

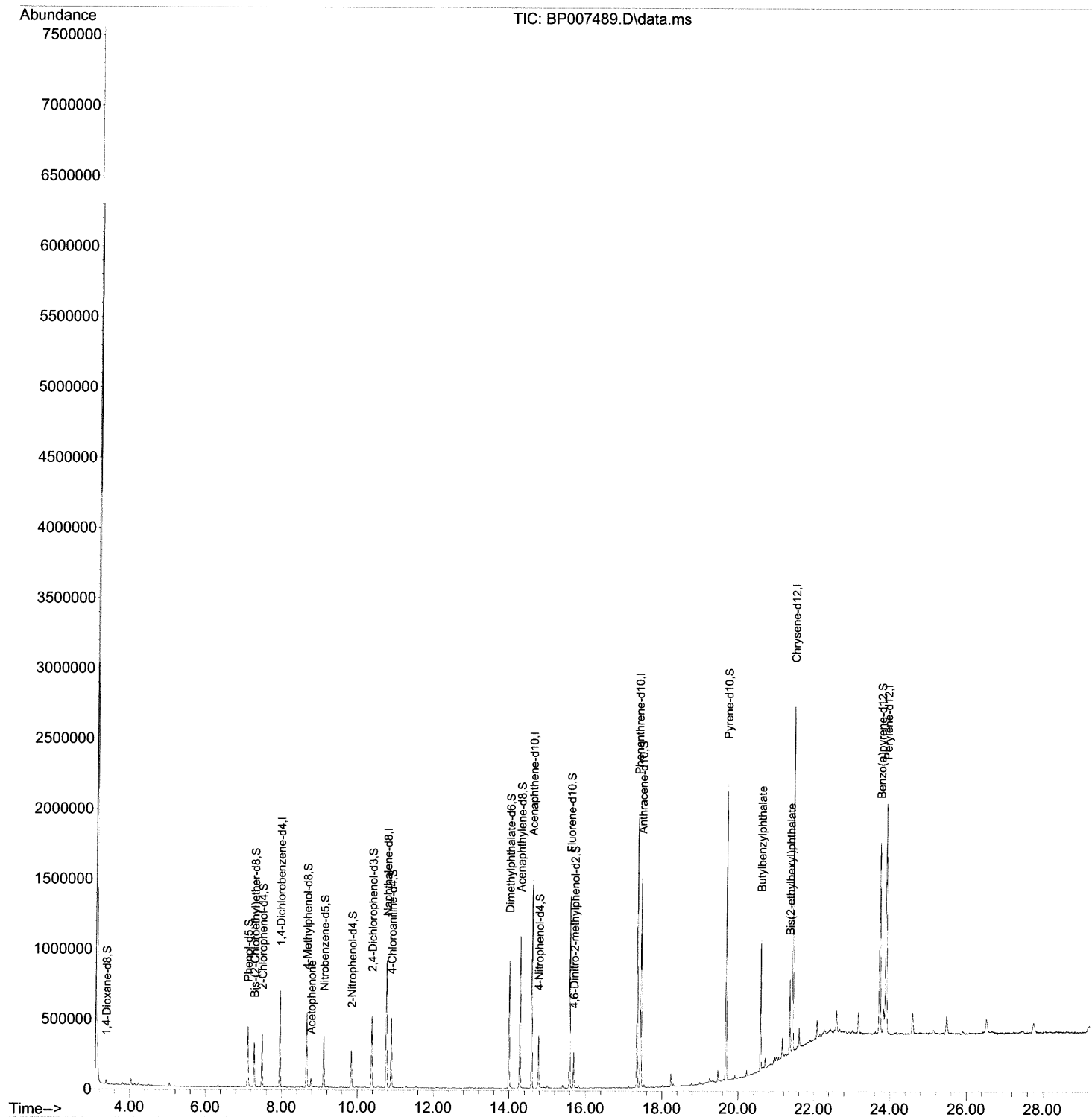
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007489.D
Acq On : 19 Oct 2021 11:57
Operator : CG/JU
Sample : M4196-19
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGD4

Manual Integrations
APPROVED

mohammad
10/20/2021 1:56:37 PM

Quant Time: Oct 19 12:27:05 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



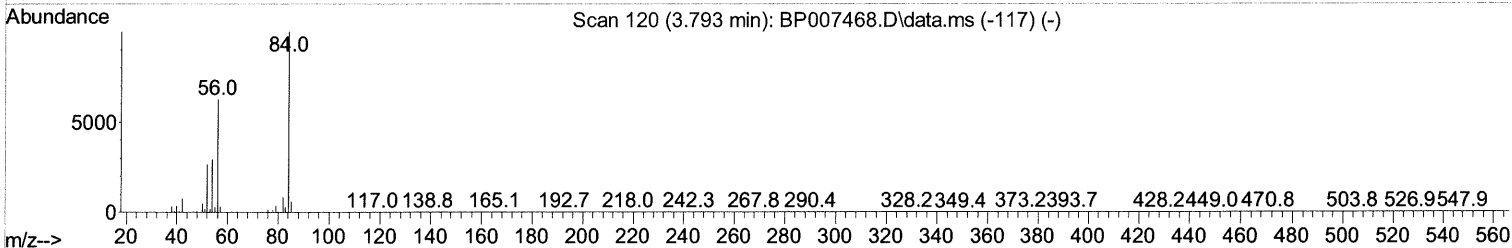
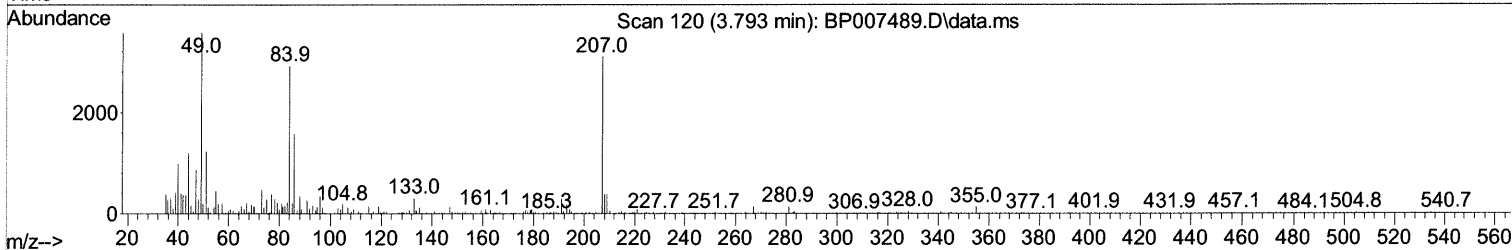
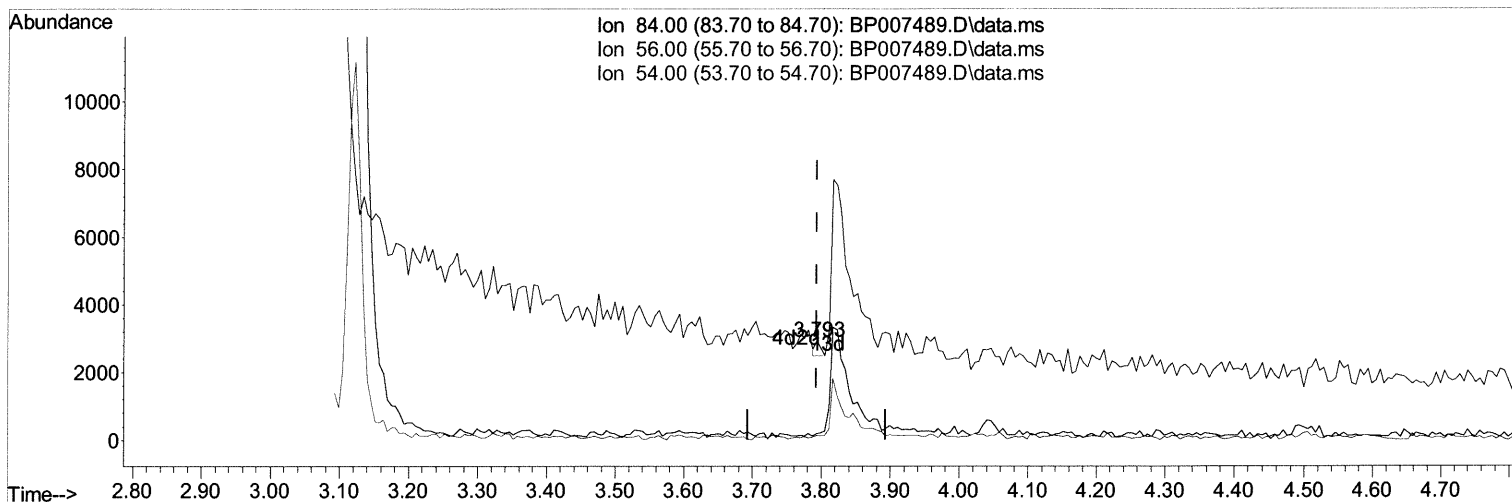
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TIC: BP007489.D\data.ms

(4) Pyridine-d5 (S)

3.793min (-0.000) 0.02 ng/ul

response 230

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	6.53#
54.00	31.70	4.51#
0.00	0.00	0.00

Quantitation Report (Qedit)

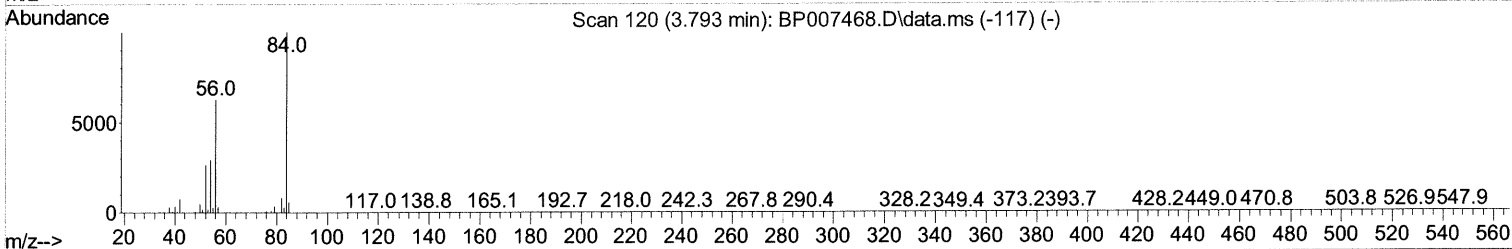
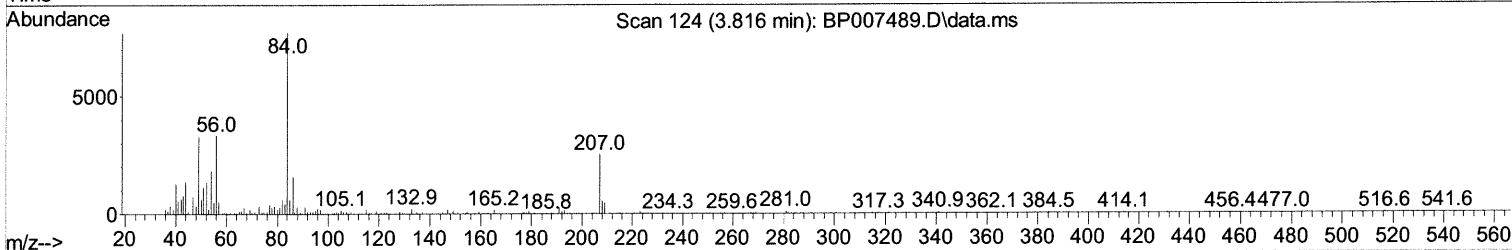
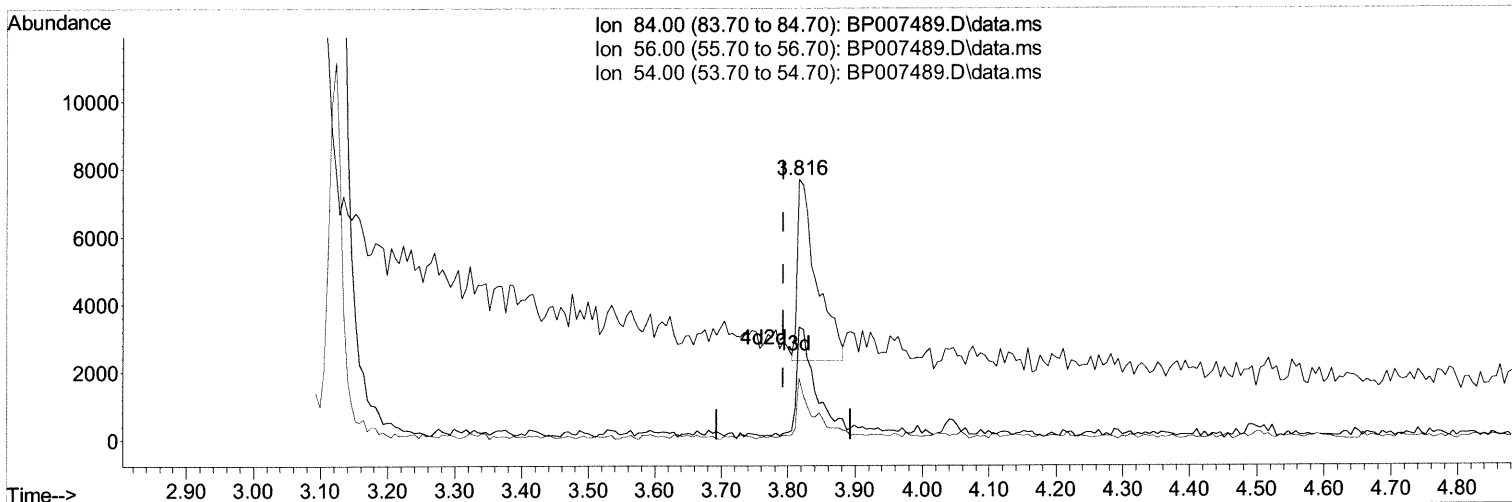
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Manual Integrations
 APPROVED

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 10/20/2021 1:56:37 PM

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 Response via : Initial Calibration



TIC: BP007489.D\data.ms

(4) Pyridine-d5 (S)

3.816min (+ 0.023) 0.79 ng/ul m 10/21/21 JU

response 10598

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	43.24#
54.00	31.70	23.71#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007489.D
 Acq On : 19 Oct 2021 11:57
 Operator : CG/JU
 Sample : M4196-19
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGD4

Manual Integrations
 APPROVED

mohammad
 10/20/2021 1:56:37 PM

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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	193008	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	786090	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	516991	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1129502	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	1227752	20.000	ng/ul	-0.01
88) Perylene-d12	23.880	264	1254726	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	14787	3.026	ng/uL	0.00
4) Pyridine-d5	3.816	84	10598m	0.792	ng/ul	0.02
7) Phenol-d5	7.110	99	240512	15.957	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	136074	15.136	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	179805	15.164	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	189543	16.325	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	88354	17.761	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	87951	18.927	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	186599	16.464	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	267442	17.450	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	623761	17.439	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	762767	16.969	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	87550	15.255	ng/ul	0.00
60) Fluorene-d10	15.586	176	552675	18.464	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	54413	12.023	ng/ul	-0.01
73) Anthracene-d10	17.445	188	886675	18.485	ng/ul	0.00
81) Pyrene-d10	19.674	212	1083022	17.645	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1064271	17.047	ng/ul	-0.01
Target Compounds						
					Qvalue	
16) Acetophenone	8.775	105	33767	1.901	ng/ul	95
83) Butylbenzylphthalate	20.580	149	264643	9.386	ng/ul	97
86) Bis(2-ethylhexyl)phtha...	21.351	149	181017	4.199	ng/ul	99

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