

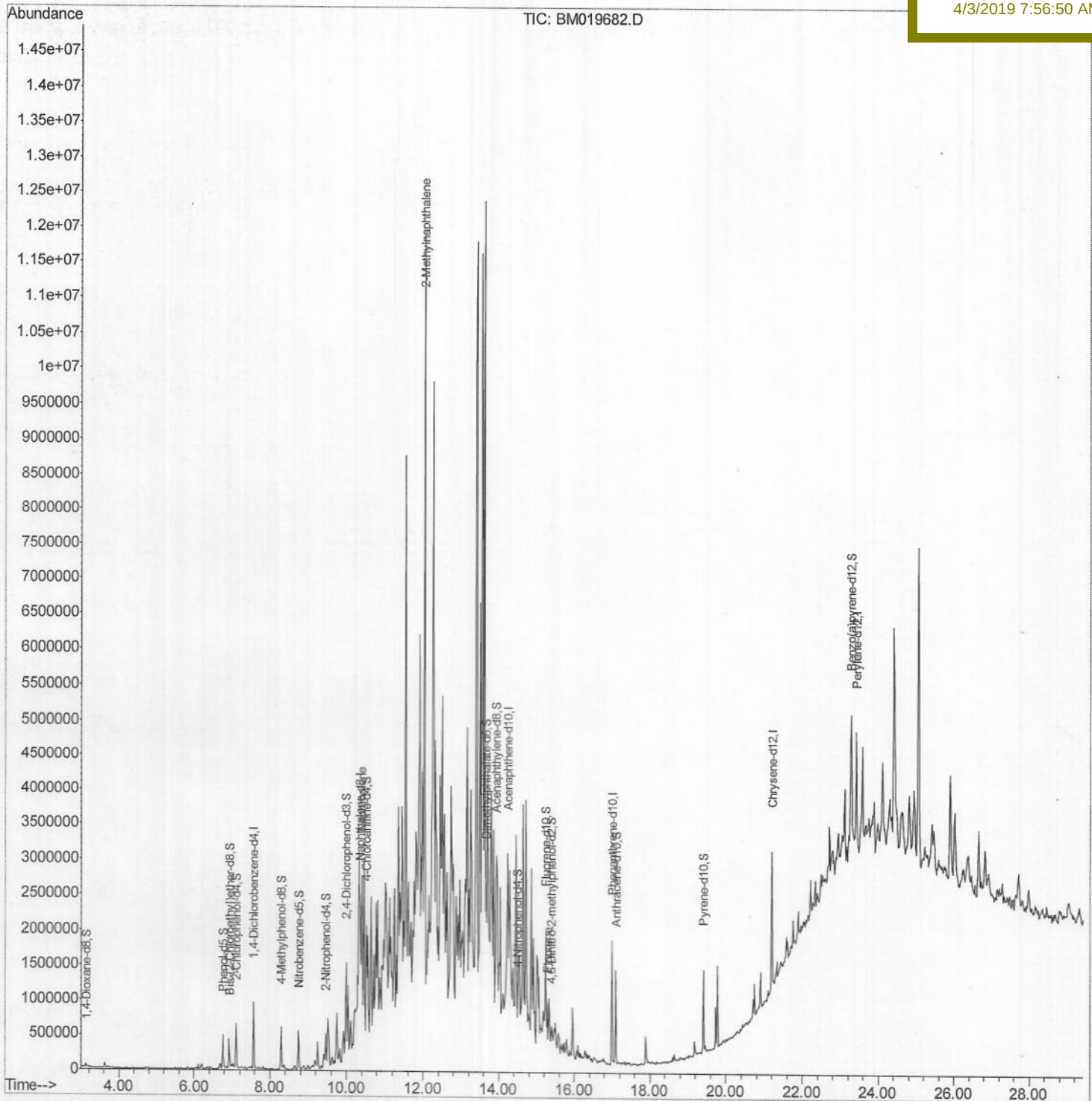
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040119\
Data File : BM019682.D
Acq On : 02 Apr 2019 07:34
Operator : JU/SJ
Sample : K2119-19
Misc :
ALS Vial : 36 Sample Multiplier: 1

Quant Time: Apr 03 05:17:16 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 28 05:07:19 2019
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
BF5B6

Manual Integrations
APPROVED

Sohil
4/3/2019 7:56:50 AM



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM040119\
Data File : BM019682.D
Acq On : 02 Apr 2019 07:34
Operator : JU/SJ
Sample : K2119-19
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ALS Vial : 36 Sample Multiplier: 1

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Instrument :

BNA_M

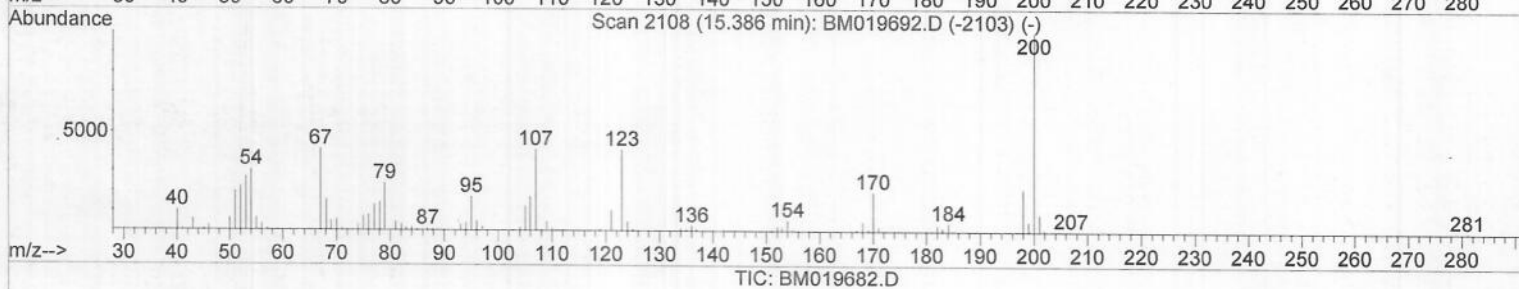
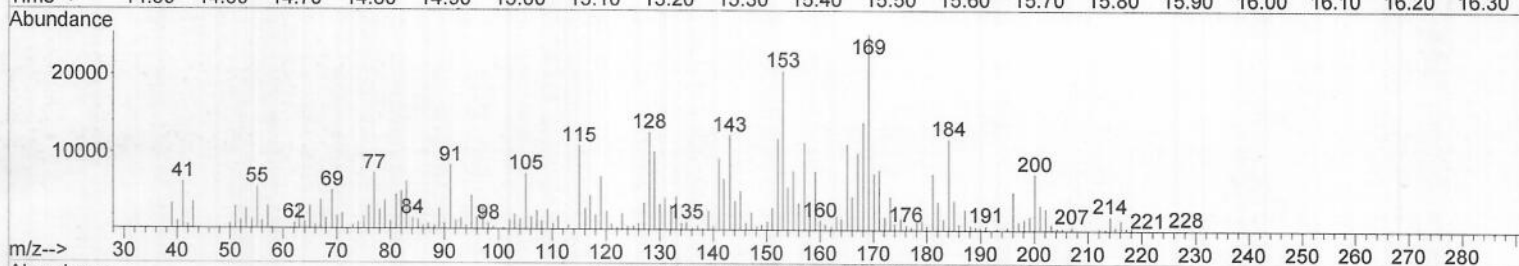
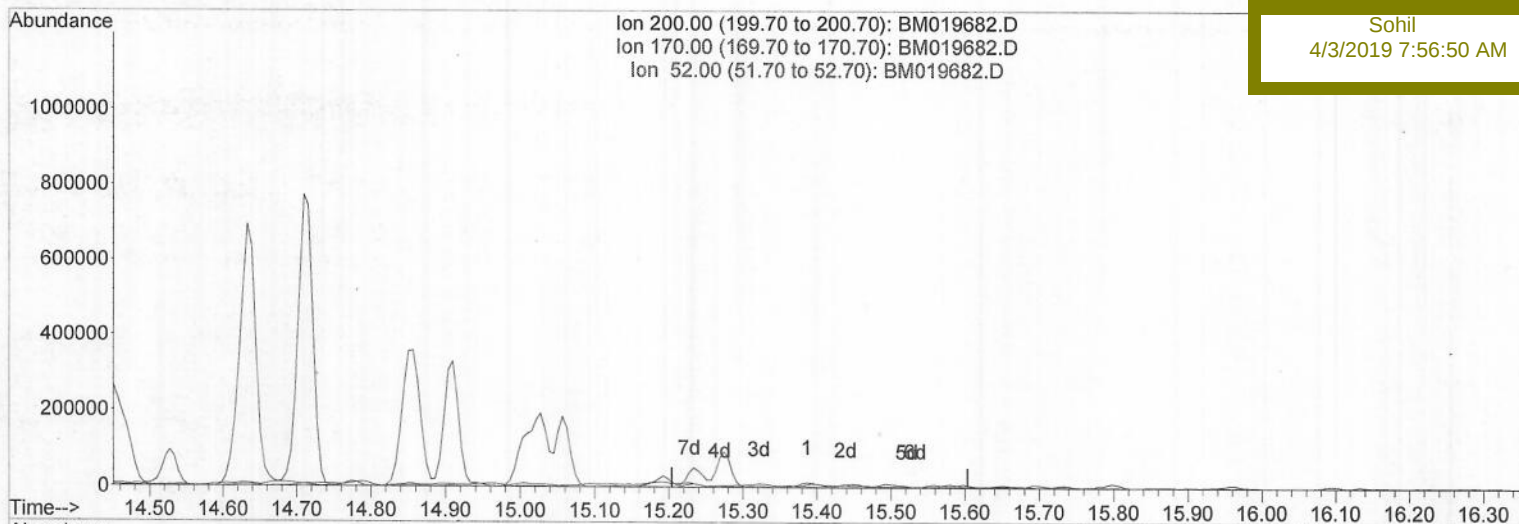
ClientSampleId :

BF5B6

Manual Integrations
APPROVED

Sohil

4/3/2019 7:56:50 AM



(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.404min (-15.404) 0.00ng/ul d

response 0

Ion	Exp%	Act%
200.00	100	0.00
170.00	17.70	0.00
52.00	31.50	0.00
0.00	0.00	0.00

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Instrument :

BNA_M

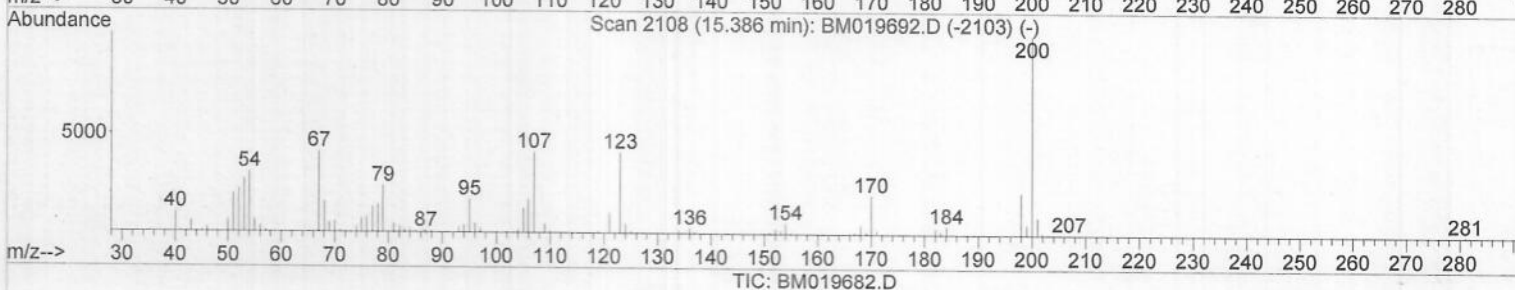
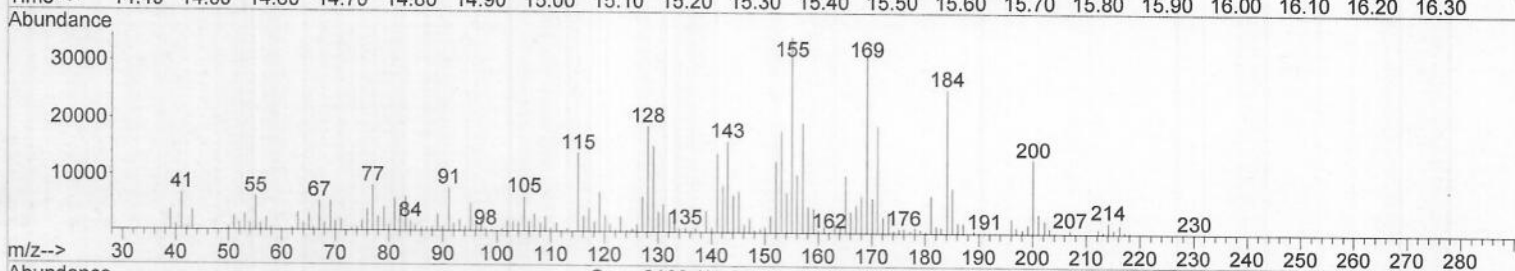
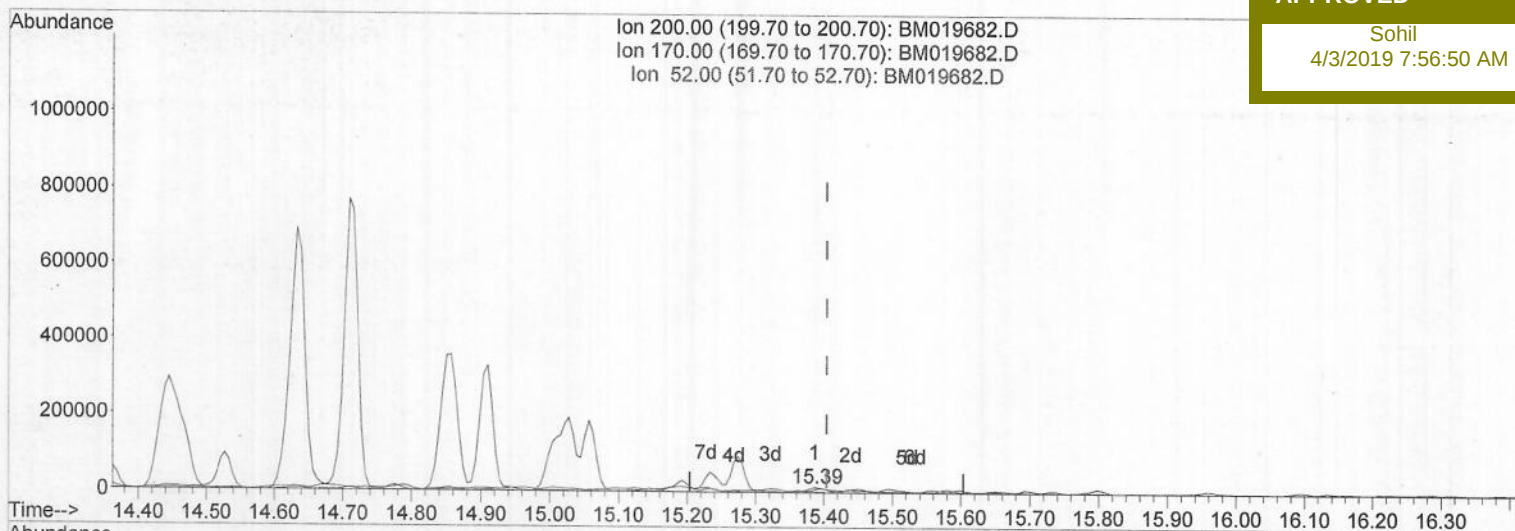
ClientSampleId :

BF5B6

Manual Integrations
APPROVED

Sohil

4/3/2019 7:56:50 AM



(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.386min (-0.017) 2.86ng/ul m

response 17533

Ion	Exp%	Act%
200.00	100	100
170.00	17.70	48.64#
52.00	31.50	13.32#
0.00	0.00	0.00

SJ
04/02/19

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040119\
 Data File : BM019682.D
 Acq On : 02 Apr 2019 07:34
 Operator : JU/SJ
 Sample : K2119-19
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 Client Sampled :
 BF5B6

Quant Time: Apr 03 05:17:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 28 05:07:19 2019
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Sohil
 4/3/2019 7:56:50 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	254486	20.00	ng/ul	-0.03
18) Naphthalene-d8	10.35	136	1073307	20.00	ng/ul	-0.03
36) Acenaphthene-d10	14.23	164	533302	20.00	ng/ul	-0.02
62) Phenanthrene-d10	16.99	188	926664	20.00	ng/ul	-0.02
78) Chrysene-d12	21.20	240	816034	20.00	ng/ul	-0.02
86) Perylene-d12	23.42	264	956022	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	28300	4.35	ng/uL	-0.02
5) Phenol-d5	6.76	99	318790	14.07	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	6.92	67	221353	14.59	ng/ul	-0.03
9) 2-Chlorophenol-d4	7.10	132	269886	15.71	ng/ul	-0.03
13) 4-Methylphenol-d8	8.29	113	251838	14.80	ng/ul	-0.02
19) Nitrobenzene-d5	8.74	128	127547	16.66	ng/ul	-0.03
22) 2-Nitrophenol-d4	9.45	143	135365	15.89	ng/ul	-0.03
26) 2,4-Dichlorophenol-d3	9.99	165	247963	15.42	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.52	131	133286	6.90	ng/ul	-0.02
44) Dimethylphthalate-d6	13.66	166	714152	16.60	ng/ul	-0.02
47) Acenaphthylene-d8	13.93	160	855869	15.91	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.49	143	24742	3.67	ng/ul	0.00
58) Fluorene-d10	15.23	176	552159	14.89	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.39	200	17533m	2.86	ng/ul	-0.02
71) Anthracene-d10	17.09	188	686053	15.43	ng/ul	-0.02
79) Pyrene-d10	19.40	212	592645	15.72	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.27	264	677390	13.37	ng/ul	-0.02

Target Compounds

				Ovalue	
28) Naphthalene	10.40	128	801063	14.407 ng/ul#	93
34) 2-Methylnaphthalene	12.04	142	5309546	136.281 ng/ul	100
59) Fluorene	15.29	166	66201	1.593 ng/ul#	97

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