

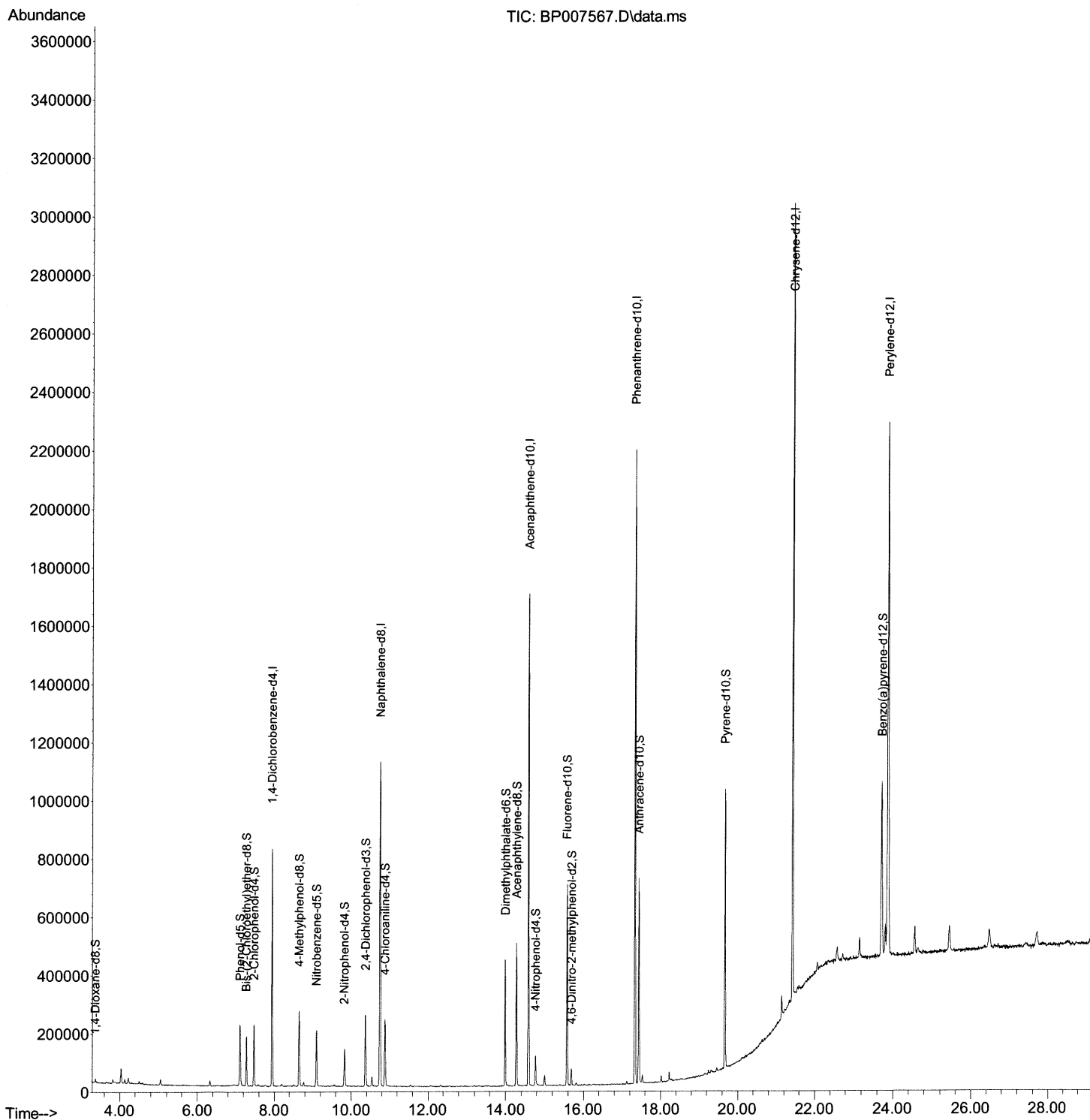
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007567.D
Acq On : 21 Oct 2021 22:37
Operator : CG/JU
Sample : M4217-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGB5

Manual Integrations
APPROVED

mohammad
10/22/2021 2:54:57 PM

Quant Time: Oct 22 01:00:55 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)

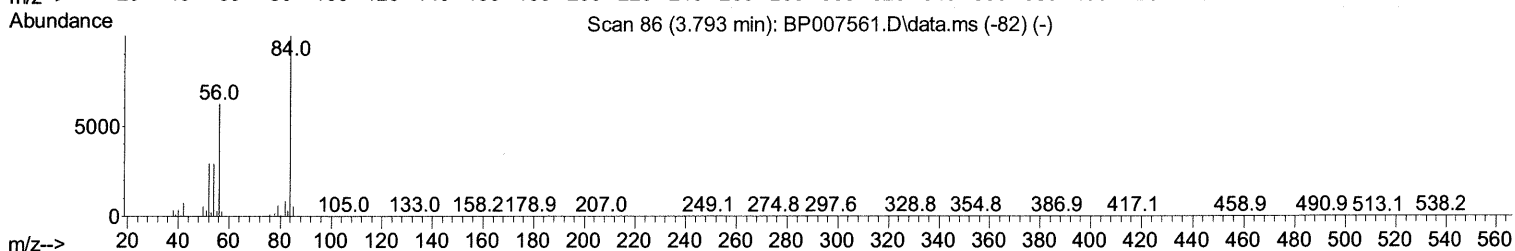
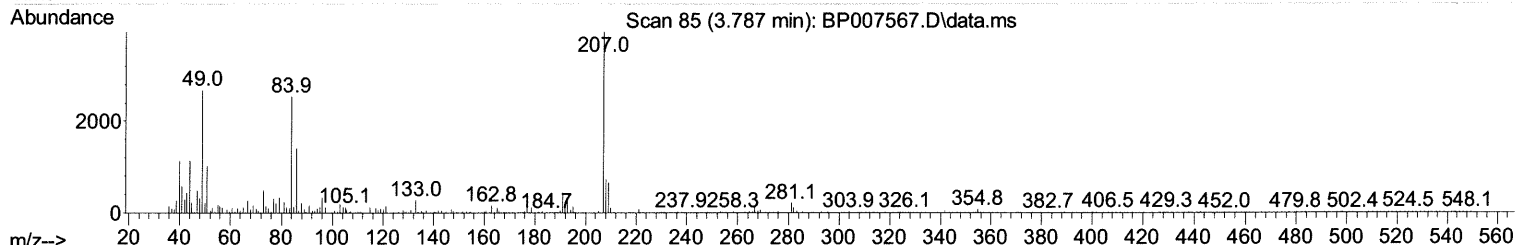
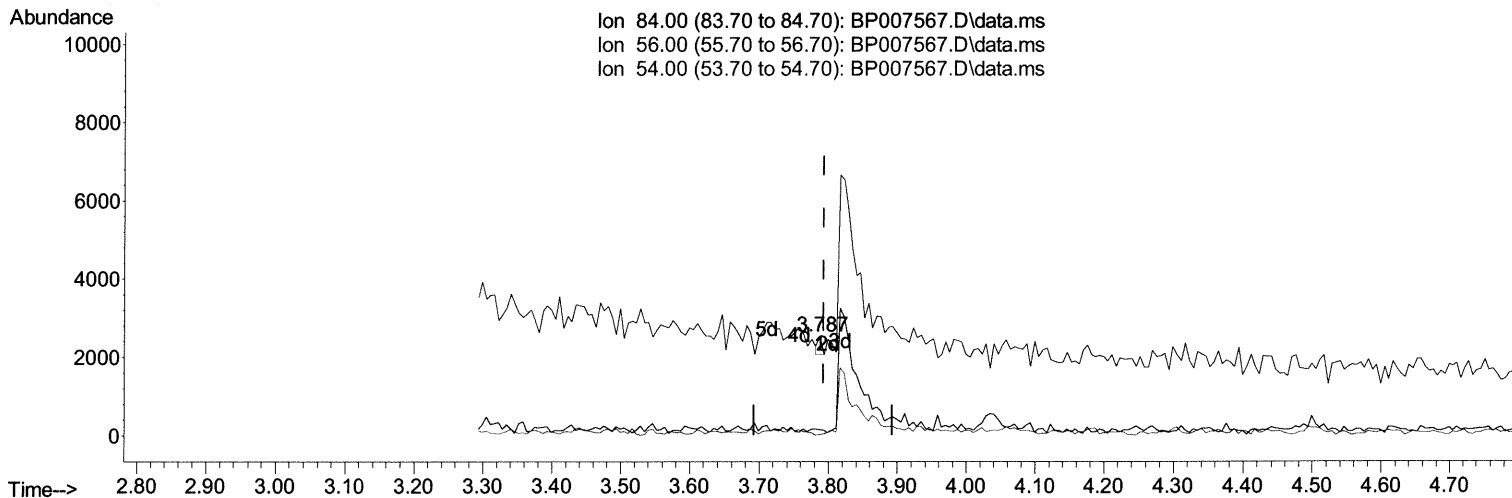
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TIC: BP007567.D\data.ms

(4) Pyridine-d5 (S)

3.787min (-0.006) 0.01 ng/ul

response 170

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	5.67#
54.00	31.70	1.02#
0.00	0.00	0.00

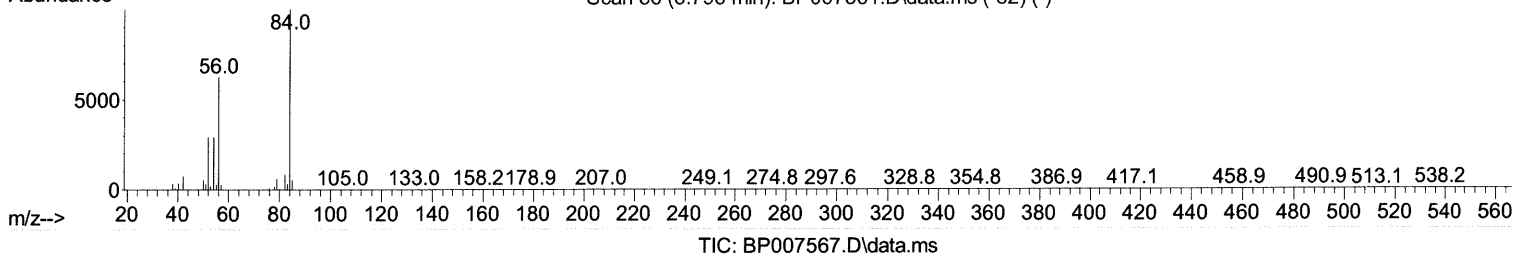
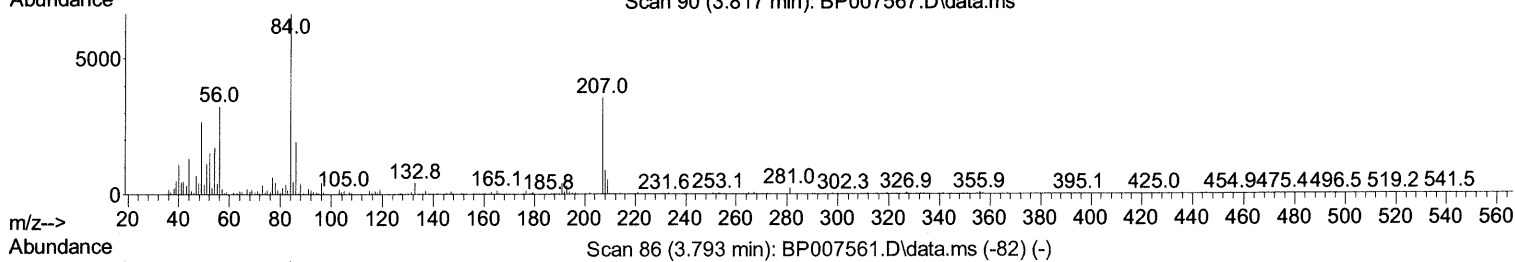
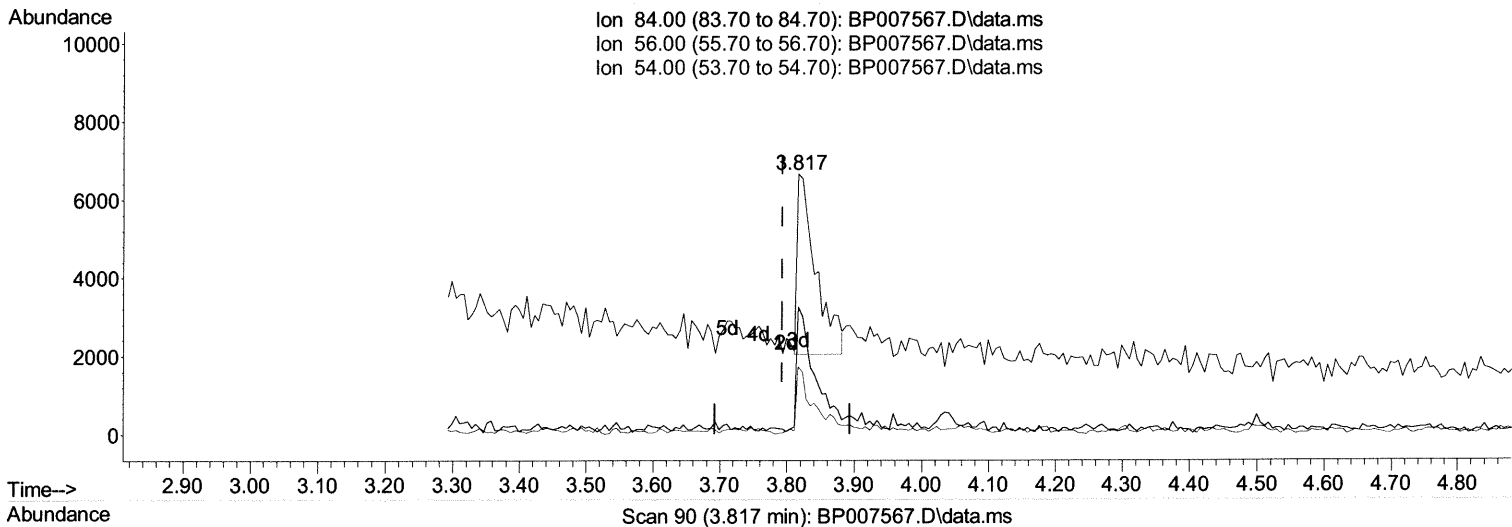
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Manual Integrations
APPROVED

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(4) Pyridine-d5 (S)

3.817min (+ 0.024) 0.57 ng/ul m

10/26/21 JU

response 8848

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	48.97#
54.00	31.70	26.02
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : M4217-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGB5

Manual Integrations
 APPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	225695	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	930747	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	611474	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	1301759	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	1337510	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	1368678	20.000	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.375	96	6687	1.170	ng/ul	0.00
4) Pyridine-d5	3.817	84	8848m	0.566	ng/ul	> 0.0210/26/21JU
7) Phenol-d5	7.105	99	120084	6.813	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	73577	6.999	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	97854	7.057	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	94827	6.985	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	44818	7.609	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	41410	7.527	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	90083	6.713	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	117232	6.460	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	288152	6.811	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	340511	6.405	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	25937	3.821	ng/ul	0.00
60) Fluorene-d10	15.575	176	261582	7.389	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	13215	2.534	ng/ul	0.00
73) Anthracene-d10	17.433	188	402360	7.278	ng/ul	0.00
81) Pyrene-d10	19.663	212	503382	7.528	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	463996	6.813	ng/ul	0.00

Target Compounds Qvalue

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