

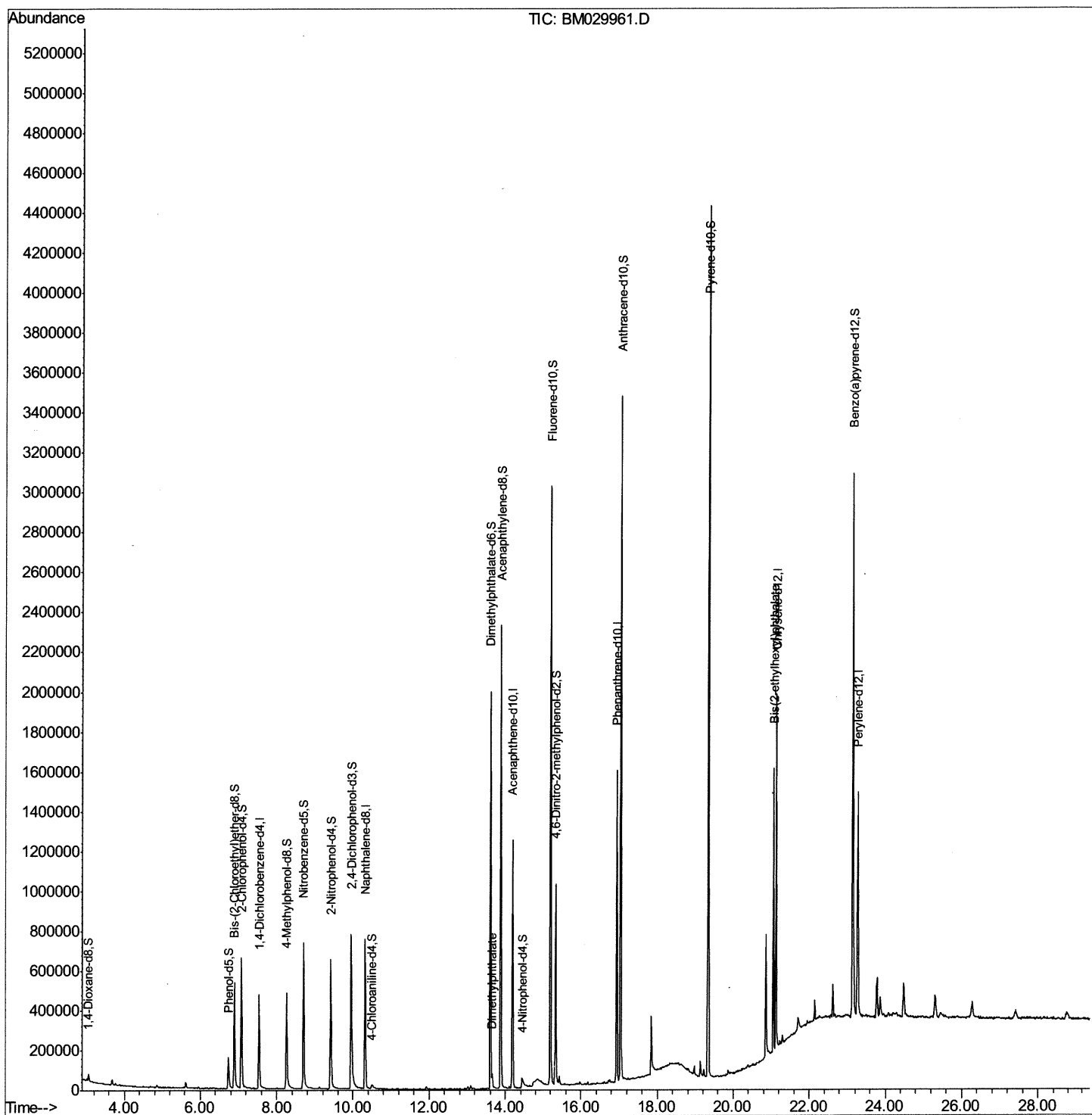
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
Data File : BM029961.D
Acq On : 12 May 2021 02:49
Operator : CG/JU
Sample : M2197-07
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YARG6

Manual Integrations
APPROVED

mohammad
5/12/2021 2:07:55 PM

Quant Time: May 12 03:32:00 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 12 03:18:52 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

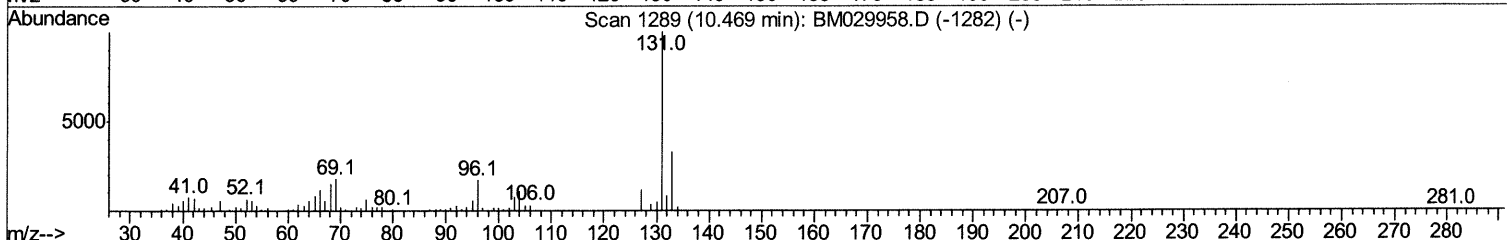
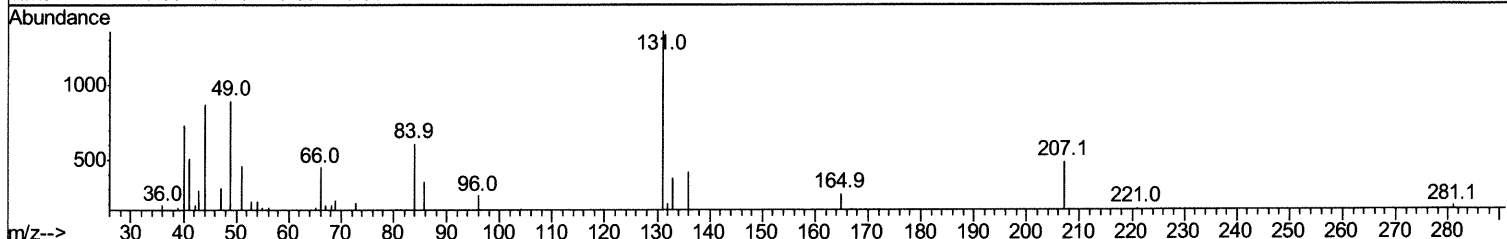
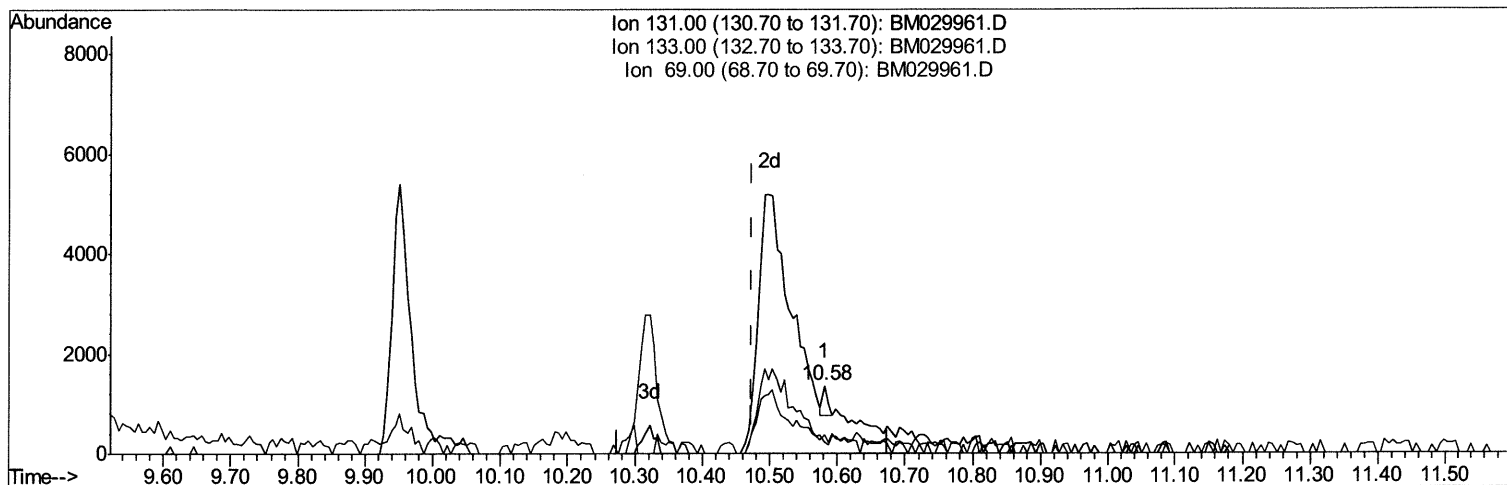
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Quant Time: Mav 12 03:29:24 2021
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TIC: BM029961.D

(31) 4-Chloroaniline-d4 (S)

10.581min (+0.106) 0.02ng/ul

response 296

Ion	Exp%	Act%
131.00	100	100
133.00	33.00	27.55
69.00	18.00	16.61
0.00	0.00	0.00

Quantitation Report (Qedit)

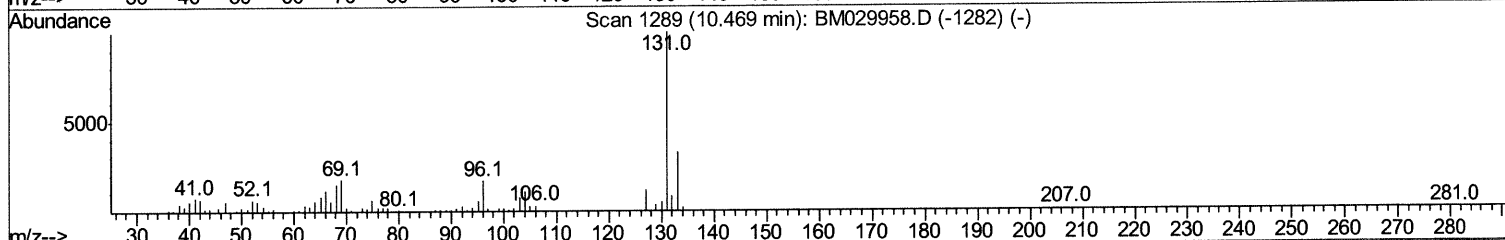
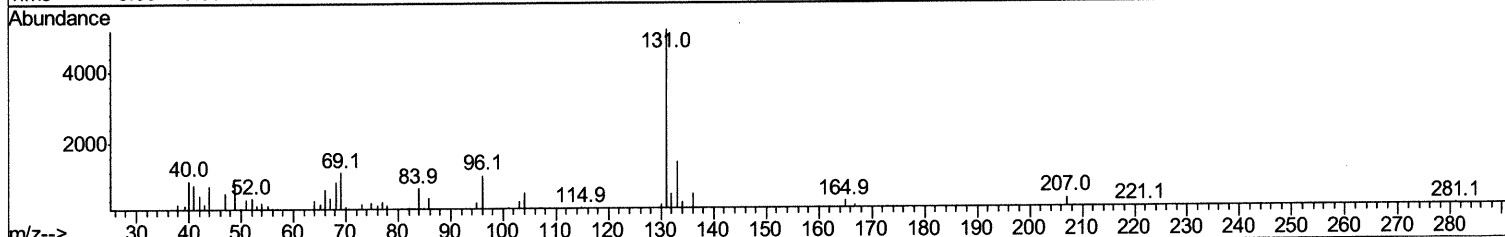
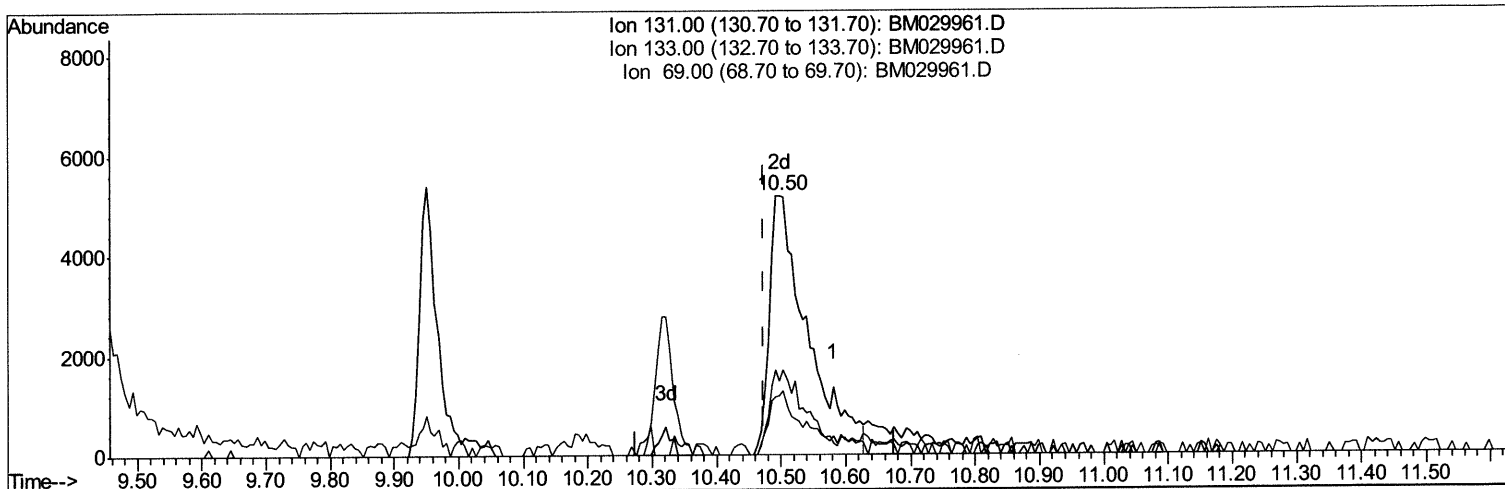
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Manual Integrations
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 Response via : Initial Calibration



TIC: BM029961.D

(31) 4-Chloroaniline-d4 (S)

10.498min (+0.024) 1.48ng/ul m 05/13/21JU

response 21201

Ion	Exp%	Act%
131.00	100	100
133.00	33.00	28.41
69.00	18.00	22.80#
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
 Data File : BM029961.D
 Acq On : 12 May 2021 02:49
 Operator : CG/JU
 Sample : M2197-07
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YARG6

Manual Integrations
 APPROVED

mohammad
 5/12/2021 2:07:55 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	132744	20.00	ng/ul	0.00
20) Naphthalene-d8	10.32	136	604102	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.19	164	404548	20.00	ng/ul	-0.01
64) Phenanthrene-d10	16.94	188	846857	20.00	ng/ul	0.00
79) Chrysene-d12	21.13	240	829401	20.00	ng/ul	0.00
88) Perylene-d12	23.28	264	741810	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	21143	6.36	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	6.73	99	103692	8.40	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.89	67	272351	35.58	ng/ul	0.00
11) 2-Chlorophenol-d4	7.08	132	297398	31.21	ng/ul	0.00
15) 4-Methylphenol-d8	8.26	113	203263	19.91	ng/ul	-0.01
21) Nitrobenzene-d5	8.70	128	169234	40.65	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.42	143	194466	41.39	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.95	165	329541	34.98	ng/ul	0.00
31) 4-Chloroaniline-d4	10.50	131	21201m	1.48	ng/ul	0.02
46) Dimethylphthalate-d6	13.62	166	1297097	41.37	ng/ul	0.00
49) Acenaphthylene-d8	13.88	160	1573603	39.60	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.44	143	23806	4.66	ng/ul	0.02
60) Fluorene-d10	15.19	176	1136014	42.68	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	204723	37.22	ng/ul	0.00
73) Anthracene-d10	17.04	188	1812436	42.50	ng/ul	0.00
81) Pyrene-d10	19.34	212	2068997	48.63	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.14	264	1767446	42.03	ng/ul	0.00

05/12/21

Target Compounds

					Ovalue	
47) Dimethylphthalate	13.66	163	40238	1.282	ng/ul	99
86) Bis(2-ethylhexyl)phthalate	21.05	149	457972	13.715	ng/ul#	98

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