

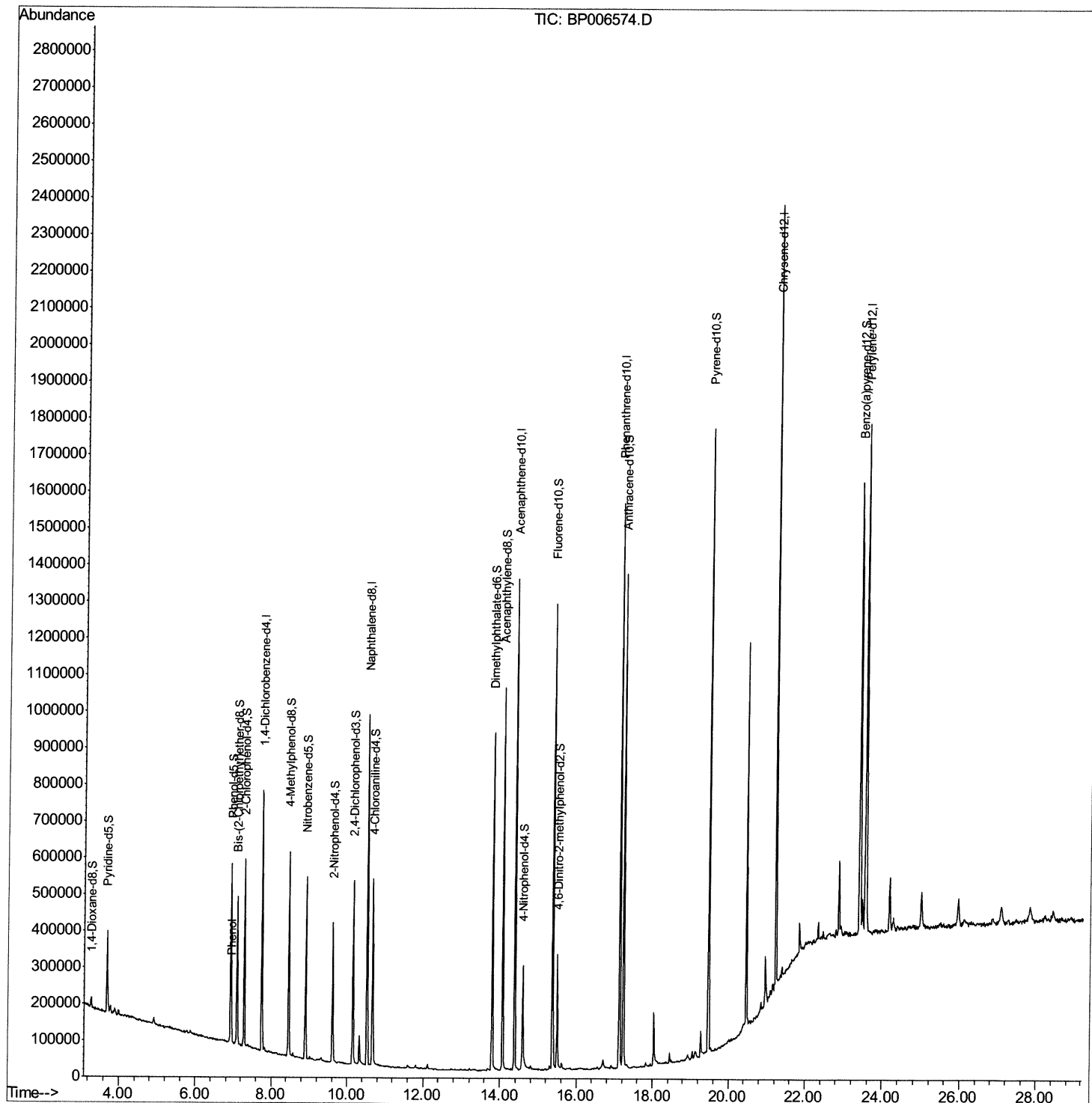
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP080721\
Data File : BP006574.D
Acq On : 07 Aug 2021 18:23
Operator : CG/JU
Sample : M3163-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGEZ0

Manual Integrations
APPROVED

mohammad
8/9/2021 12:16:45 PM

Quant Time: Aug 08 02:53:52 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 07 13:59:25 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

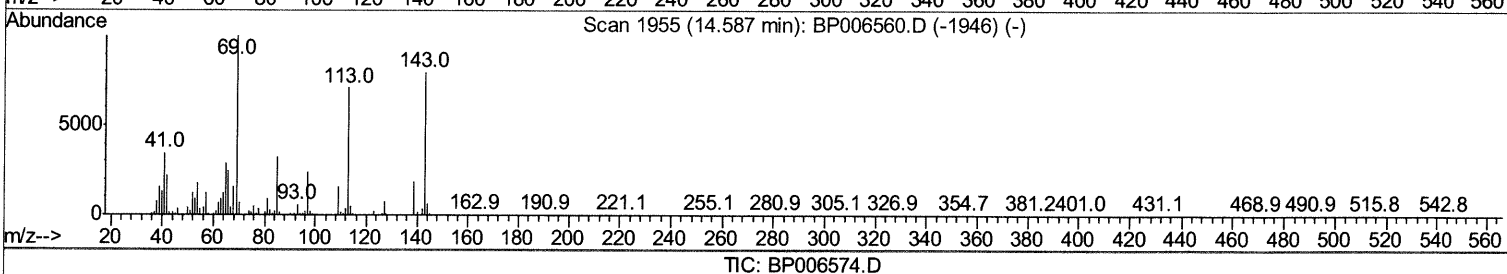
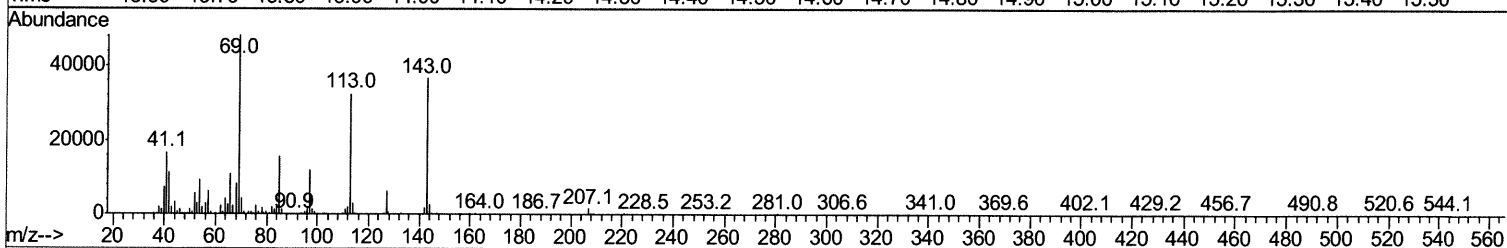
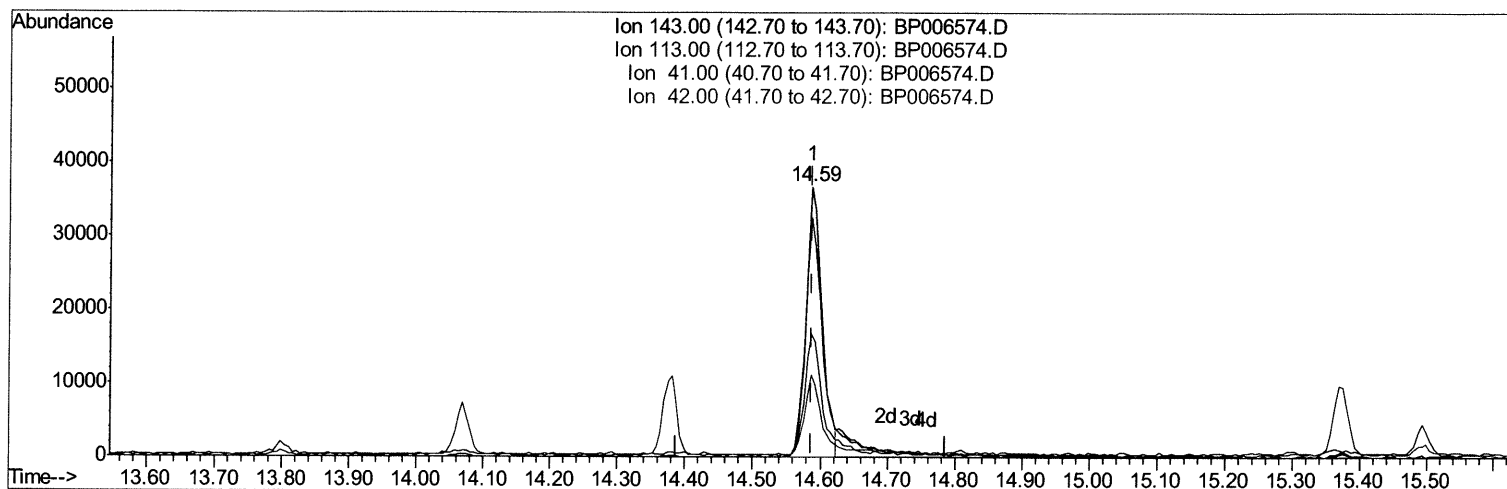
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(54) 4-Nitrophenol-d4 (S)

14.587min (-0.000) 10.64ng/ul

response 61062

Ion	Exp%	Act%
143.00	100	100
113.00	89.10	88.54
41.00	43.50	45.32
42.00	27.90	30.29

Quantitation Report (Qedit)

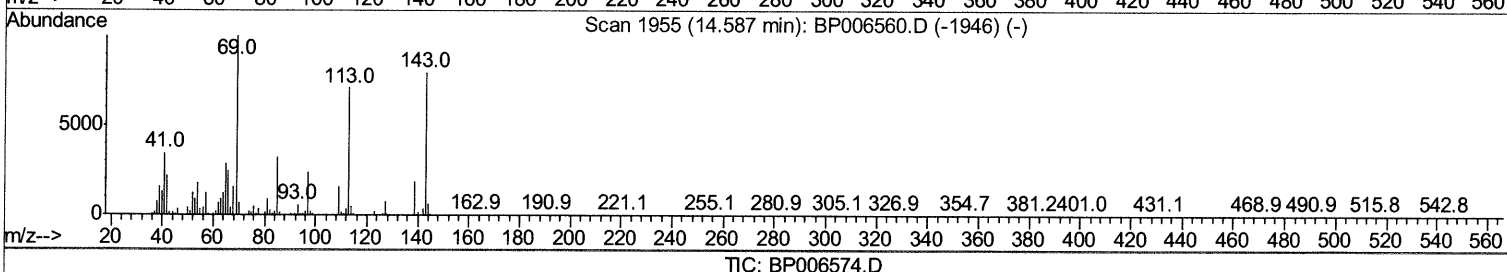
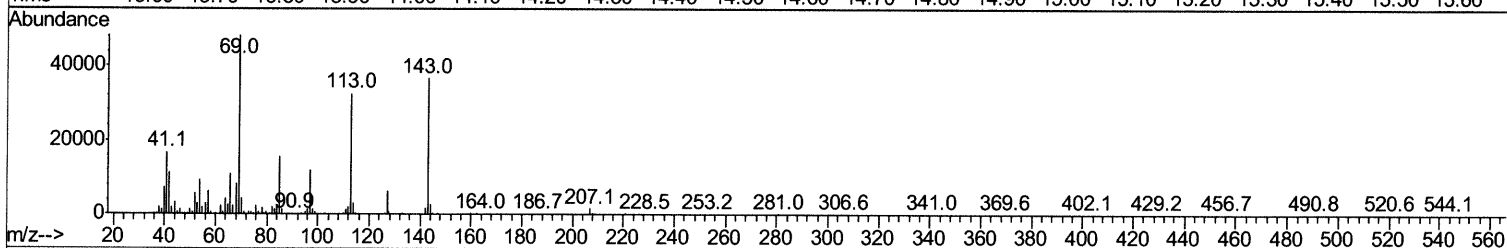
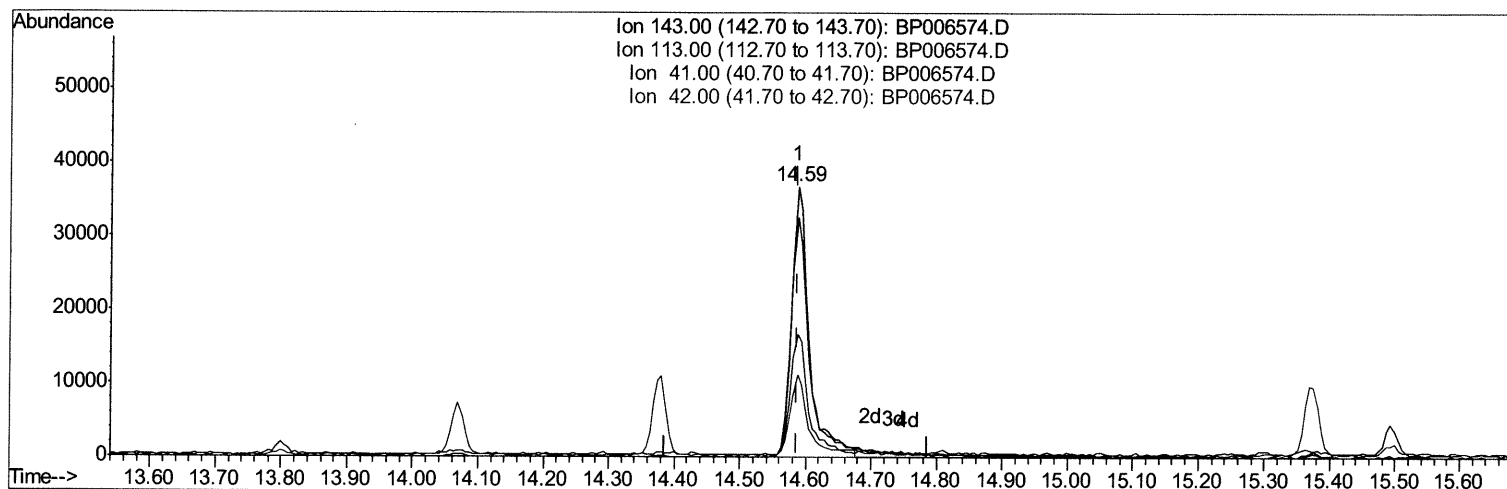
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Instrument :
 BNA_P
 ClientSampled :
 BGEZ0

Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Aug 07 13:59:25 2021
 Response via : Initial Calibration



(54) 4-Nitrophenol-d4 (S)

14.587min (-0.000) 11.89ng/ul m

response 68268

Ion	Exp%	Act%
143.00	100	100
113.00	89.10	88.54
41.00	43.50	45.32
42.00	27.90	30.29

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
 Data File : BP006574.D
 Acq On : 07 Aug 2021 18:23
 Operator : CG/JU
 Sample : M3163-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGEZ0

Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	174913	20.00	ng/ul	0.00
20) Naphthalene-d8	10.53	136	732862	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.38	164	406246	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.13	188	807081	20.00	ng/ul	0.00
79) Chrysene-d12	21.23	240	838218	20.00	ng/ul	0.00
88) Perylene-d12	23.55	264	901957	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	17680	3.77	ng/uL	0.00
4) Pyridine-d5	3.68	84	137662	9.90	ng/ul	0.00
7) Phenol-d5	6.93	99	268648	16.35	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.09	67	185155	18.11	ng/ul	0.00
11) 2-Chlorophenol-d4	7.28	132	211371	17.32	ng/ul	0.00
15) 4-Methylphenol-d8	8.46	113	192723	14.92	ng/ul	0.00
21) Nitrobenzene-d5	8.90	128	104393	19.13	ng/ul	0.00
24) 2-Nitrophenol-d4	9.62	143	101657	19.51	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.16	165	158227	15.50	ng/ul	0.00
31) 4-Chloroaniline-d4	10.68	131	255298	15.19	ng/ul	0.00
46) Dimethylphthalate-d6	13.80	166	537283	18.58	ng/ul	0.00
49) Acenaphthylene-d8	14.07	160	647813	18.24	ng/ul	0.00
54) 4-Nitrophenol-d4	14.59	143	68268m	11.89	ng/ul	0.00
60) Fluorene-d10	15.37	176	440582	18.13	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.49	200	48866	10.86	ng/ul	0.00
73) Anthracene-d10	17.23	188	681541	18.71	ng/ul	0.00
81) Pyrene-d10	19.47	212	792193	18.37	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.40	264	802431	17.51	ng/ul	-0.01

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
8) Phenol	6.95	94	21083	1.256	ng/ul	89

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