

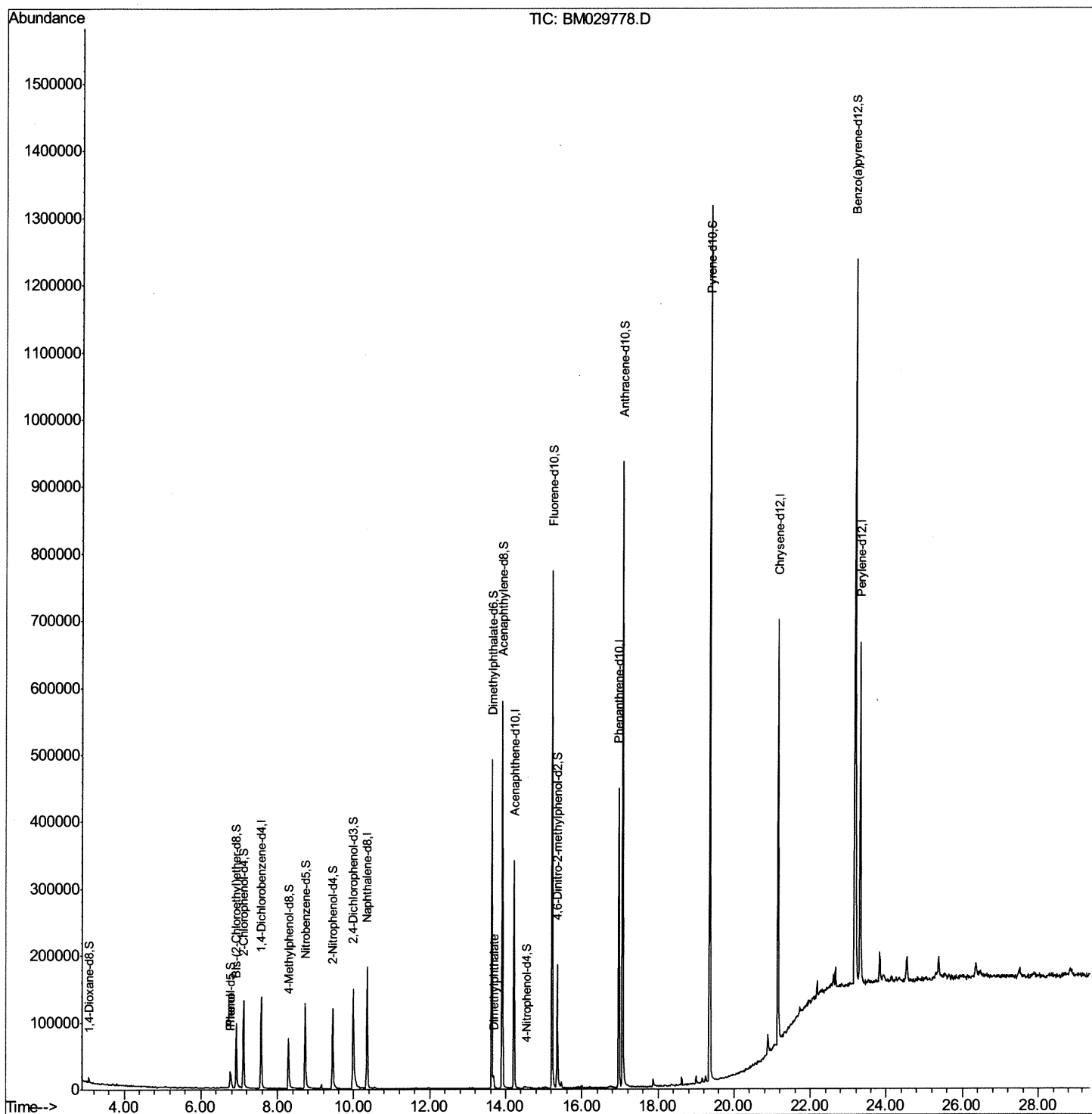
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\
Data File : BM029778.D
Acq On : 02 May 2021 11:04
Operator : CG/JU
Sample : M2144-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0GH4

Manual Integrations
APPROVED

mohammad
5/3/2021 12:25:31 PM

Quant Time: May 03 02:15:38 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050121.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat May 01 16:23:05 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

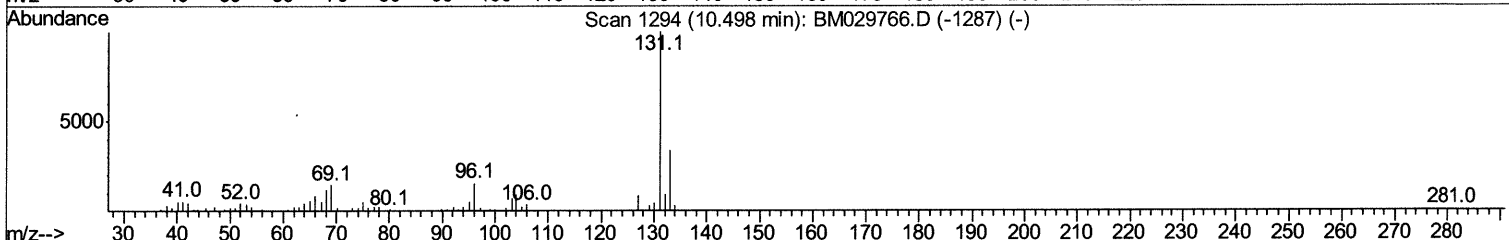
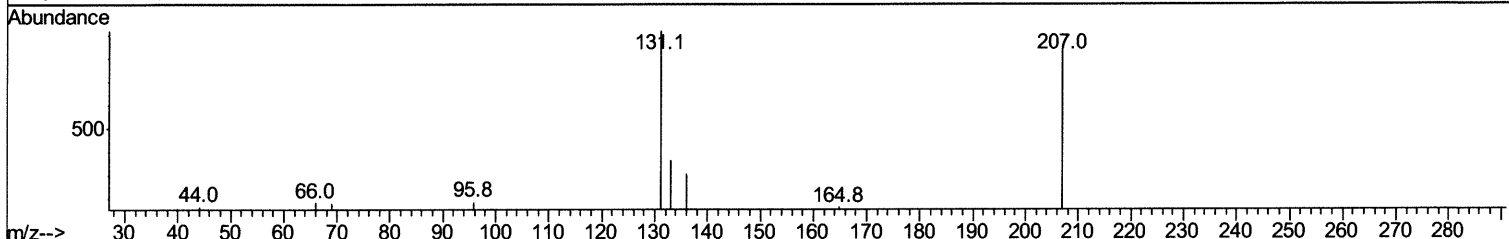
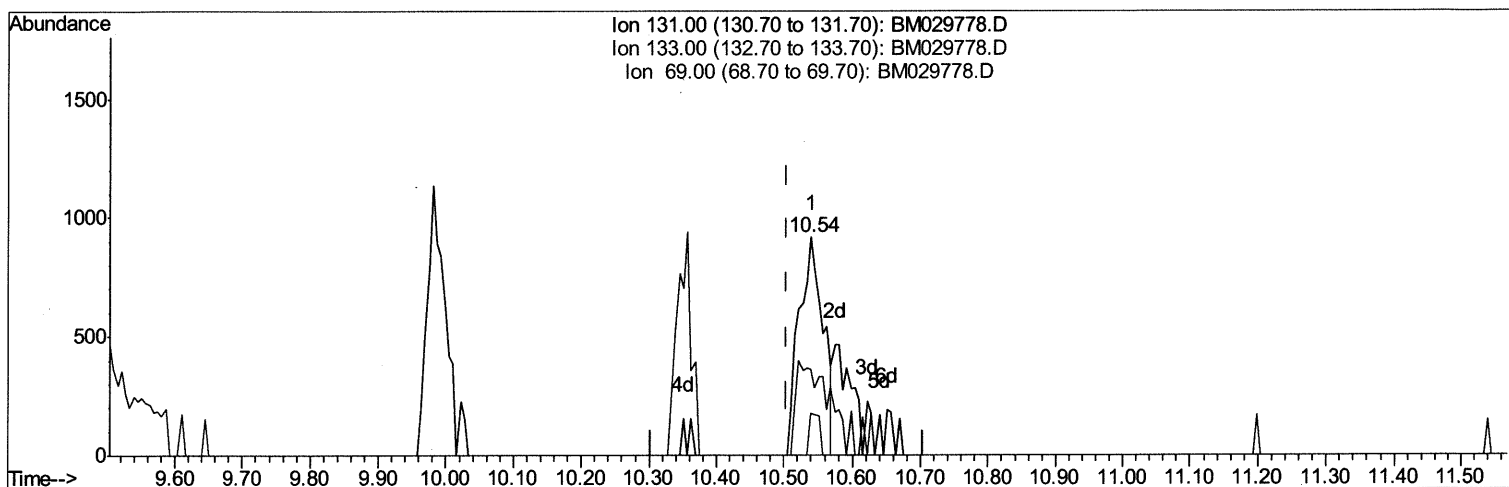
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TIC: BM029778.D

(31) 4-Chloroaniline-d4 (S)
 10.539min (+0.035) 0.62ng/ul
 response 2280

Ion	Exp%	Act%
131.00	100	100
133.00	33.40	39.39
69.00	16.40	19.37
0.00	0.00	0.00

Quantitation Report (Qedit)

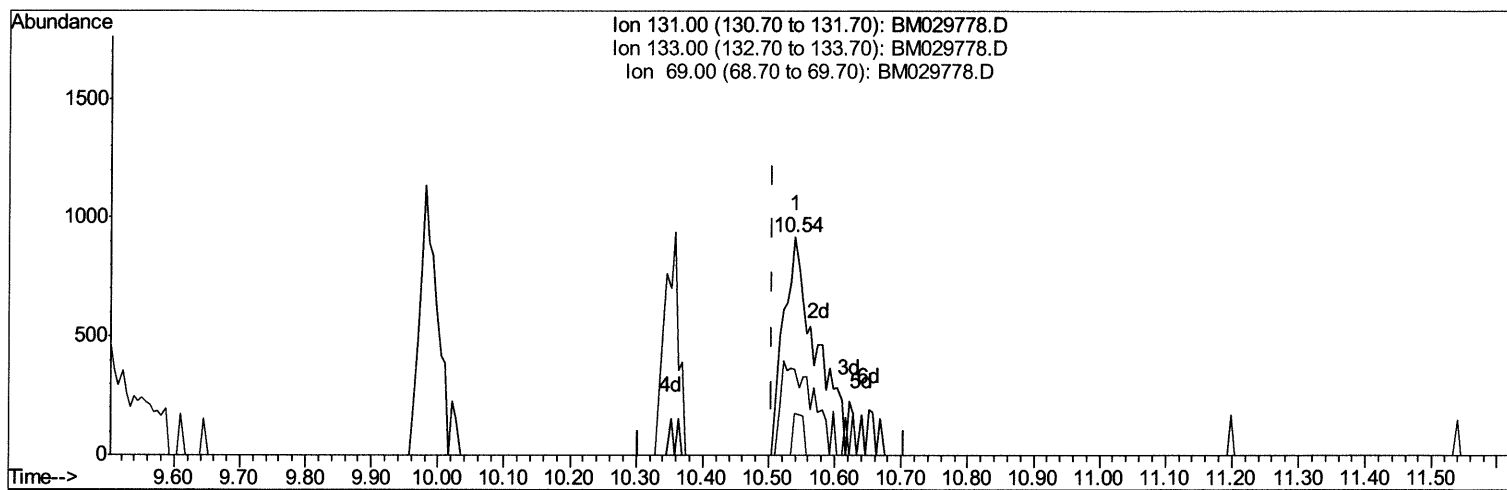
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\
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Manual Integrations
APPROVED

mohammad
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Quantitation Report (Qedit)

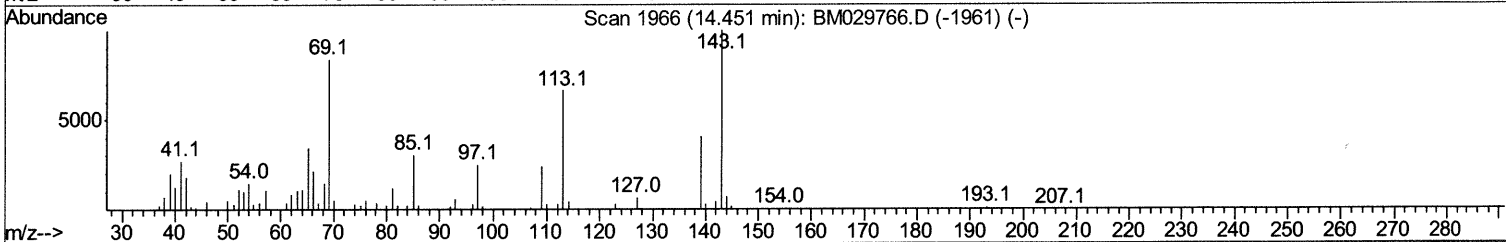
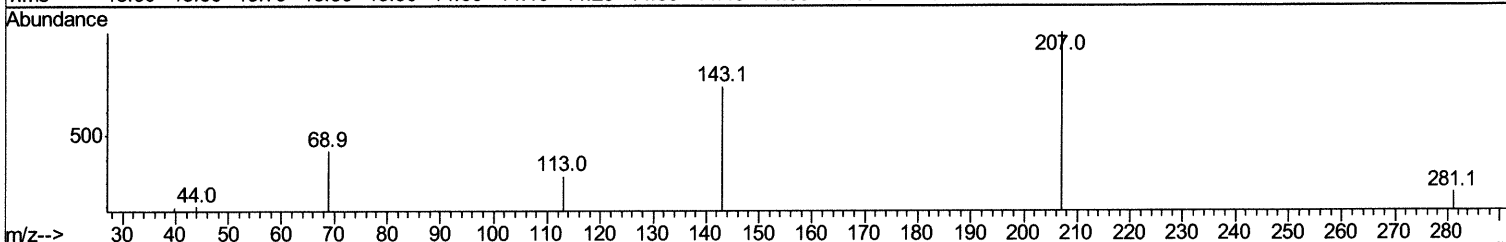
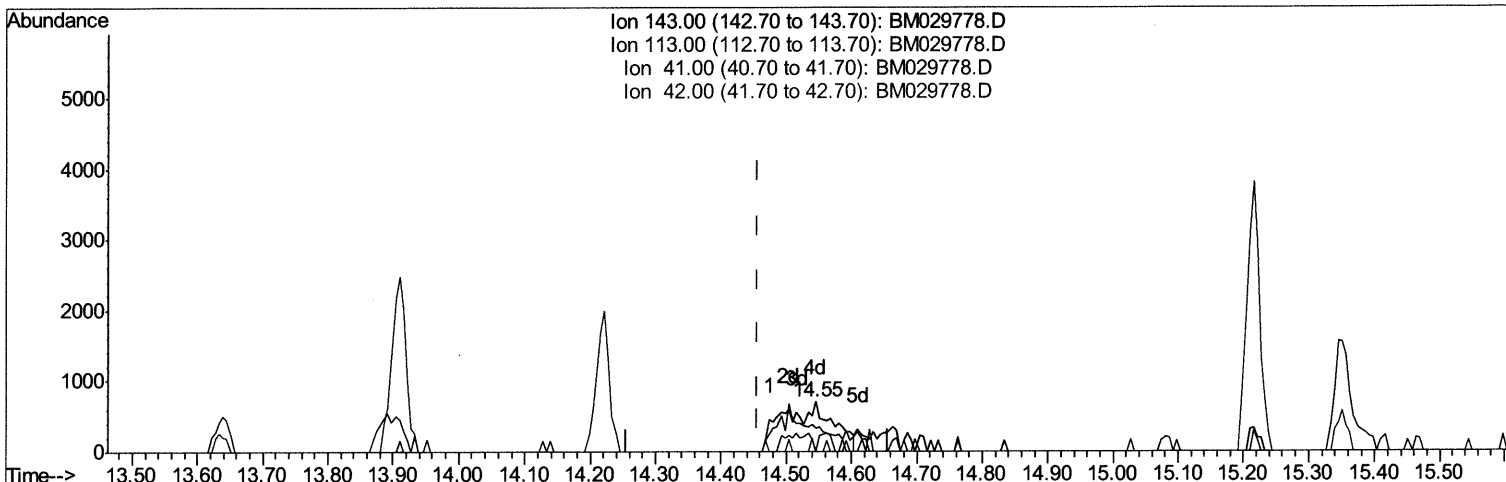
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\
 Data File : BM029778.D
 Acq On : 02 May 2021 11:04
 Operator : CG/JU
 Sample : M2144-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0GH4

Manual Integrations
 APPROVED

mohammad
 5/3/2021 12:25:31 PM

Quant Time: Mav 03 02:14:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050121.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat May 01 16:23:05 2021
 Response via : Initial Calibration



TIC: BM029778.D

(54) 4-Nitrophenol-d4 (S)

14.545min (+0.088) 2.74ng/ul m 05/04/2021

response 3727

Ion	Exp%	Act%
143.00	100	100
113.00	71.30	46.11#
41.00	32.40	0.00#
42.00	22.10	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	7.57	152	45796	20.00	ng/ul	0.00
20) Naphthalene-d8	10.35	136	168967	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.22	164	129896	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.96	188	305613	20.00	ng/ul	-0.01
79) Chrysene-d12	21.15	240	358650	20.00	ng/ul	0.00
88) Perylene-d12	23.32	264	393802	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	4814	5.14	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	6.76	99	17378	5.64	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.92	67	45973	28.46	ng/ul	0.00
11) 2-Chlorophenol-d4	7.11	132	71431	25.61	ng/ul	0.00
15) 4-Methylphenol-d8	8.29	113	36340	13.76	ng/ul	0.00
21) Nitrobenzene-d5	8.73	128	37155	30.20	ng/ul	0.00
24) 2-Nitrophenol-d4	9.45	143	51182	32.86	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.98	165	91954	28.25	ng/ul	0.00
31) 4-Chloroaniline-d4	10.54	131	3114m>	0.85	ng/ul>	0.04
46) Dimethylphthalate-d6	13.64	166	380592	36.99	ng/ul	0.00
49) Acenaphthylene-d8	13.91	160	444319	35.02	ng/ul	0.00
54) 4-Nitrophenol-d4	14.55	143	3727m>	2.74	ng/ul>	0.09
60) Fluorene-d10	15.22	176	343661	38.61	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.35	200	59889	26.95	ng/ul	0.00
73) Anthracene-d10	17.06	188	604553	39.24	ng/ul	0.00
81) Pyrene-d10	19.36	212	761721	41.59	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.18	264	887947	39.30	ng/ul	0.00

05/04/21 JU

05/04/21 JU

Target Compounds					Ovalue
8) Phenol	6.79	94	8664	2.835	ng/ul 94
47) Dimethylphthalate	13.69	163	13022	1.305	ng/ul# 98

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