

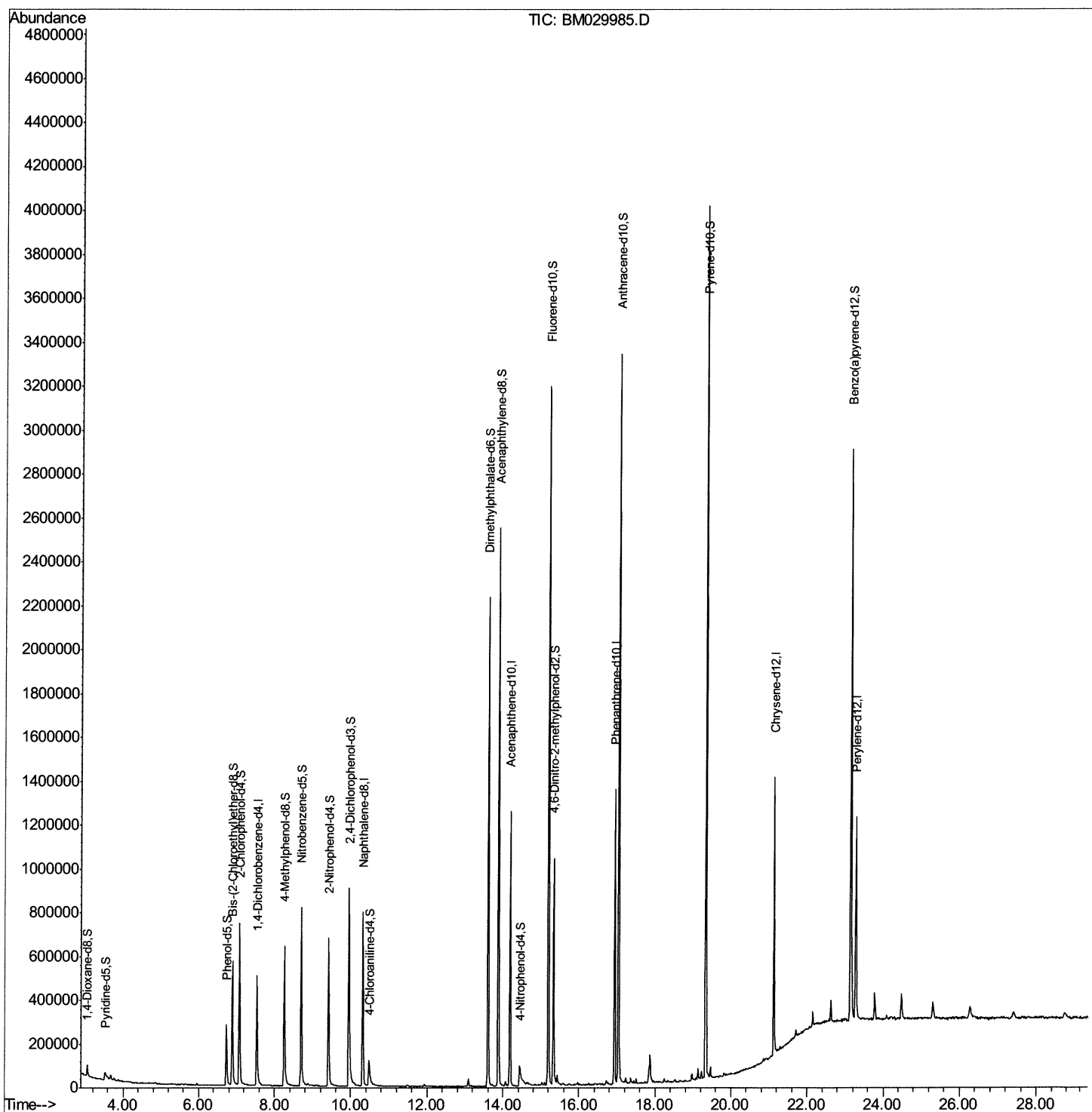
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
Data File : BM029985.D
Acq On : 12 May 2021 17:22
Operator : CG/JU
Sample : M2197-16
Misc :
ALS Vial : 84 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YARR9

Manual Integrations
APPROVED

Sohil
5/13/2021 10:30:47 AM

Quant Time: May 12 18:01:53 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 11 15:21:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

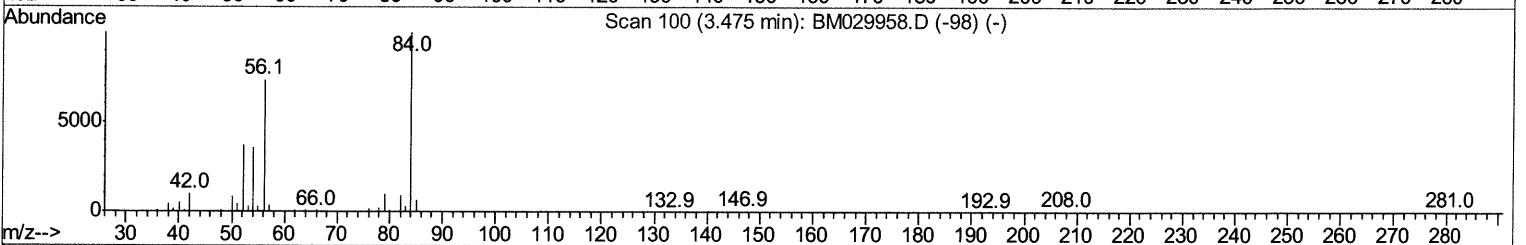
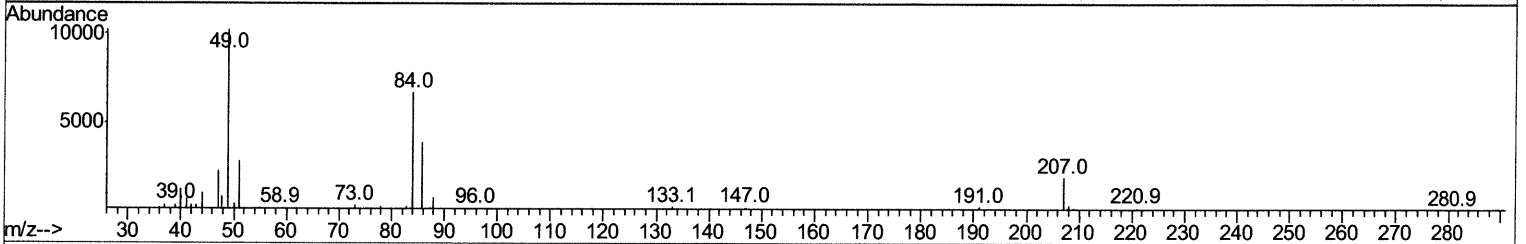
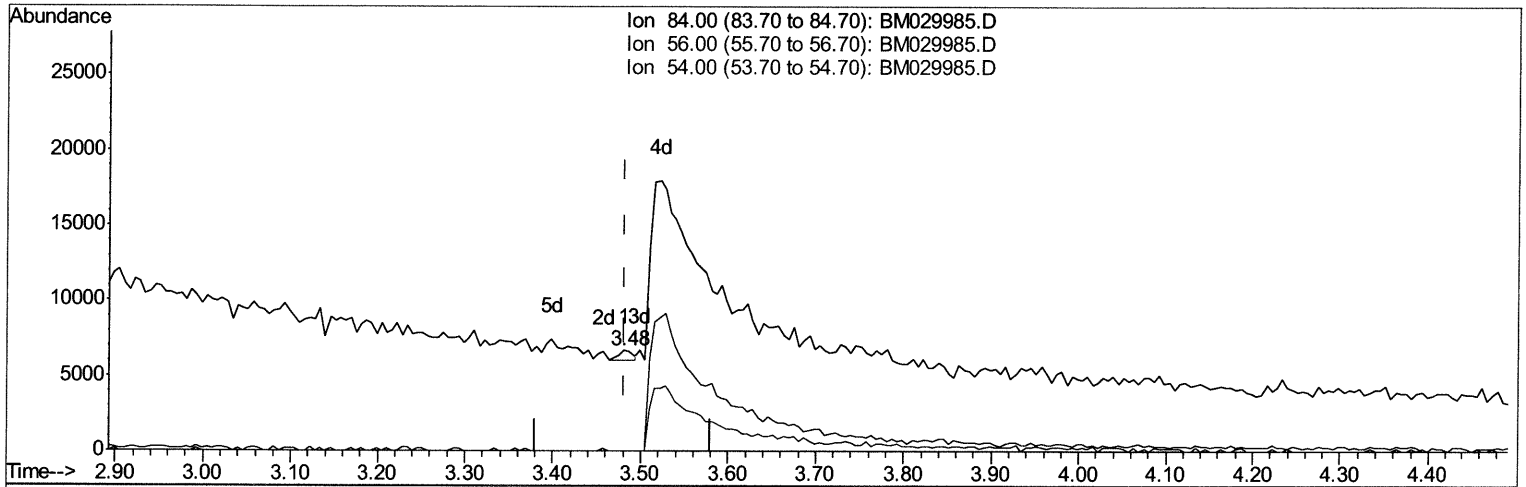
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TIC: BM029985.D

(4) Pyridine-d5 (S)
 3.481min (-0.000) 0.08ng/ul
 response 704

Ion	Exp%	Act%
84.00	100	100
56.00	67.00	0.00#
54.00	32.50	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

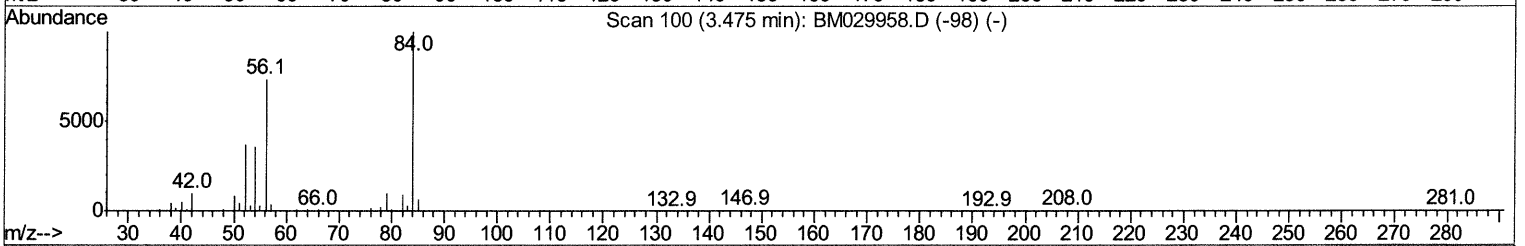
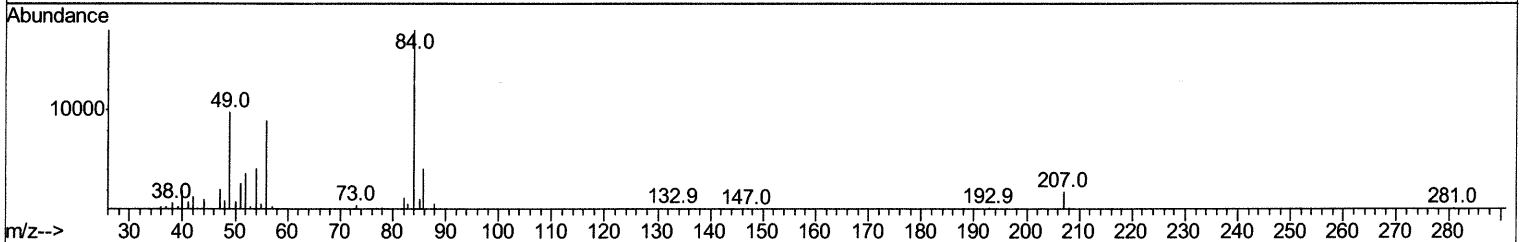
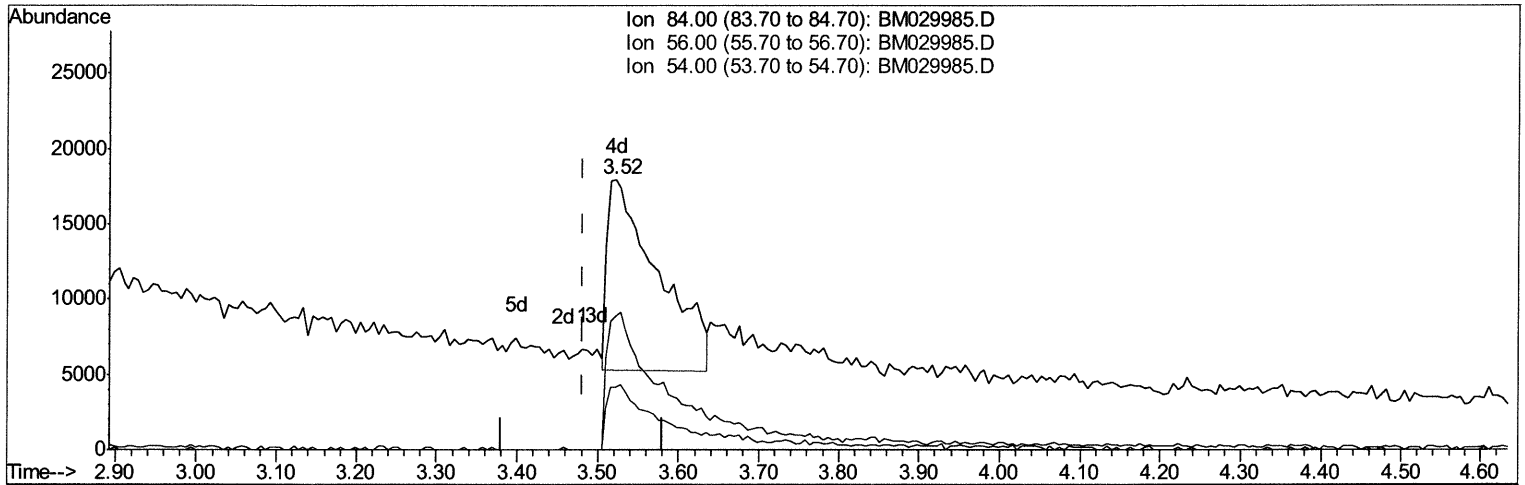
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Manual Integrations
 APPROVED

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TIC: BM029985.D

(4) Pyridine-d5 (S)

3.522min (+0.041) 6.33ng/ul m 5/13/21 JU

response 55207

Ion	Exp%	Act%
84.00	100	100
56.00	67.00	49.48#
54.00	32.50	23.27#
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
 Data File : BM029985.D
 Acq On : 12 May 2021 17:22
 Operator : CG/JU
 Sample : M2197-16
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YARR9

Manual Integrations
 APPROVED

Sohil
 5/13/2021 10:30:47 AM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	139009	20.00	ng/ul	-0.01
20) Naphthalene-d8	10.32	136	632287	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.19	164	406263	20.00	ng/ul	-0.01
64) Phenanthrene-d10	16.94	188	781060	20.00	ng/ul	0.00
79) Chrysene-d12	21.13	240	618612	20.00	ng/ul	0.00
88) Perylene-d12	23.28	264	605402	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	22547	6.48	ng/uL	0.00
4) Pyridine-d5	3.52	84	55207m>	6.33	ng/ul>	0.0405/13/21JU
7) Phenol-d5	6.73	99	180495	13.96	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.89	67	302989	37.80	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.08	132	336782	33.76	ng/ul	0.00
15) 4-Methylphenol-d8	8.26	113	259981	24.32	ng/ul	-0.01
21) Nitrobenzene-d5	8.70	128	187772	43.09	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.41	143	209619	42.62	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.95	165	373412	37.87	ng/ul	0.00
31) 4-Chloroaniline-d4	10.48	131	87288	5.82	ng/ul	0.00
46) Dimethylphthalate-d6	13.61	166	1427950	45.35	ng/ul	-0.01
49) Acenaphthylene-d8	13.88	160	1677462	42.03	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.44	143	37244	7.26	ng/ul	0.02
60) Fluorene-d10	15.19	176	1204090	45.04	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	210865	41.56	ng/ul	0.00
73) Anthracene-d10	17.04	188	1871600	47.59	ng/ul	0.00
81) Pyrene-d10	19.33	212	1940889	61.16	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.14	264	1663699	48.48	ng/ul	0.00

Target Compounds Ovalue

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