

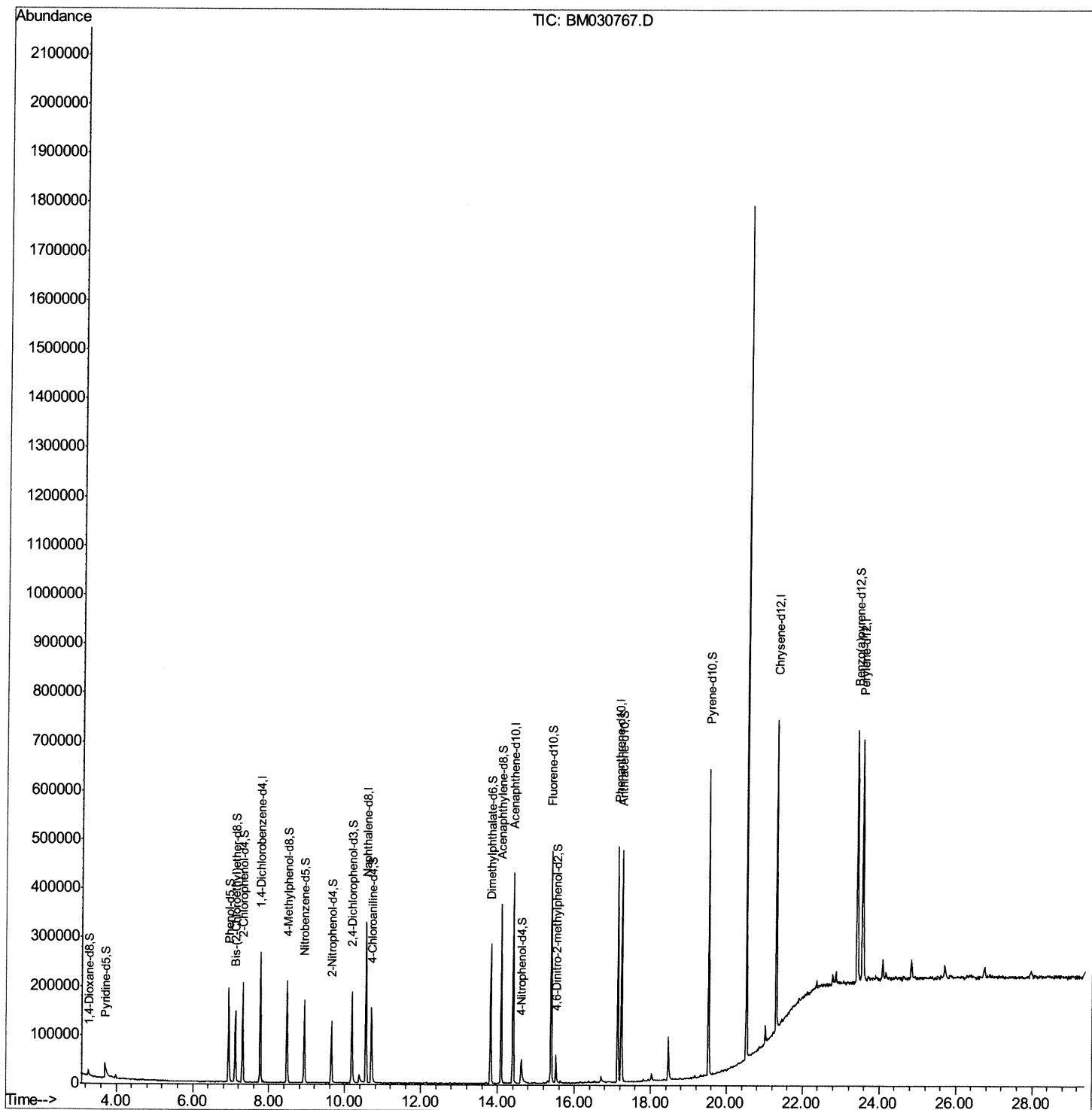
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM063021\
Data File : BM030767.D
Acq On : 02 Jul 2021 06:16
Operator : CG/JU
Sample : M2825-21
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGDS0

Manual Integrations
APPROVED

mohammad
7/2/2021 12:04:22 PM

Quant Time: Jul 02 07:35:11 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 02 04:13:39 2021
Response via : Initial Calibration



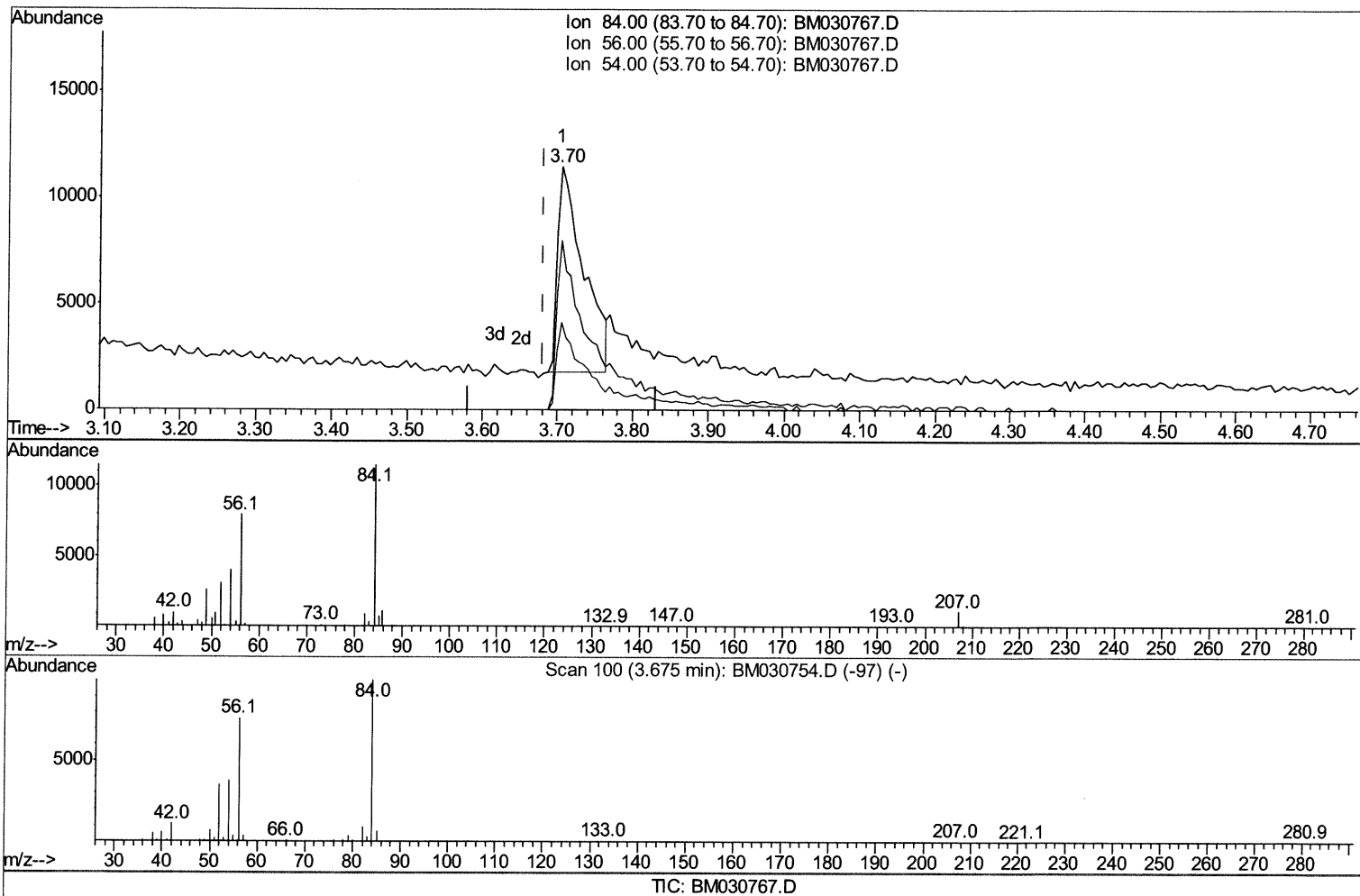
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(4) Pyridine-d5 (S)

3.705min (+0.023) 5.02ng/ul

response 23317

Ion	Exp%	Act%
84.00	100	100
56.00	73.70	69.44
54.00	35.80	36.03
0.00	0.00	0.00

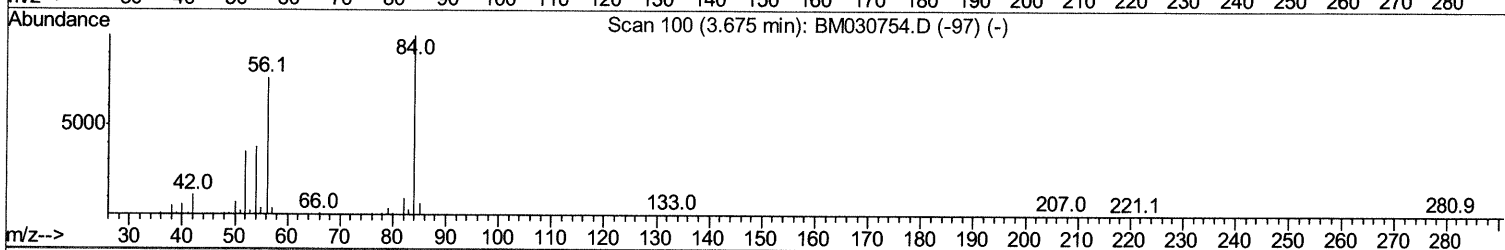
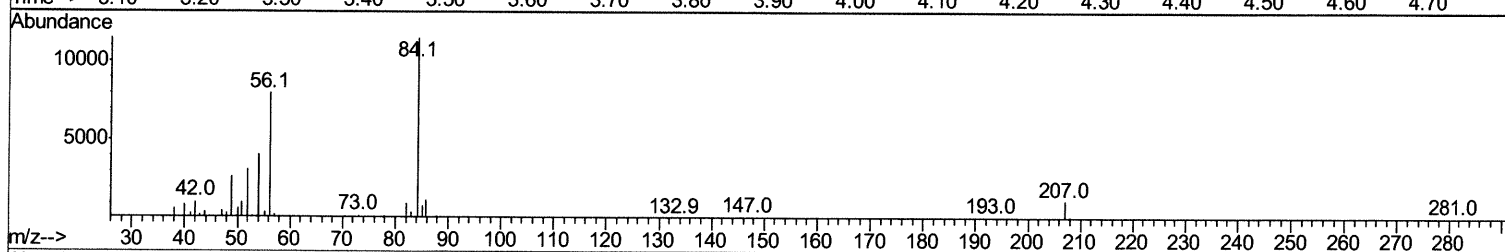
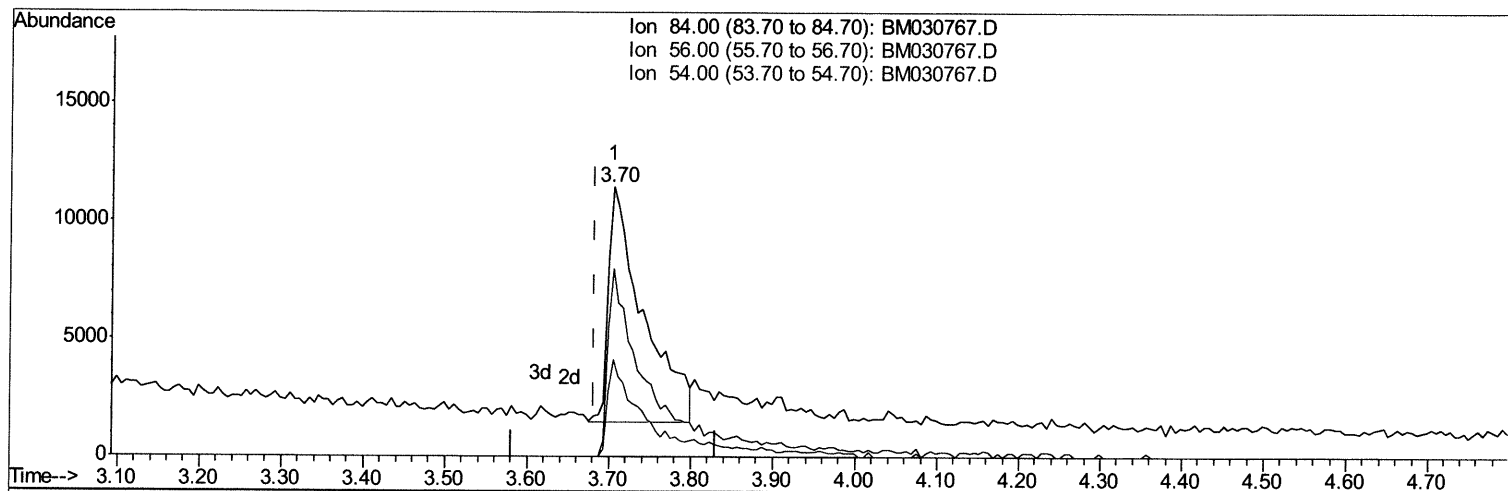
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TIC: BM030767.D

(4) Pyridine-d5 (S)

3.705min (+0.023) 6.33ng/ul m

response 29407

Ion	Exp%	Act%
84.00	100	100
56.00	73.70	69.44
54.00	35.80	36.03
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	70092	20.00	ng/ul	0.00
20) Naphthalene-d8	10.56	136	267008	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.40	164	145983	20.00	ng/ul	-0.01
64) Phenanthrene-d10	17.14	188	281383	20.00	ng/ul	0.00
79) Chrysene-d12	21.31	240	295053	20.00	ng/ul	0.00
88) Perylene-d12	23.56	264	345214	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	6537	3.78	ng/uL	0.00
4) Pyridine-d5	3.70	84	29407m	6.33	ng/ul	0.02
7) Phenol-d5	6.94	99	101311	16.76	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.10	67	69385	18.26	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.30	132	84467	18.10	ng/ul	0.00
15) 4-Methylphenol-d8	8.47	113	78437	16.67	ng/ul	0.00
21) Nitrobenzene-d5	8.93	128	36599	18.37	ng/ul	0.00
24) 2-Nitrophenol-d4	9.65	143	36746	16.60	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.18	165	70884	16.71	ng/ul	0.00
31) 4-Chloroaniline-d4	10.70	131	91884	15.53	ng/ul	0.00
46) Dimethylphthalate-d6	13.82	166	199506	19.79	ng/ul	0.00
49) Acenaphthylene-d8	14.09	160	256458	19.60	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.62	143	19274	10.09	ng/ul	0.00
60) Fluorene-d10	15.39	176	179731	20.10	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.52	200	14565	7.91	ng/ul	0.00
73) Anthracene-d10	17.24	188	264642	20.15	ng/ul	0.00
81) Pyrene-d10	19.53	212	293648	18.99	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.41	264	364963	20.33	ng/ul	-0.02

Target Compounds

Ovalue

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