

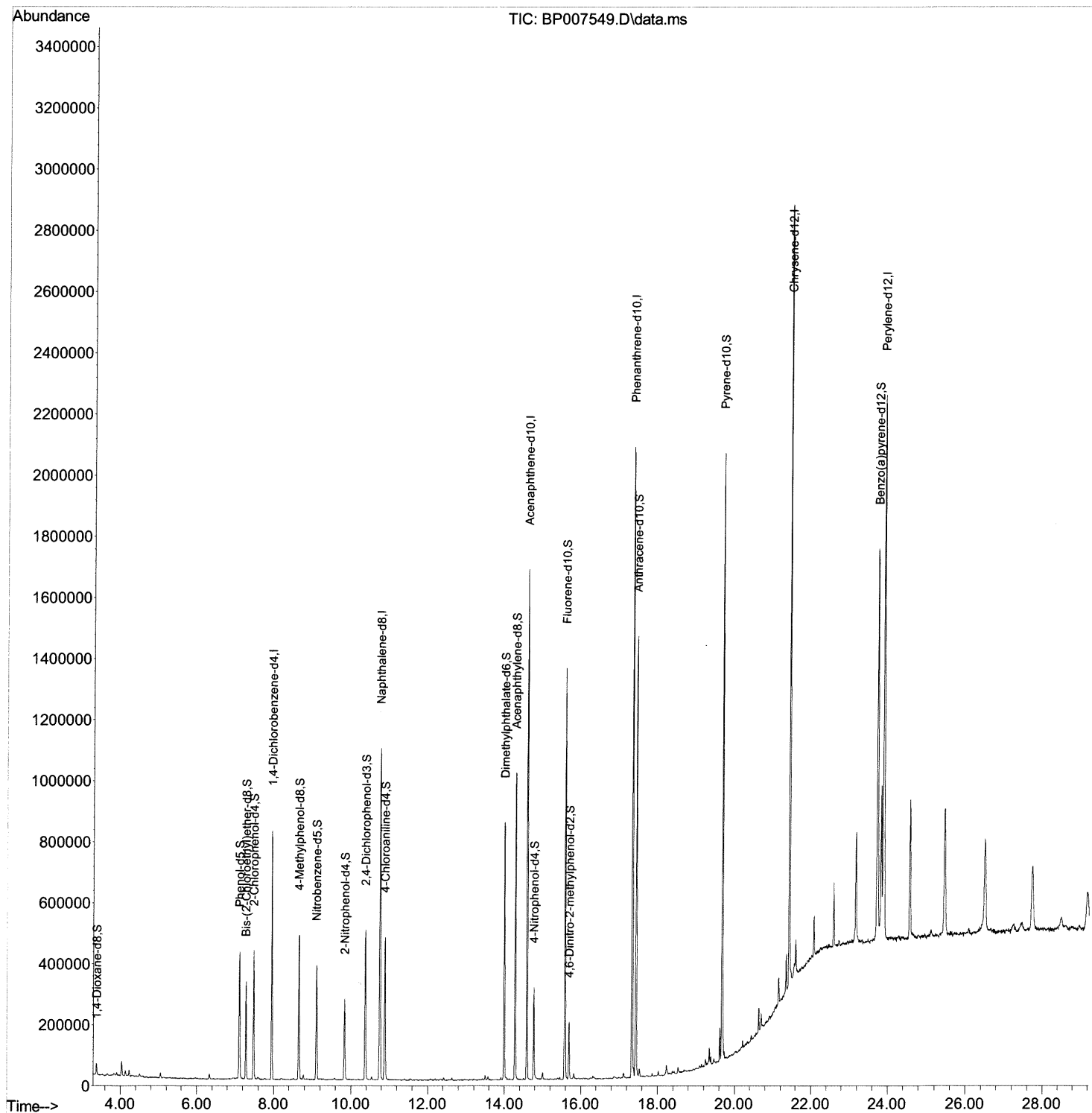
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007549.D
Acq On : 21 Oct 2021 11:47
Operator : CG/JU
Sample : M4182-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B38

Manual Integrations
APPROVED

mohammad
10/22/2021 2:54:47 PM

Quant Time: Oct 21 12:12:53 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)

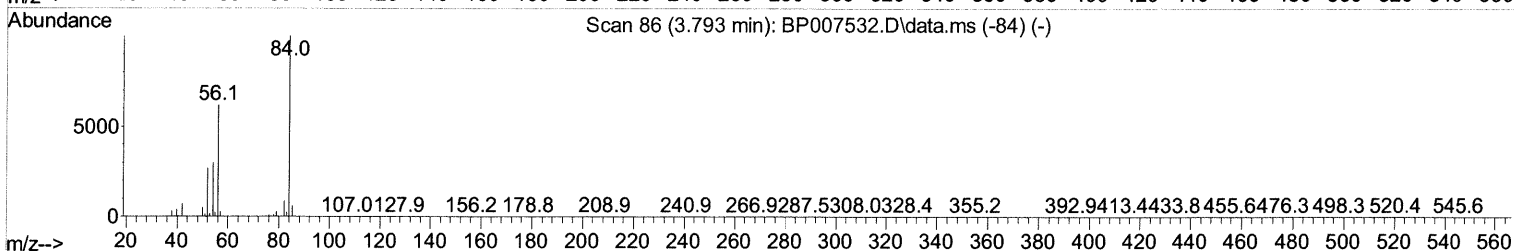
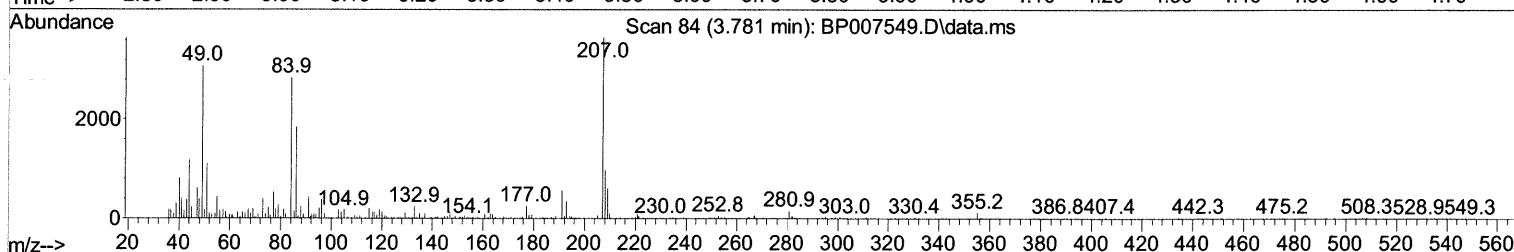
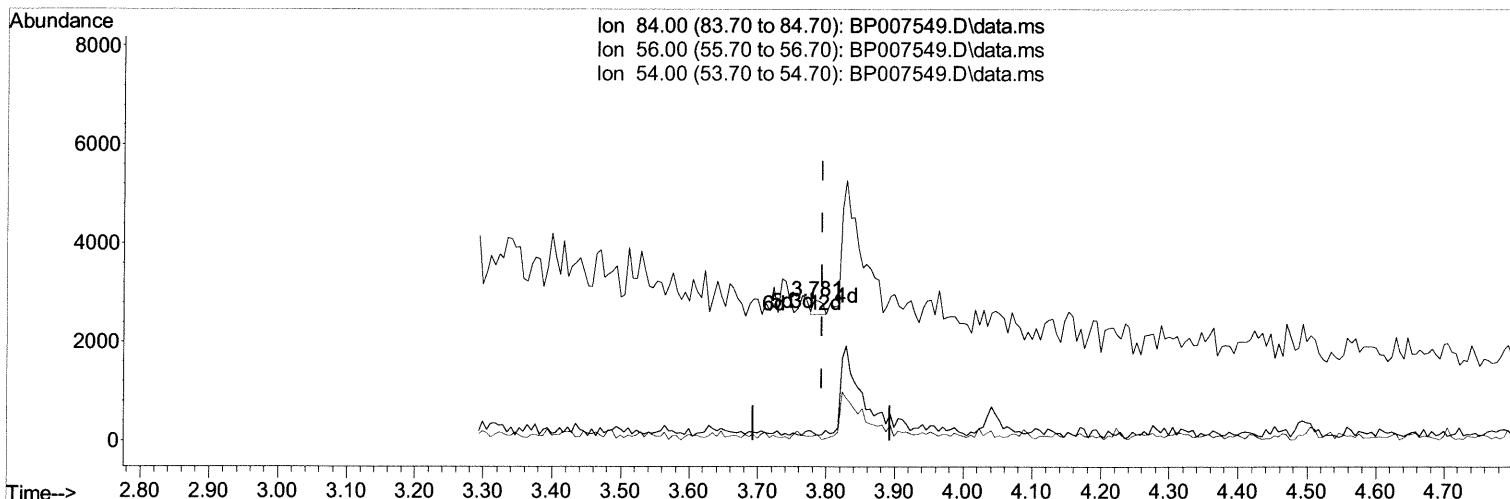
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TIC: BP007549.D\data.ms

(4) Pyridine-d5 (S)

3.781min (-0.012) 0.02 ng/ul

response 274

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	5.49#
54.00	31.70	3.38#
0.00	0.00	0.00

Quantitation Report (Qedit)

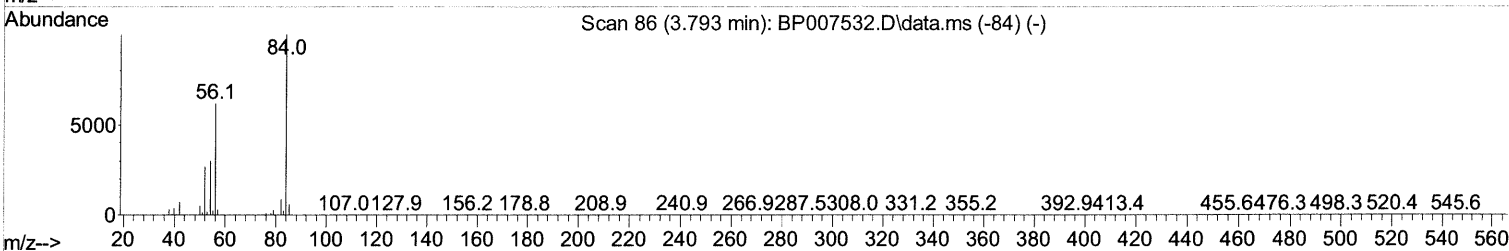
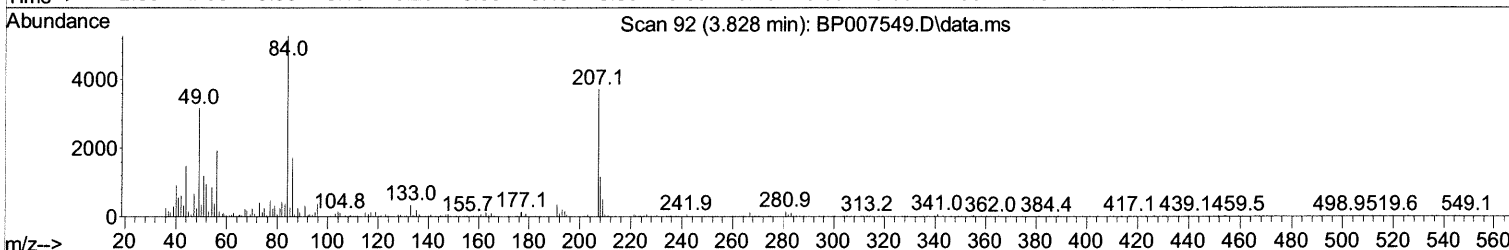
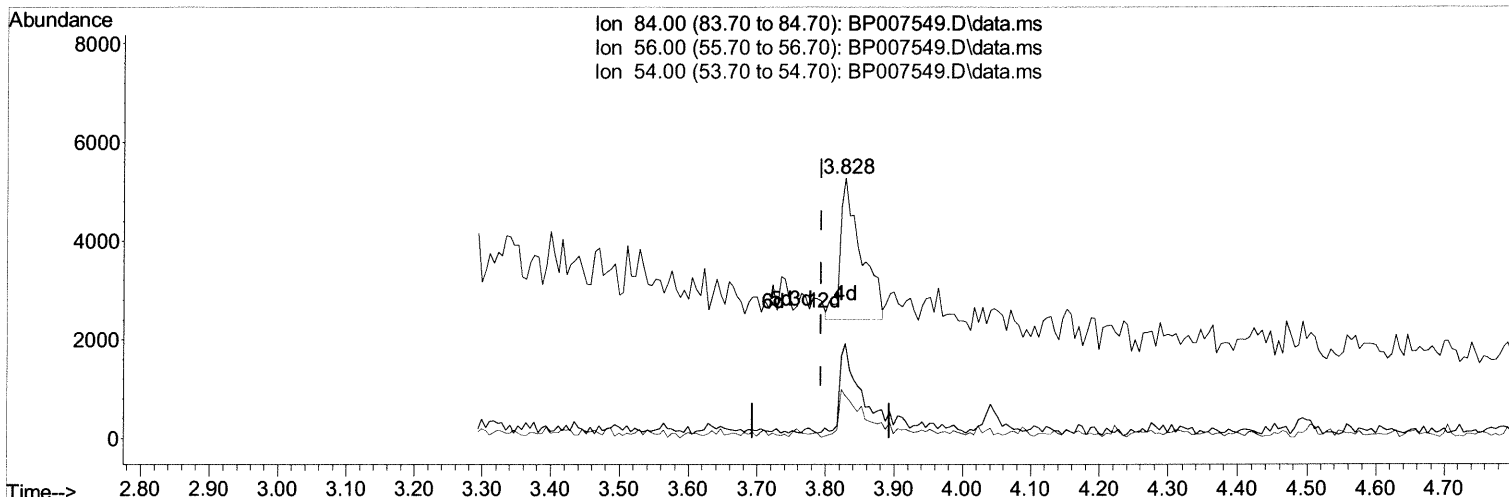
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Manual Integrations
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TIC: BP007549.D\data.ms

(4) Pyridine-d5 (S)

3.828min (+ 0.035) 0.40 ng/ul m 10/26/21JU

response 6027

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	36.47#
54.00	31.70	16.38#
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : M4182-01
 Misc :
 ALS Vial : 7 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	218264	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	886176	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	578956	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1249401	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	1315914	20.000	ng/ul	0.00
88) Perylene-d12	23.874	264	1359178	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	21324	3.859	ng/ul	0.00
4) Pyridine-d5	3.828	84	6027m >	0.398	ng/ul >	0.04 10/26/21 JU
7) Phenol-d5	7.111	99	240995	14.139	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	140487	13.819	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	190589	14.213	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	181014	13.787	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	86836	15.484	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	85542	16.330	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	184280	14.423	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	247793	14.342	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	584720	14.598	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	707628	14.058	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	69640	10.835	ng/ul	0.00
60) Fluorene-d10	15.581	176	525058	15.664	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	43067	8.603	ng/ul	0.00
73) Anthracene-d10	17.439	188	834265	15.723	ng/ul	0.00
81) Pyrene-d10	19.669	212	1027117	15.613	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	985149	14.567	ng/ul	0.00

Target Compounds Qvalue

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