

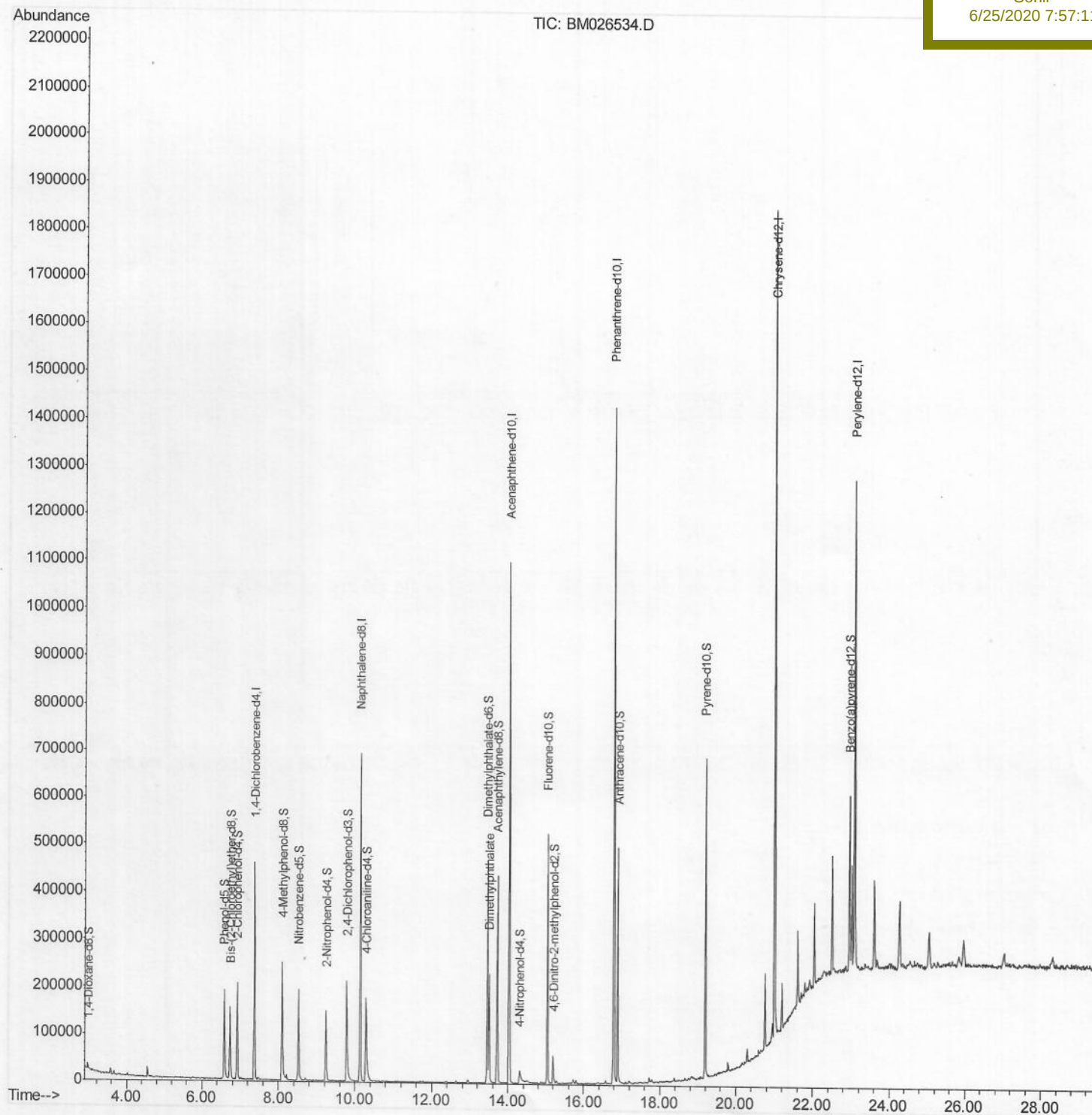
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062320\
Data File : BM026534.D
Acq On : 24 Jun 2020 16:08
Operator : JU/CG
Sample : L3064-01
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Jun 24 16:56:12 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 23 10:25:31 2020
Response via : Initial Calibration

Instrument :
BNA_M
Client Sampled :
ET154

Manual Integrations
APPROVED

Sohil
6/25/2020 7:57:11 AM



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062320\

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Operator : JU/CG

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ClientSampleId :

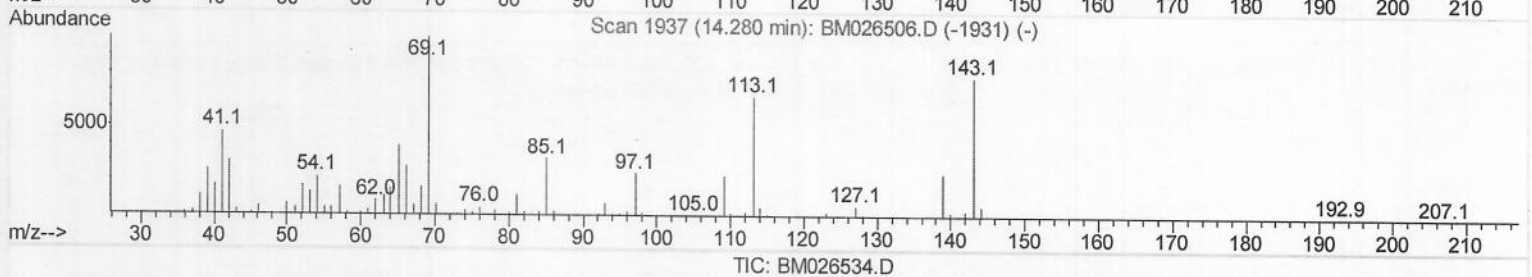
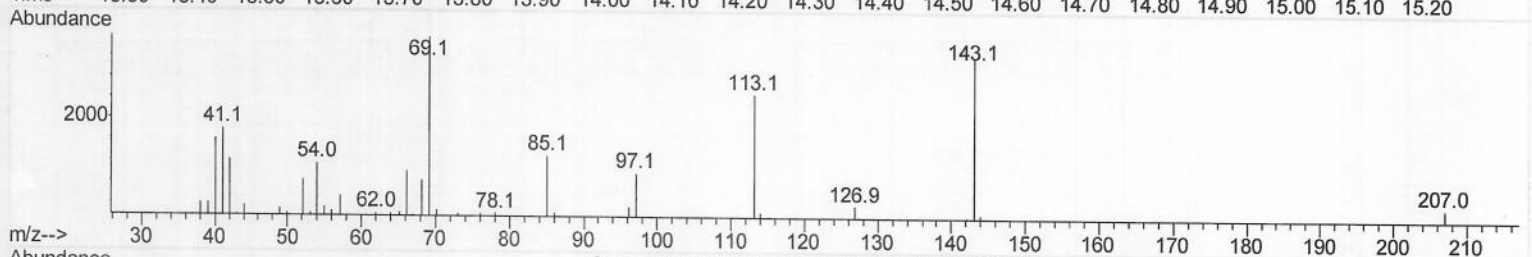
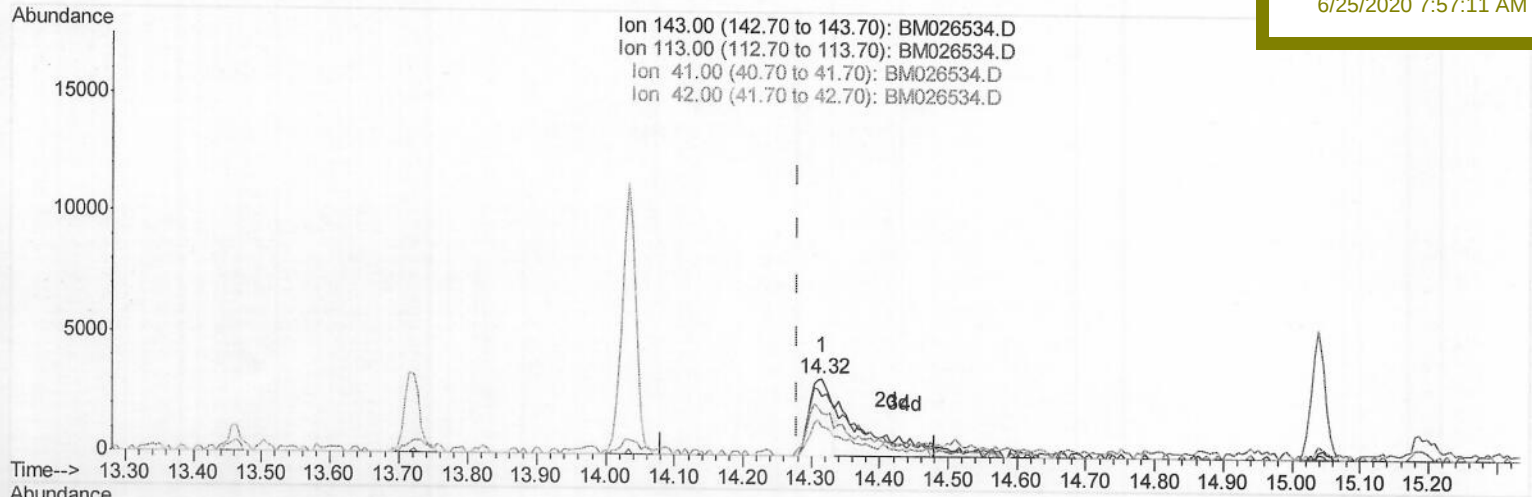
ET154

Manual Integrations

APPROVED

Sohil

6/25/2020 7:57:11 AM



(52) 4-Nitrophenol-d4 (S)

14.316min (+0.035) 1.56ng/ul

response 7200

Ion	Exp%	Act%
143.00	100	100
113.00	81.70	77.13
41.00	49.50	55.71
42.00	34.20	37.37

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BNA_M

ClientSampleId :

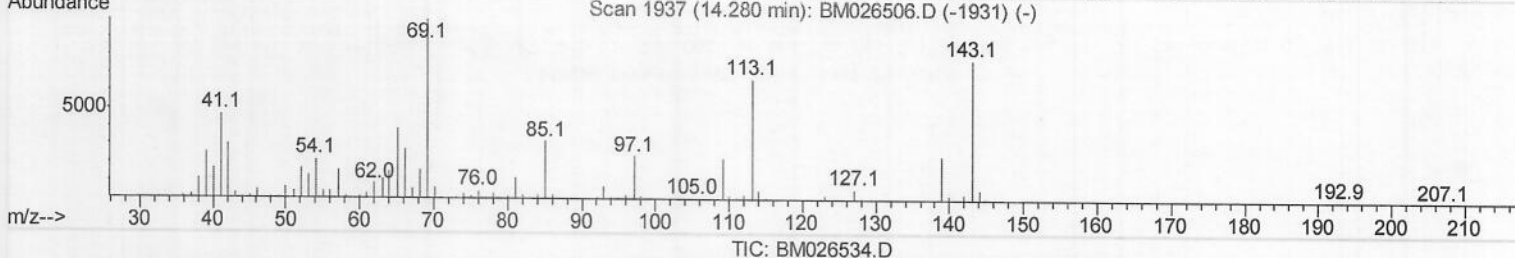
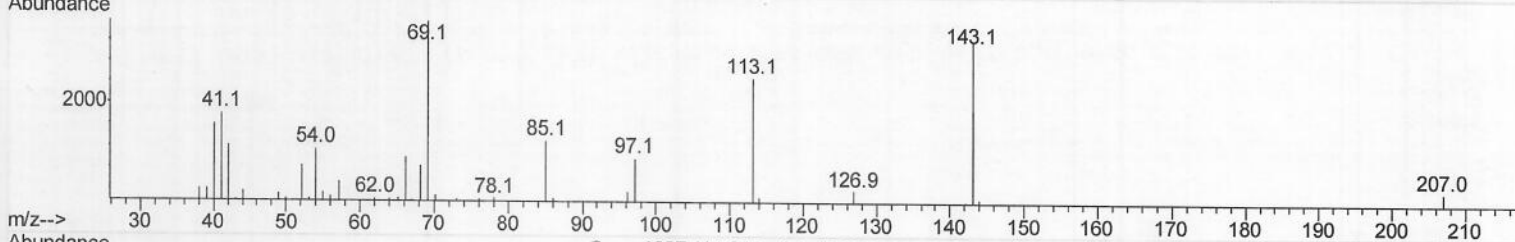
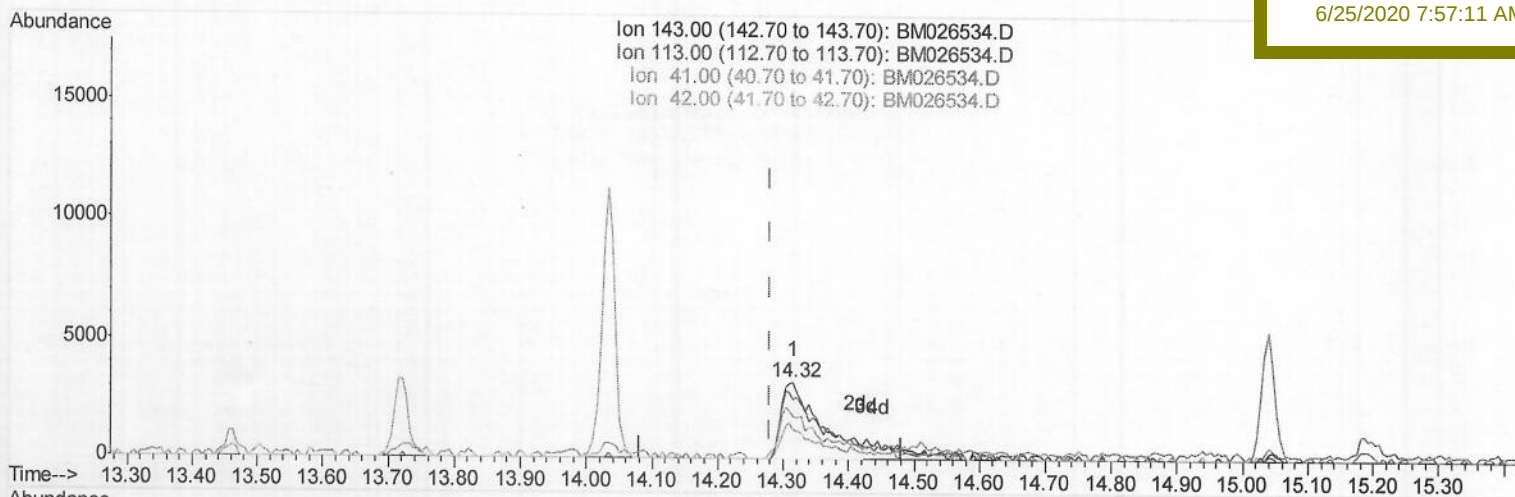
ET154

Manual Integrations

APPROVED

Sohil

6/25/2020 7:57:11 AM



(52) 4-Nitrophenol-d4 (S)

14.316min (+0.035) 2.91ng/ul m> 06/26/20 JW

response 13426

Ion	Exp%	Act%
143.00	100	100
113.00	81.70	77.13
41.00	49.50	55.71
42.00	34.20	37.37

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 Operator : JU/CG
 Sample : L3064-01
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Quant Time: Jun 24 16:56:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 23 10:25:31 2020
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 ET154

Manual Integrations
 APPROVED

Sohil
 6/25/2020 7:57:11 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	106676	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	484965	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	331345	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	741778	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	750490	20.00	ng/ul	0.00
86) Perylene-d12	23.09	264	675469	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.98	96	6192	2.04	ng/uL	0.00
5) Phenol-d5	6.59	99	98081	9.98	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.73	67	71429	11.49	ng/ul	0.00
9) 2-Chlorophenol-d4	6.92	132	78752	9.83	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	83261	10.76	ng/ul	0.00
19) Nitrobenzene-d5	8.53	128	37542	8.71	ng/ul	0.00
22) 2-Nitrophenol-d4	9.24	143	38461	8.11	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.78	165	77190	8.84	ng/ul	0.00
29) 4-Chloroaniline-d4	10.29	131	102292	11.24	ng/ul	0.00
44) Dimethylphthalate-d6	13.46	166	270434	9.71	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	271878	8.14	ng/ul	0.00
52) 4-Nitrophenol-d4	14.32	143	13426m	2.91	ng/ul	0.04
58) Fluorene-d10	15.04	176	190120	7.95	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.19	200	11243	2.32	ng/ul	0.00
71) Anthracene-d10	16.89	188	269830	7.15	ng/ul	0.00
79) Pyrene-d10	19.20	212	316364	7.69	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.96	264	234071	6.27	ng/ul	0.00

Target Compounds					Ovalue
45) Dimethylphthalate	13.51	163	133458	4.613	ng/ul 98

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

06/26/20 JU