

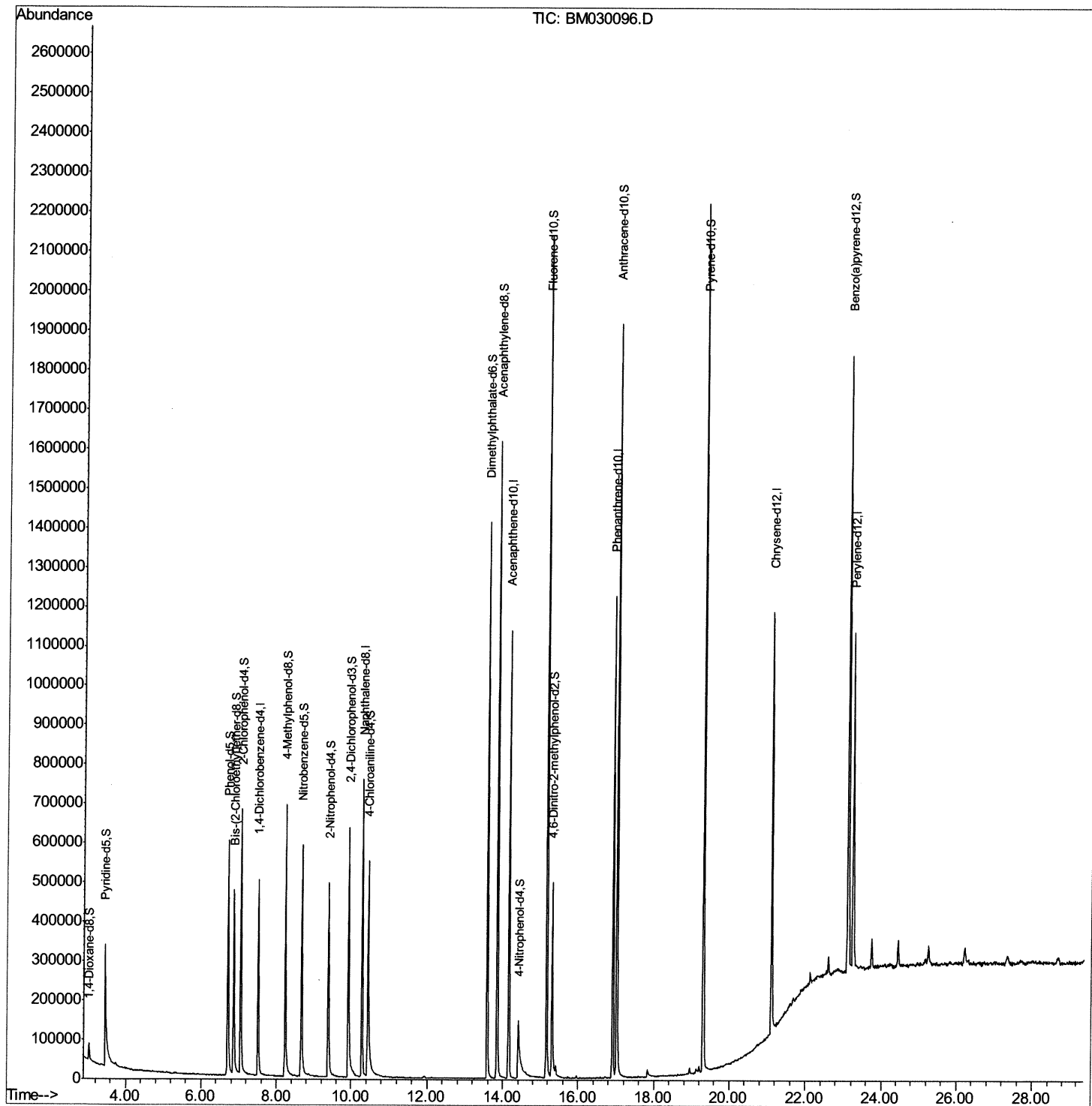
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052021\
Data File : BM030096.D
Acq On : 20 May 2021 19:07
Operator : CG/JU
Sample : PB136551BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK551

Manual Integrations
APPROVED

mohammad
5/21/2021 3:38:04 PM

Quant Time: May 21 01:10:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 20 12:00:20 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

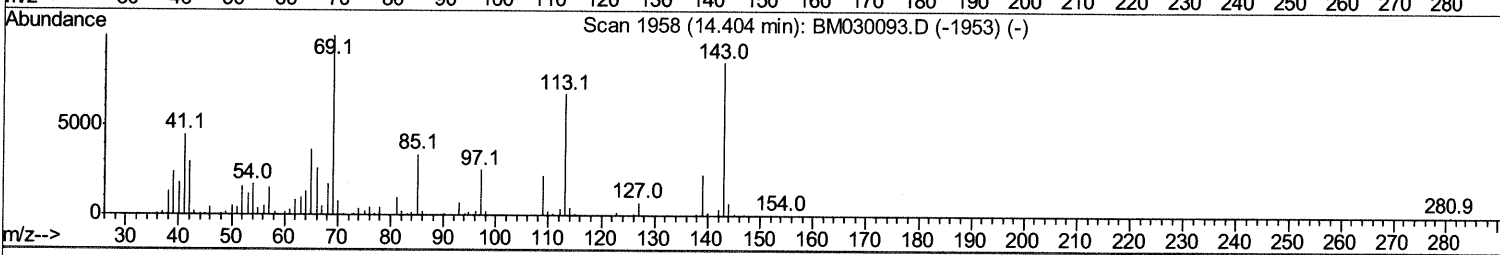
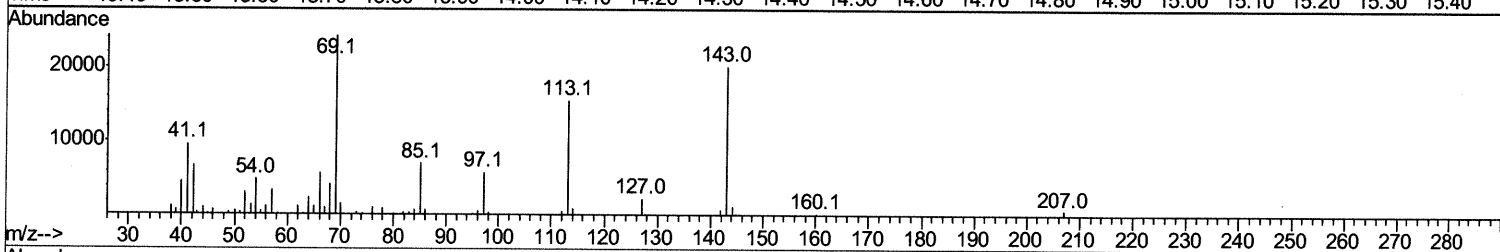
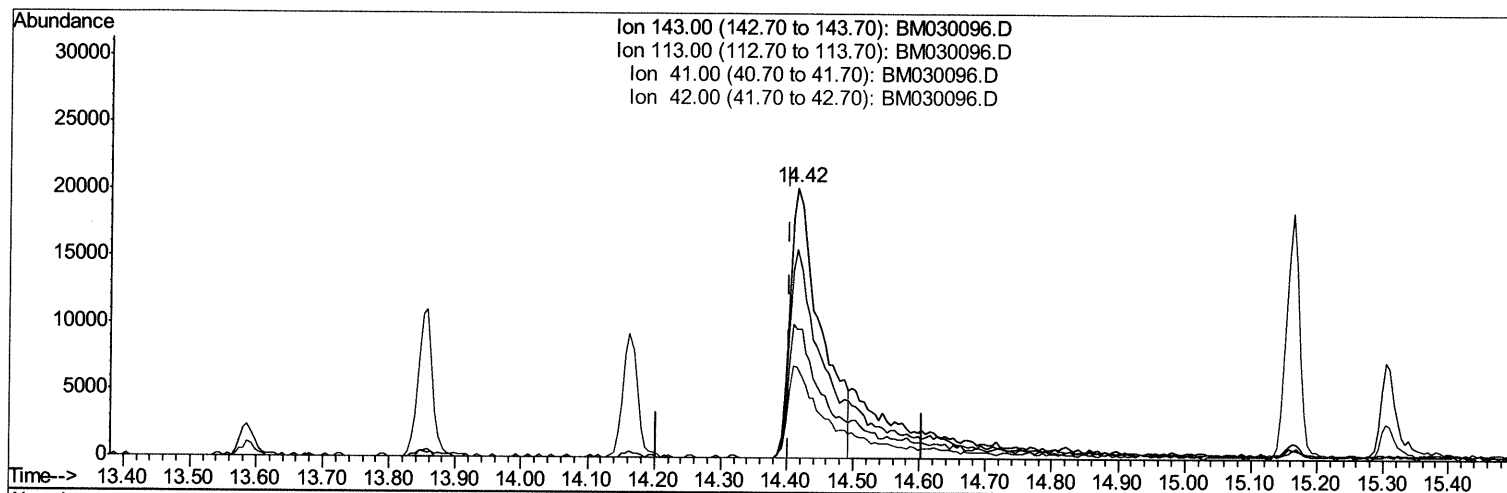
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TIC: BM030096.D

(54) 4-Nitrophenol-d4 (S)

14.416min (+0.012) 13.63ng/ul

response 64562

Ion	Exp%	Act%
143.00	100	100
113.00	71.10	77.27
41.00	44.70	47.82
42.00	28.30	33.77

Quantitation Report (Qedit)

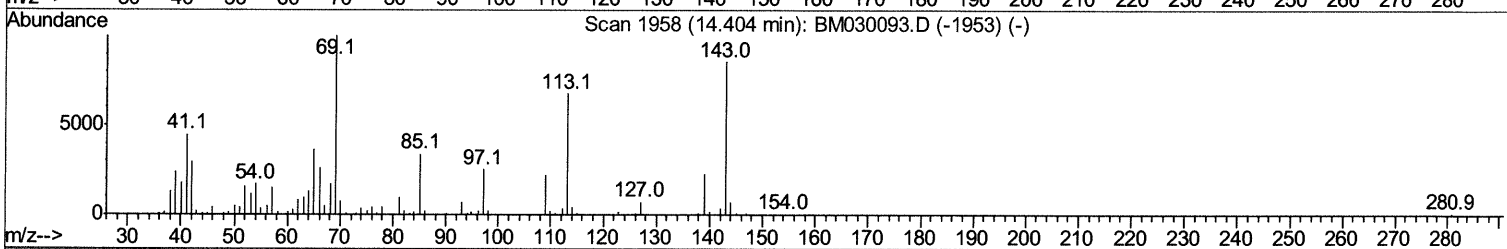
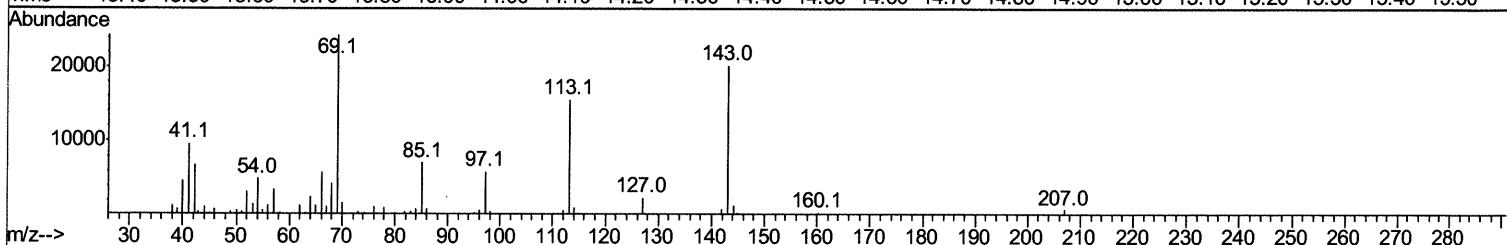
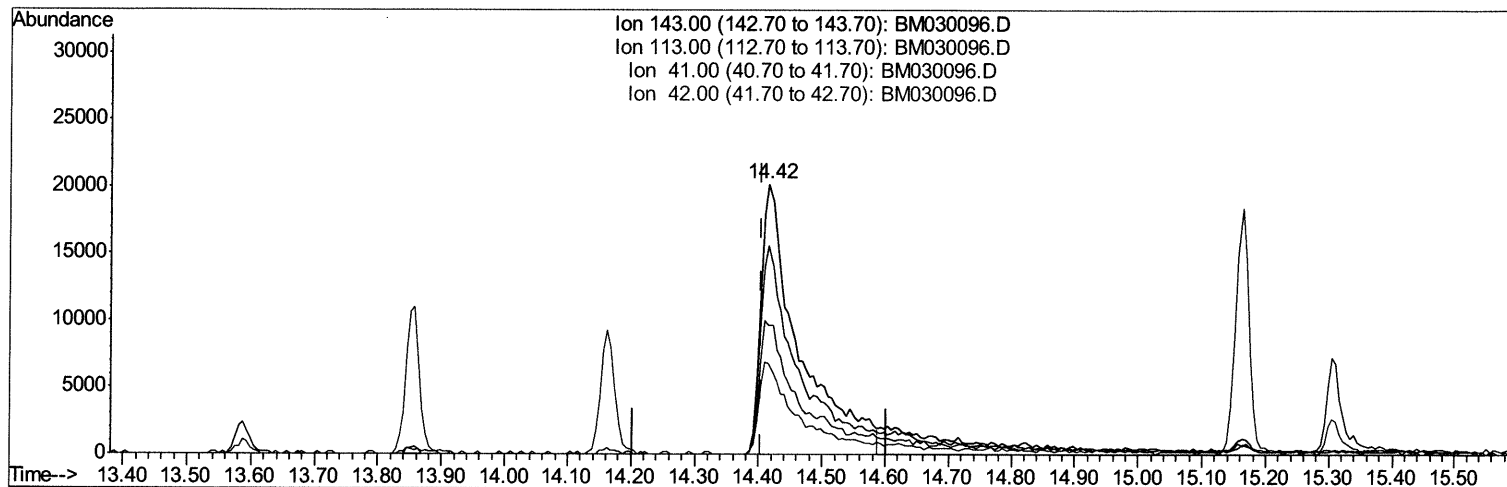
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Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



TIC: BM030096.D

(54) 4-Nitrophenol-d4 (S)

14.416min (+0.012) 17.58ng/ul m 05/24/21 JU

response 83247

Ion	Exp%	Act%
143.00	100	100
113.00	71.10	77.27
41.00	44.70	47.82
42.00	28.30	33.77

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM052021\
 Data File : BM030096.D
 Acq On : 20 May 2021 19:07
 Operator : CG/JU
 Sample : PB136551BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK551

Manual Integrations
 APPROVED

mohammad
 5/21/2021 3:38:04 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	142270	20.00	ng/ul	0.00
20) Naphthalene-d8	10.29	136	593414	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.16	164	375121	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.92	188	696053	20.00	ng/ul	0.00
79) Chrysene-d12	21.10	240	567520	20.00	ng/ul	0.00
88) Perylene-d12	23.24	264	593410	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	24288	6.82	ng/uL	0.00
4) Pyridine-d5	3.45	84	230682	25.86	ng/ul	0.00
7) Phenol-d5	6.70	99	351714	26.58	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.86	67	239776	29.23	ng/ul	0.00
11) 2-Chlorophenol-d4	7.05	132	310698	30.43	ng/ul	0.00
15) 4-Methylphenol-d8	8.23	113	284892	26.04	ng/ul	0.00
21) Nitrobenzene-d5	8.67	128	142373	34.82	ng/ul	0.00
24) 2-Nitrophenol-d4	9.39	143	157195	34.06	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.92	165	252374	27.27	ng/ul	0.00
31) 4-Chloroaniline-d4	10.44	131	385765	27.39	ng/ul	0.00
46) Dimethylphthalate-d6	13.59	166	918827	31.60	ng/ul	0.00
49) Acenaphthylene-d8	13.86	160	1140210	30.94	ng/ul	0.00
54) 4-Nitrophenol-d4	14.42	143	83247m	17.58	ng/ul>	0.0105724/21JU
60) Fluorene-d10	15.16	176	786769	31.87	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.31	200	115382	25.52	ng/ul	0.00
73) Anthracene-d10	17.01	188	1115542	31.83	ng/ul	0.00
81) Pyrene-d10	19.31	212	1116731	38.36	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.11	264	1049806	31.21	ng/ul	0.00

Target Compounds Ovalue

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