

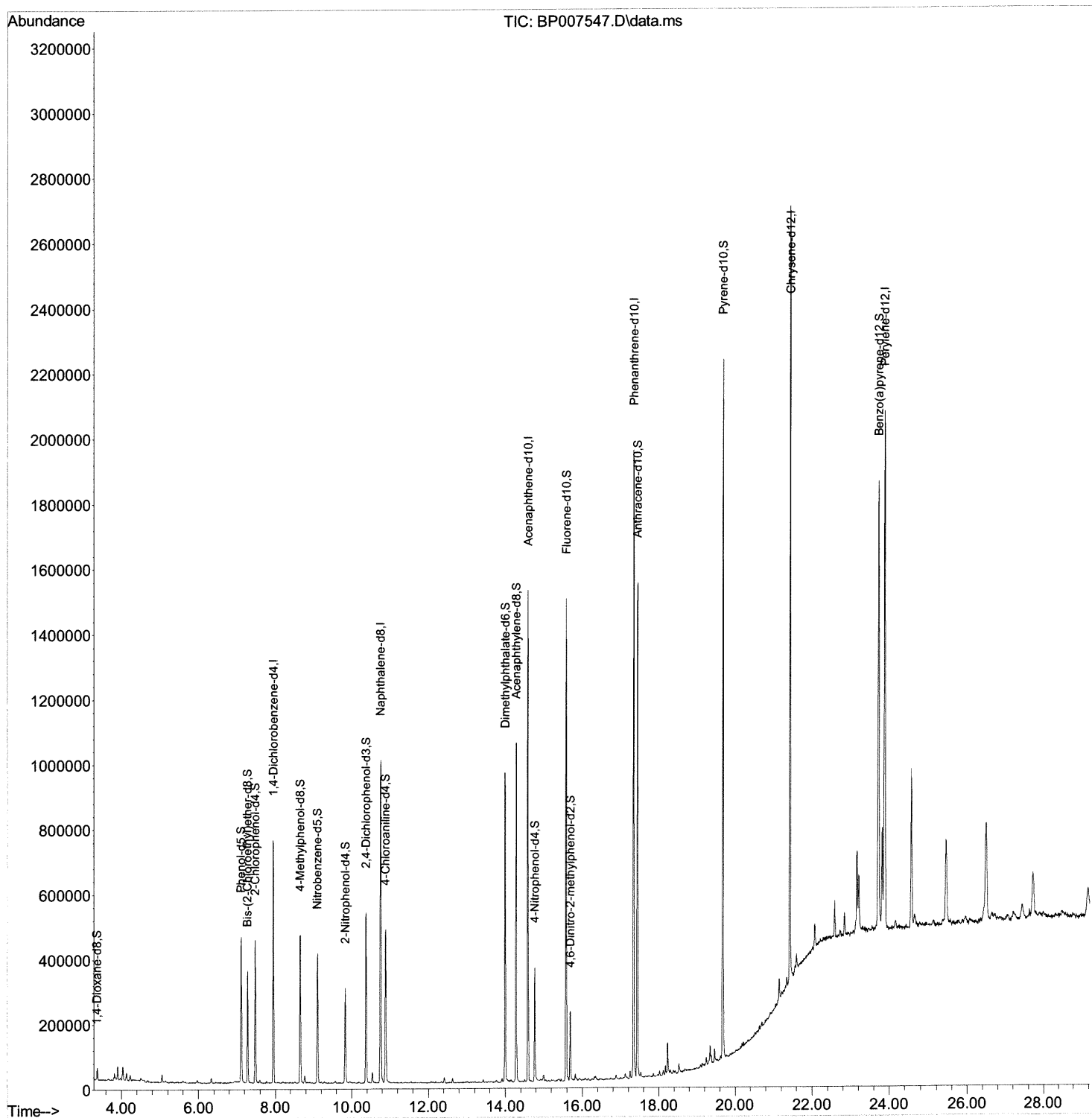
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007547.D
Acq On : 21 Oct 2021 10:39
Operator : CG/JU
Sample : M4182-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B52

Manual Integrations
APPROVED

mohammad
10/22/2021 2:54:44 PM

Quant Time: Oct 21 11:25:07 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



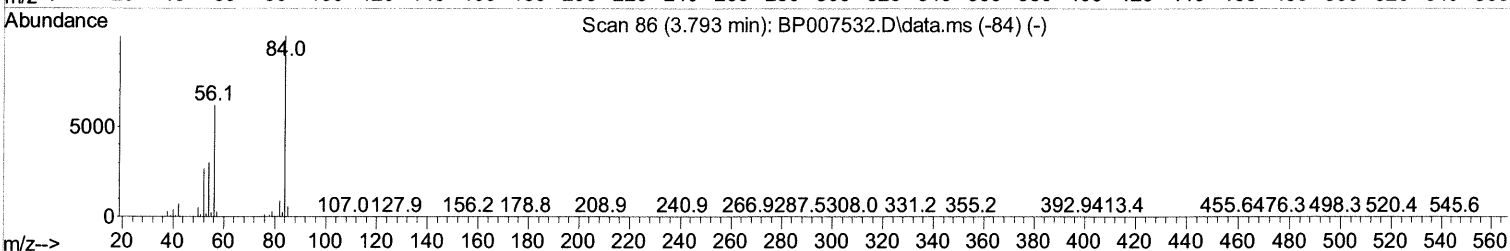
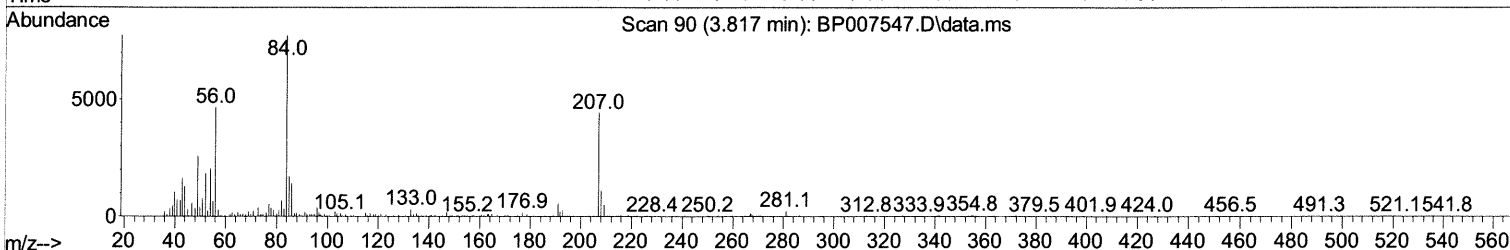
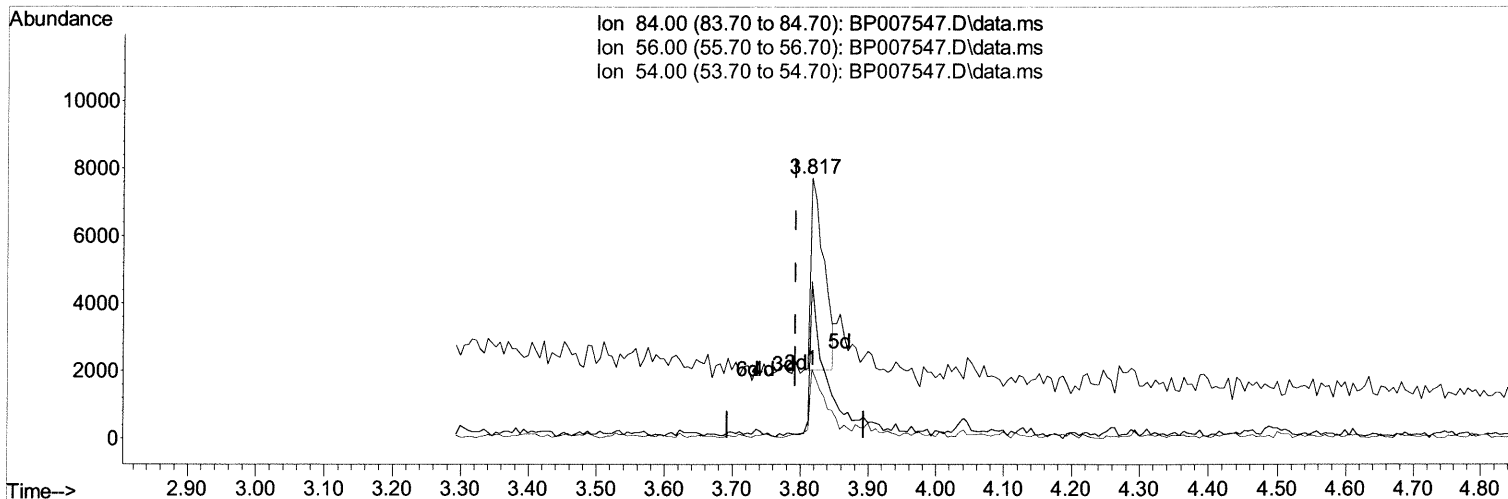
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TIC: BP007547.D\data.ms

(4) Pyridine-d5 (S)

3.817min (+ 0.024) 0.54 ng/ul

response 7473

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	60.24
54.00	31.70	26.58
0.00	0.00	0.00

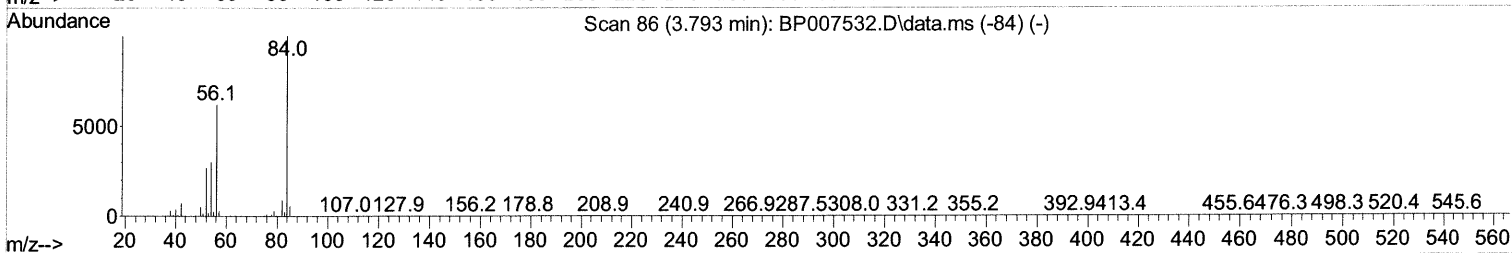
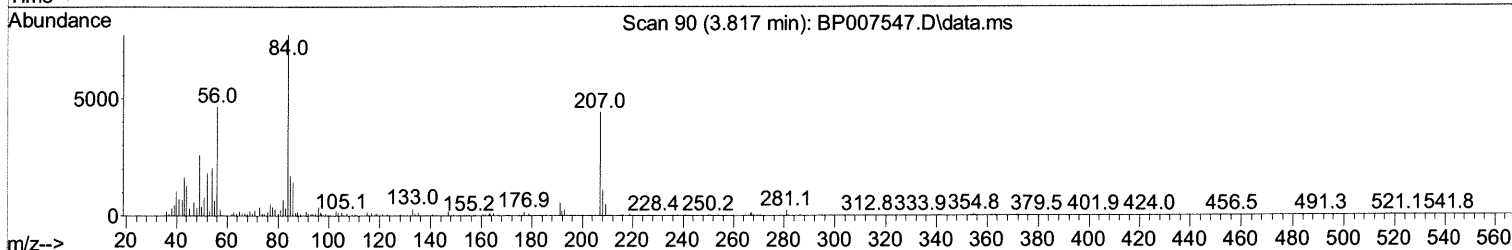
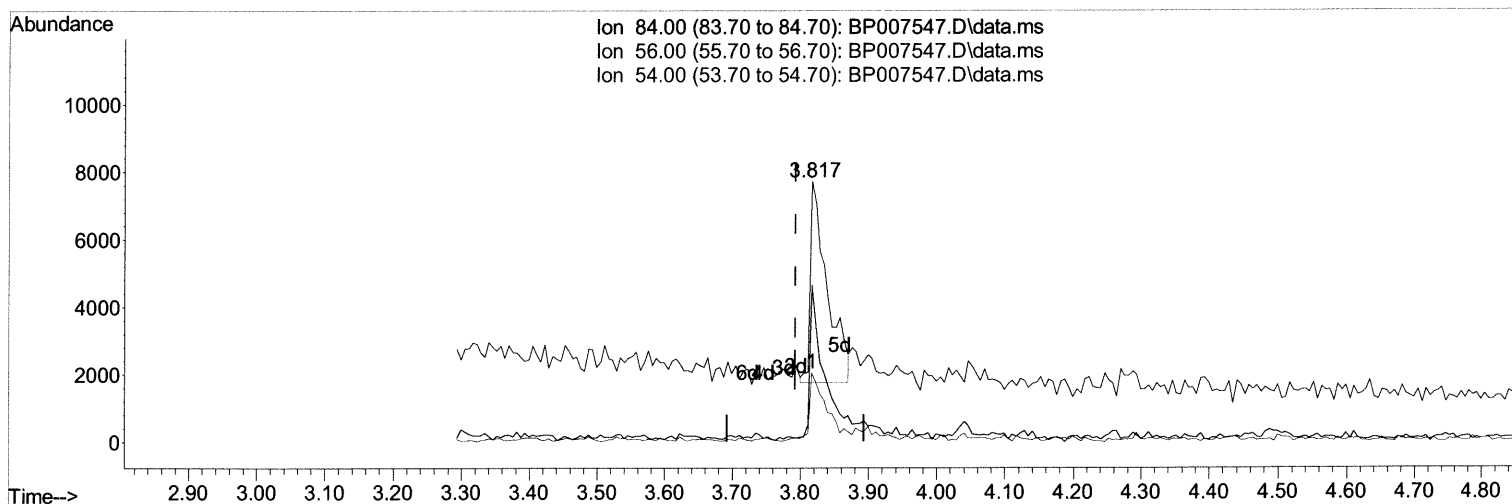
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TIC: BP007547.D\data.ms

(4) Pyridine-d5 (S)

3.817min (+ 0.024) 0.74 ng/ul m 10/21/21 JU

response 10263

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	60.24
54.00	31.70	26.58
0.00	0.00	0.00

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Manual Integrations
 APPROVED

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.952	152	200861	20.000 ng/ul	0.00
20) Naphthalene-d8	10.757	136	822886	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.587	164	525182	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1129779	20.000 ng/ul	0.00
79) Chrysene-d12	21.421	240	1188364	20.000 ng/ul	0.00
88) Perylene-d12	23.874	264	1218362	20.000 ng/ul	0.00

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.375	96	19726	3.879 ng/ul	0.00
4) Pyridine-d5	3.817	84	10263m	0.737 ng/ul	0.02
7) Phenol-d5	7.110	99	255521	16.290 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	149532	15.983 ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	202936	16.445 ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	174092	14.408 ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	94454	18.138 ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	94294	19.385 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	197487	16.645 ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	253007	15.770 ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	635005	17.477 ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	769122	16.844 ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	78356	13.440 ng/ul	0.00
60) Fluorene-d10	15.581	176	561148	18.455 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	48751	10.769 ng/ul	0.00
73) Anthracene-d10	17.439	188	884621	18.437 ng/ul	0.00
81) Pyrene-d10	19.669	212	1087829	18.311 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.715	264	1050132	17.322 ng/ul	0.00

Target Compounds Qvalue

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