

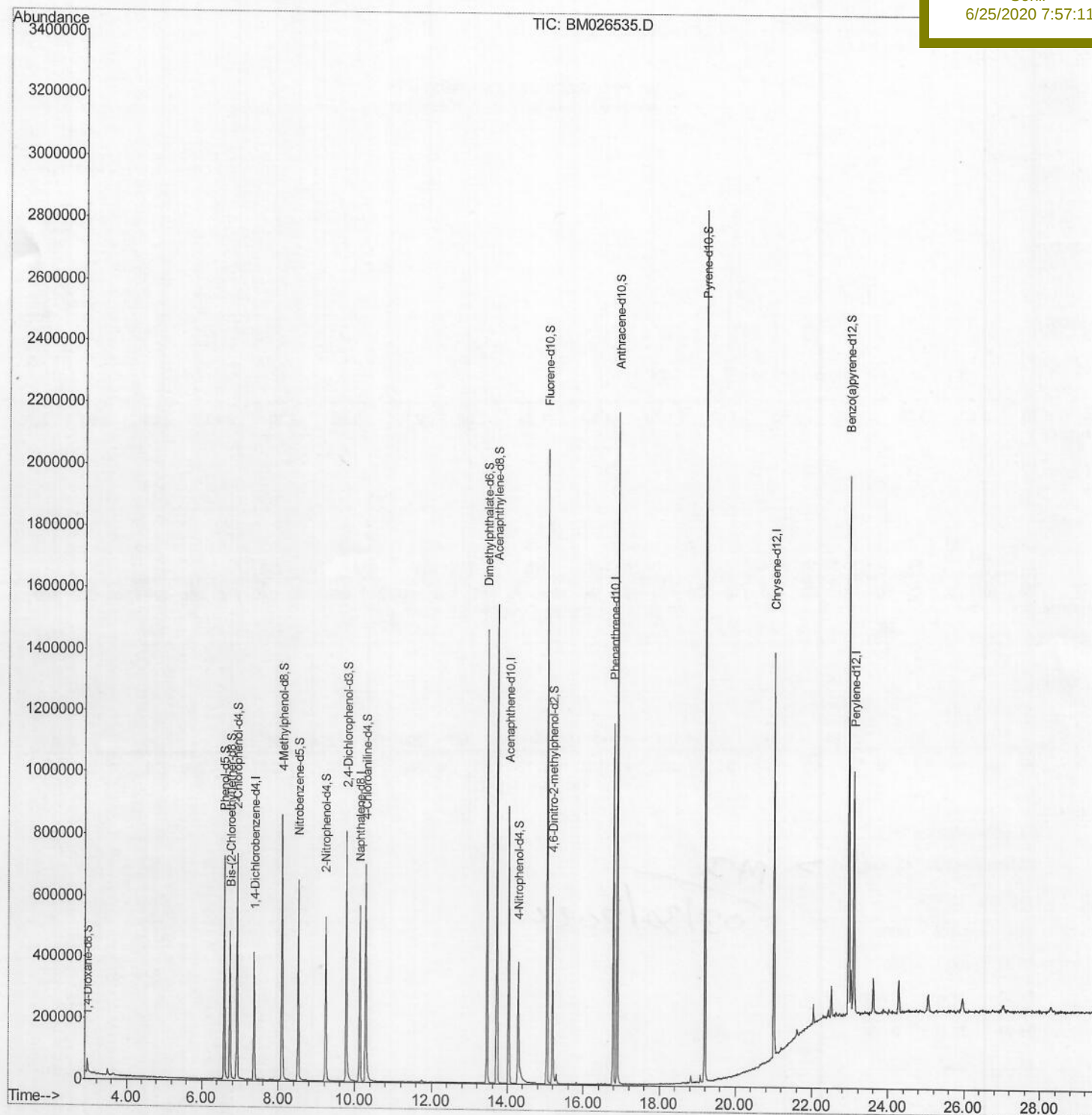
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062320\
Data File : BM026535.D
Acq On : 24 Jun 2020 16:45
Operator : JU/CG
Sample : PB129633BL
Misc :
ALS Vial : 45 Sample Multiplier: 1

Quant Time: Jun 24 17:51:07 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 :
SBLK33

Manual Integrations
APPROVED

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



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Acq On : 24 Jun 2020 16:45

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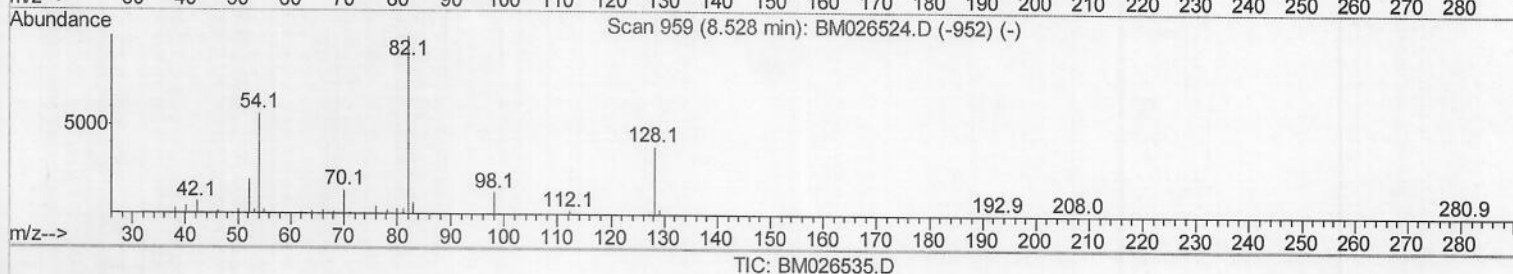
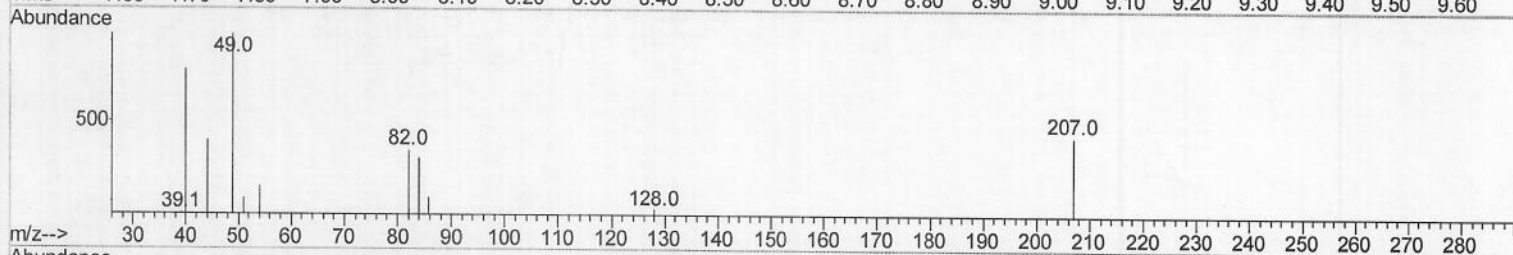
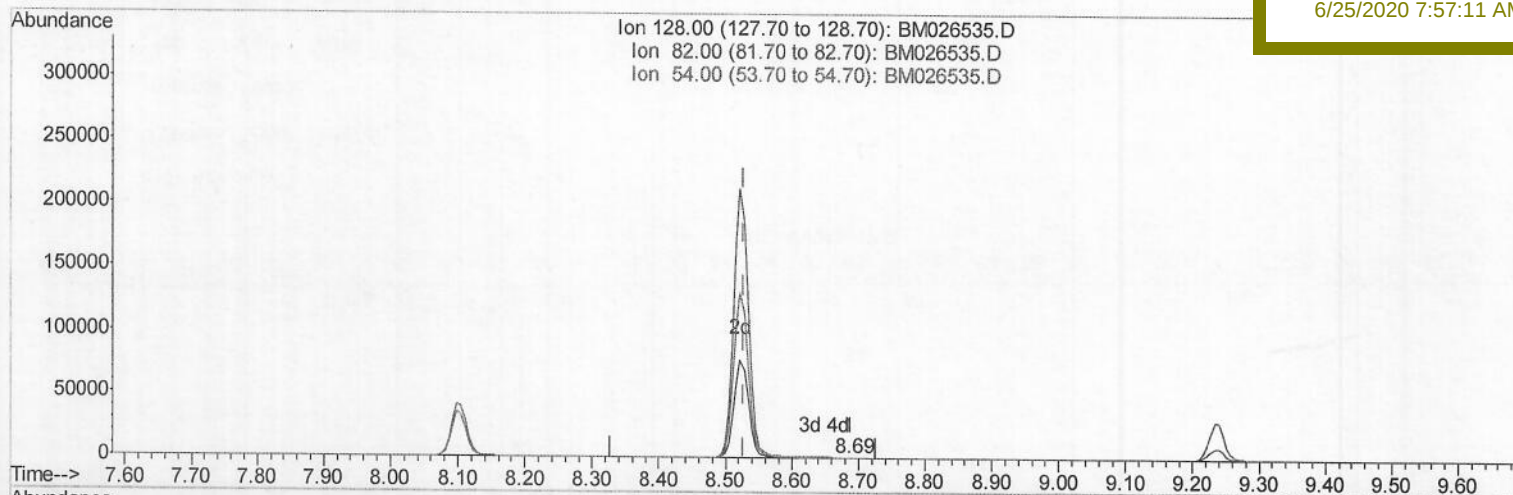
SBLK33

Manual Integrations

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Sohil

6/25/2020 7:57:11 AM



TIC: BM026535.D

(19) Nitrobenzene-d5 (S)

8.687min (+0.159) 0.03ng/ul

response 123

Ion	Exp%	Act%
128.00	100	100
82.00	257.80	225.14
54.00	139.90	147.43
0.00	0.00	0.00

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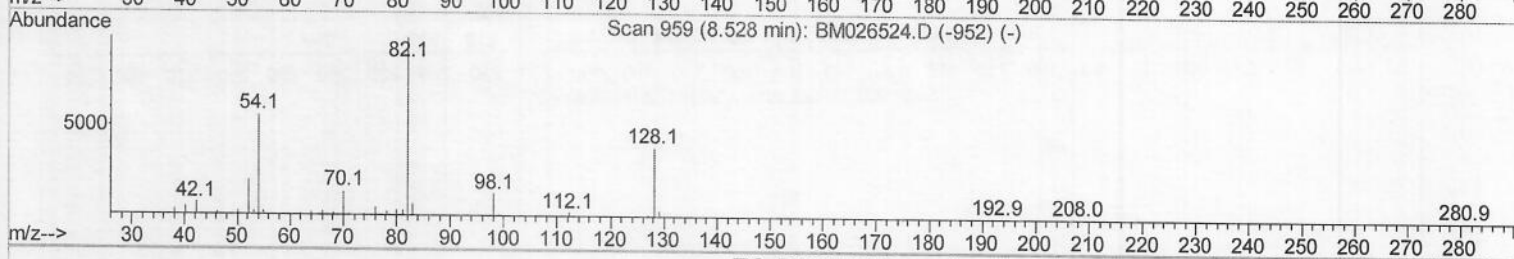
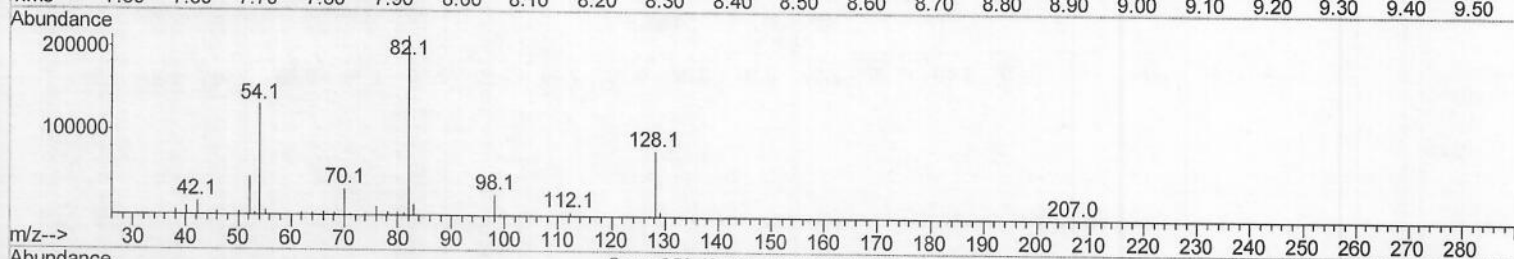
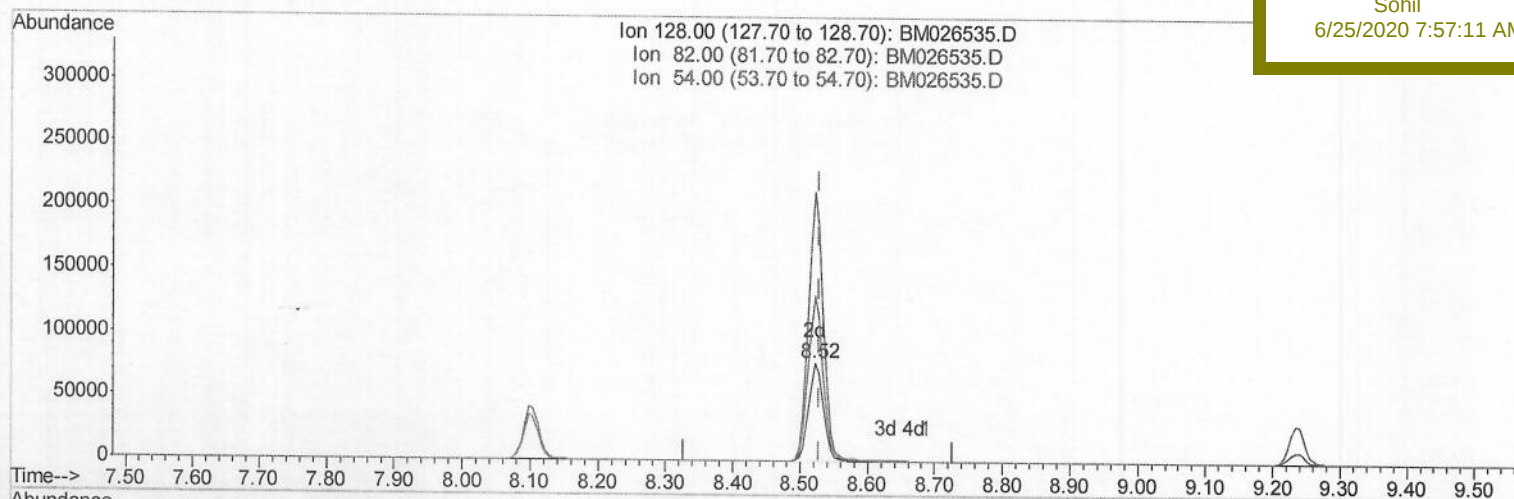
SBLK33

Manual Integrations

APPROVED

Sohil

6/25/2020 7:57:11 AM



TIC: BM026535.D

(19) Nitrobenzene-d5 (S)

8.522min (-0.006) 33.40ng/ul m >SV 06/26/20

response 117815

Ion	Exp%	Act%
128.00	100	100
82.00	257.80	274.17
54.00	139.90	168.57#
0.00	0.00	0.00

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Instrument :
 BNA_M
 Client Sample Id :
 SBLK33

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	95568	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	396957	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	270139	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	621346	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	604685	20.00	ng/ul	0.00
86) Perylene-d12	23.09	264	531976	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.96	96	20513	7.53	ng/uL	0.00
5) Phenol-d5	6.59	99	358476	40.72	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.73	67	214809	38.57	ng/ul	0.00
9) 2-Chlorophenol-d4	6.92	132	265420	36.98	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	288716	41.66	ng/ul	0.00
19) Nitrobenzene-d5	8.52	128	117815m)	33.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.24	143	136224	35.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.77	165	263806	36.93	ng/ul	0.00
29) 4-Chloroaniline-d4	10.29	131	352297	47.29	ng/ul	0.00
44) Dimethylphthalate-d6	13.46	166	843244	37.12	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	964532	35.40	ng/ul	0.00
52) 4-Nitrophenol-d4	14.28	143	111058	29.48	ng/ul	0.00
58) Fluorene-d10	15.04	176	719947	36.94	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.18	200	109975	27.04	ng/ul	0.00
71) Anthracene-d10	16.89	188	1141050	36.08	ng/ul	0.00
79) Pyrene-d10	19.20	212	1297075	39.12	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.96	264	1079394	36.70	ng/ul	0.00

JU-06/26/20

Target Compounds

Ovalue

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