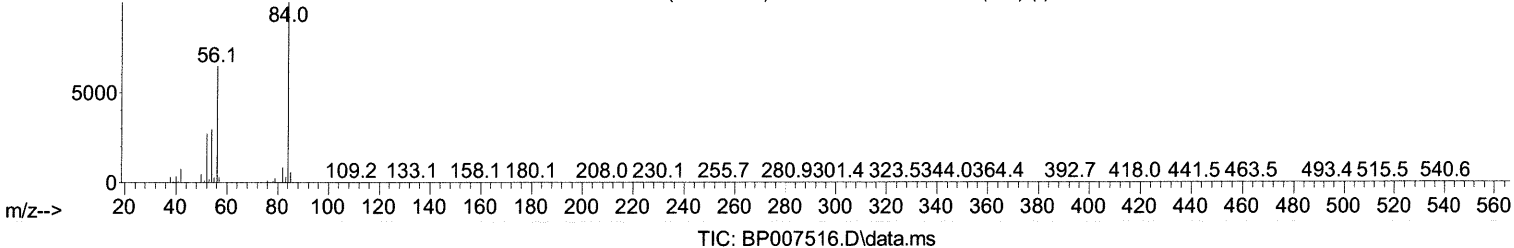
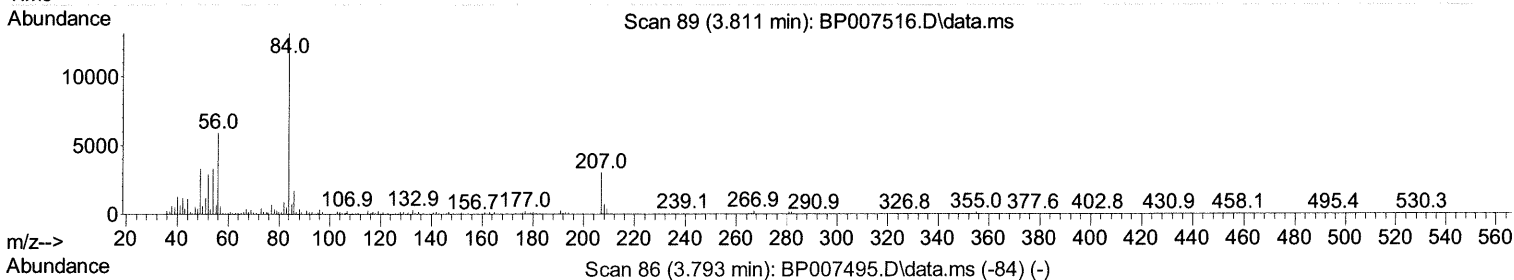
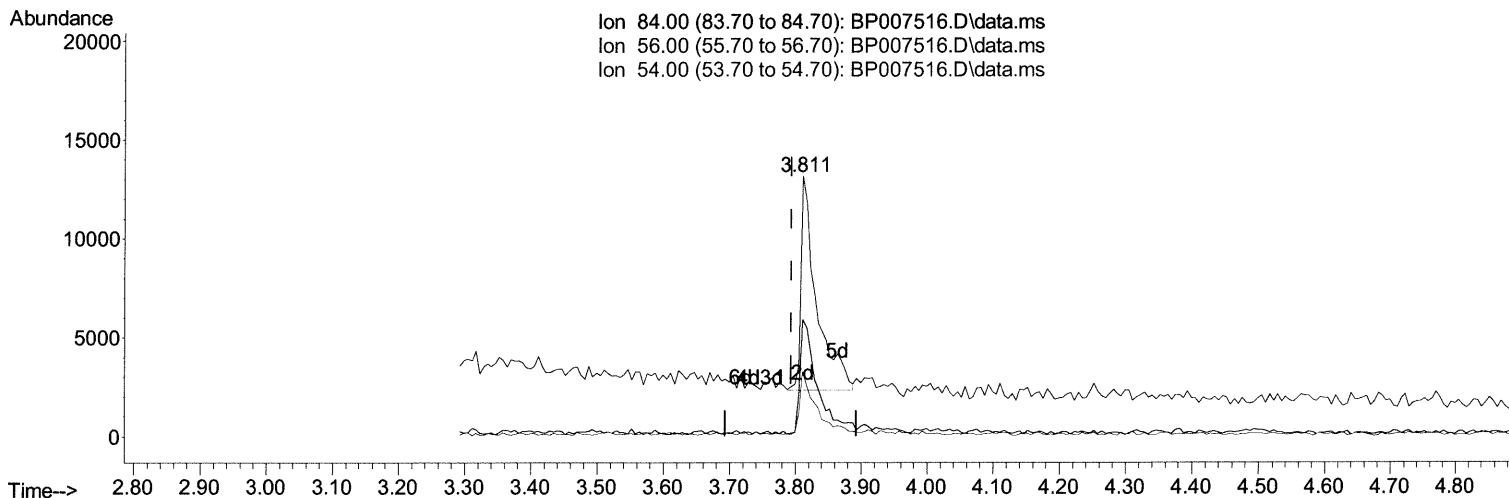
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Manual Integrations
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(4) Pyridine-d5 (S)

3.811min (+ 0.018) 1.16 ng/ul m 10/20/21 JU

response 17756

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	45.05#
54.00	31.70	25.01#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007516.D
 Acq On : 20 Oct 2021 07:46
 Operator : CG/JU
 Sample : M4218-04
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	221374	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	900309	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.586	164	586908	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.339	188	1265813	20.000	ng/ul	-0.01
79) Chrysene-d12	21.427	240	1347889	20.000	ng/ul	-0.01
88) Perylene-d12	23.874	264	1379419	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	24261	4.328	ng/uL	0.00
4) Pyridine-d5	3.811	84	17756m>	1.157	ng/ul >	0.02
7) Phenol-d5	7.110	99	451349	26.107	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	258704	25.089	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	360221	26.486	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	362089	27.191	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	172105	30.207	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.834	143	179936	33.810	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	355077	27.354	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	434722	24.766	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	1159262	28.550	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1403749	27.509	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	180780	27.747	ng/ul	0.00
60) Fluorene-d10	15.581	176	998625	29.388	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	117439	23.154	ng/ul	-0.01
73) Anthracene-d10	17.439	188	1622657	30.185	ng/ul	-0.01
81) Pyrene-d10	19.669	212	1942991	28.835	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.715	264	1927086	28.076	ng/ul	-0.02

Target Compounds Qvalue

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