

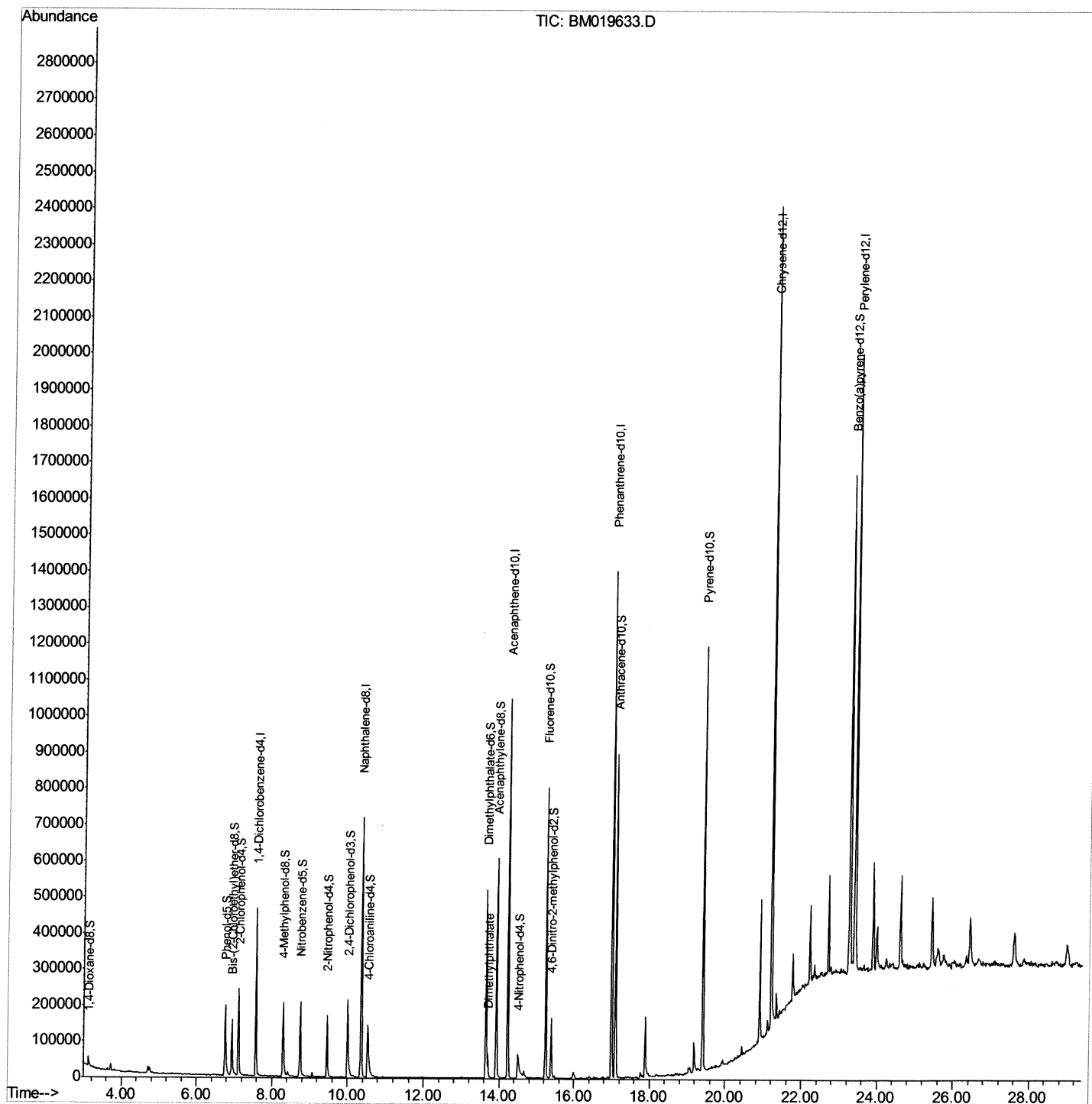
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033019\
Data File : BM019633.D
Acq On : 30 Mar 2019 23:31
Operator : JU/SJ
Sample : K2068-10
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BF5K6

Manual Integrations
APPROVED

Sohil
4/1/2019 5:50:55 PM

Quant Time: Mar 31 02:24:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 29 14:33:17 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

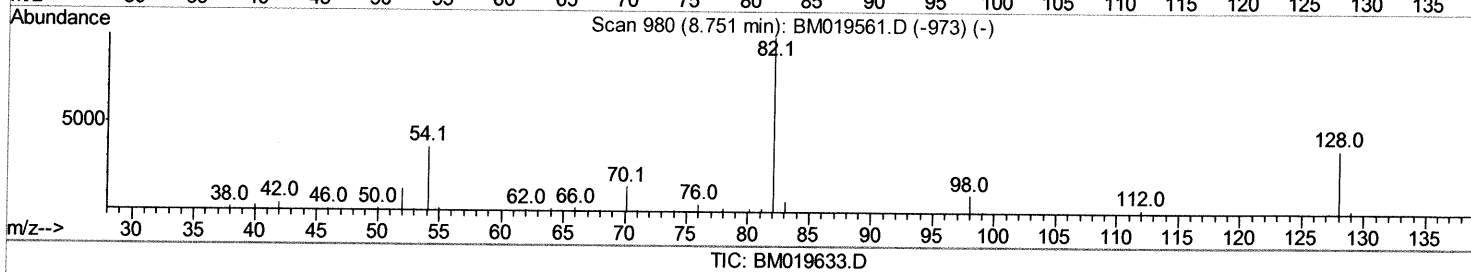
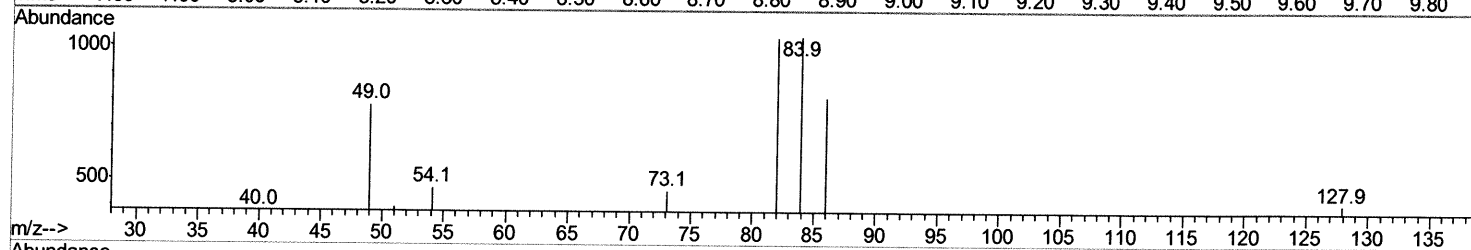
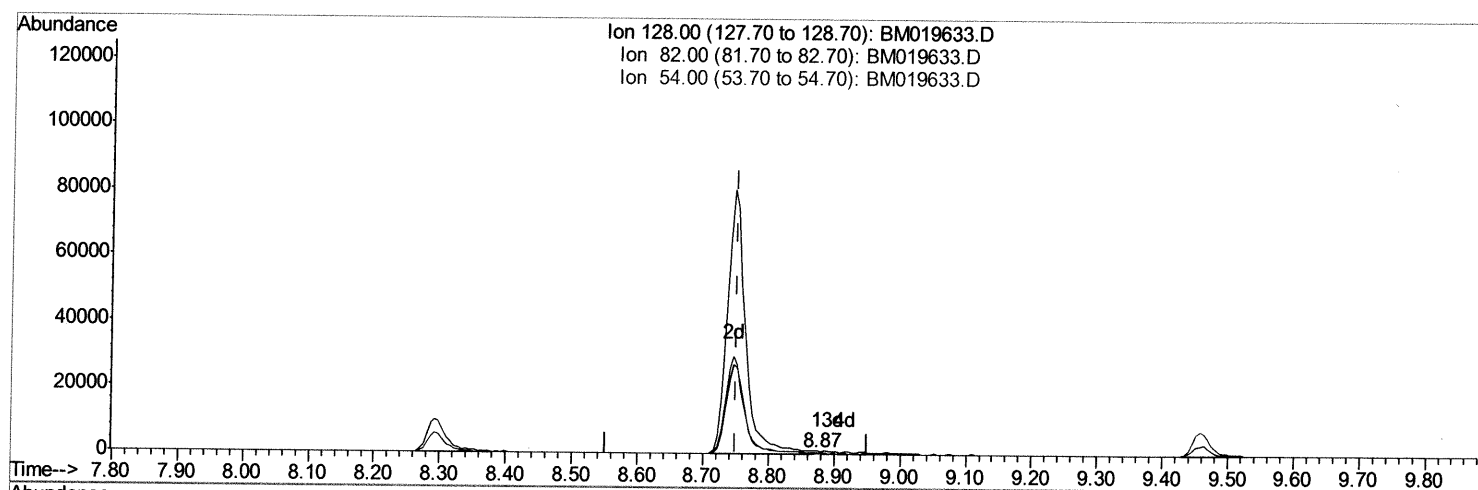
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(19) Nitrobenzene-d5 (S)

8.875min (+0.124) 0.04ng/ul

response 146

Ion	Exp%	Act%
128.00	100	100
82.00	244.60	250.72
54.00	108.20	113.53
0.00	0.00	0.00

Quantitation Report (Qedit)

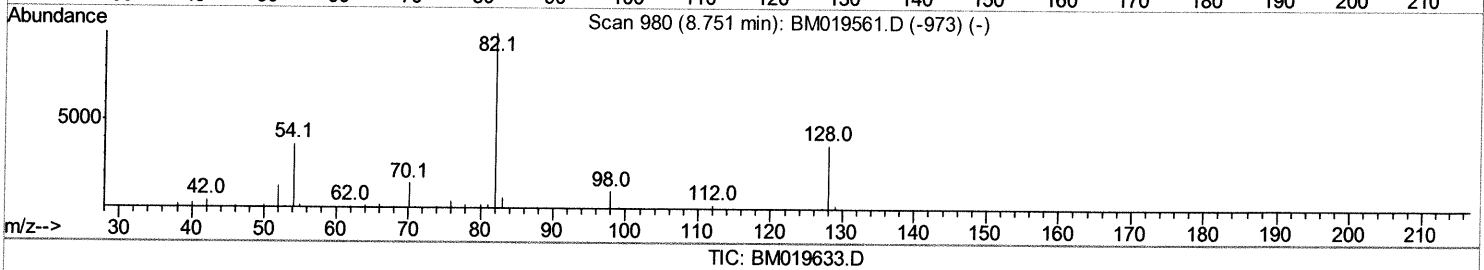
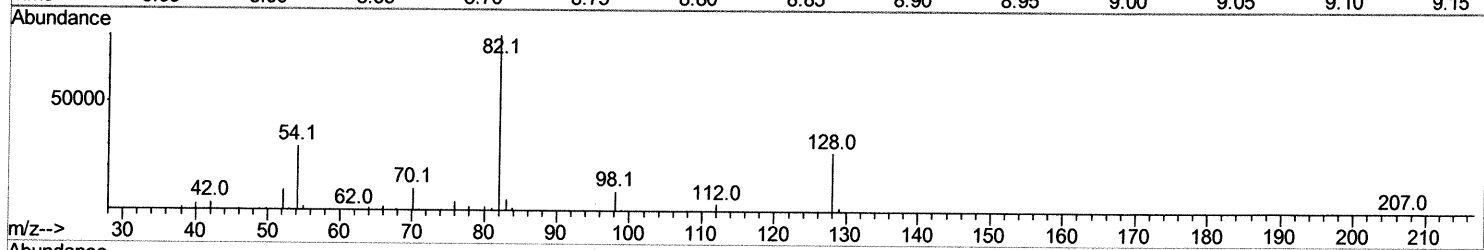
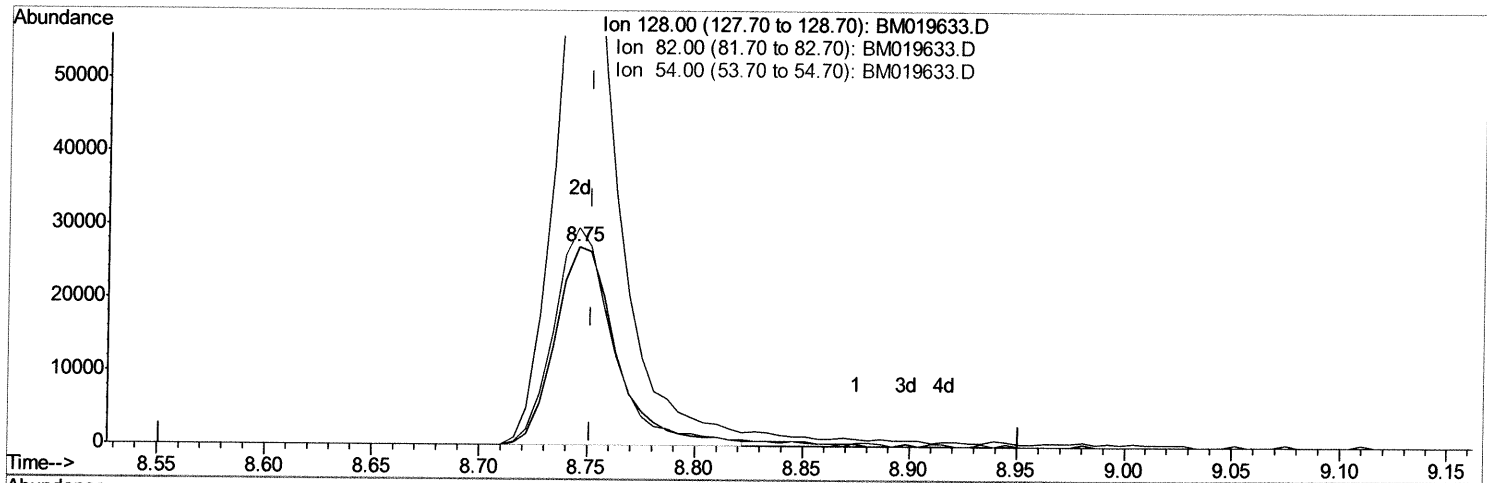
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Manual Integrations
 APPROVED

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 4/1/2019 5:50:55 PM

Quant Time: Mar 31 02:22:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 29 14:33:17 2019
 Response via : Initial Calibration



(19) Nitrobenzene-d5 (S)

8.745min (-0.006) 13.24ng/ul m> SJ 4/2/19

response 54101

Ion	Exp%	Act%
128.00	100	100
82.00	244.60	298.40#
54.00	108.20	109.86
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033019\
 Data File : BM019633.D
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 Operator : JU/SJ
 Sample : K2068-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BF5K6

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:50:55 PM

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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	128119	20.00	nq/ul	0.00
18) Naphthalene-d8	10.35	136	572615	20.00	nq/ul	0.00
36) Acenaphthene-d10	14.23	164	326018	20.00	nq/ul	0.00
62) Phenanthrene-d10	16.99	188	742196	20.00	nq/ul	0.00
78) Chrysene-d12	21.20	240	933191	20.00	nq/ul	0.00
86) Perylene-d12	23.40	264	1139650	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	17961	5.49	nq/uL	0.00
5) Phenol-d5	6.76	99	141201	12.38	nq/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.93	67	99630	13.04	nq/ul	0.00
9) 2-Chlorophenol-d4	7.11	132	117503	13.59	nq/ul	0.00
13) 4-Methylphenol-d8	8.29	113	96919	11.32	nq/ul	0.00
19) Nitrobenzene-d5	8.75	128	54101m	13.24	nq/ul	0.00
22) 2-Nitrophenol-d4	9.46	143	59735	13.14	nq/ul	0.00
26) 2,4-Dichlorophenol-d3	9.99	165	99550	11.61	nq/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	129938	12.61	nq/ul	0.00
44) Dimethylphthalate-d6	13.66	166	366550	13.93	nq/ul	0.00
47) Acenaphthylene-d8	13.93	160	426012	12.95	nq/ul	0.00
52) 4-Nitrophenol-d4	14.51	143	38178	9.26	nq/ul	0.03
58) Fluorene-d10	15.23	176	311069	13.73	nq/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	38040	7.75	nq/ul	0.00
71) Anthracene-d10	17.09	188	470966	13.22	nq/ul	0.00
79) Pyrene-d10	19.40	212	549513	12.75	nq/ul	0.00
90) Benzo(a)pyrene-d12	23.26	264	723505	11.98	ng/ul	0.00

52 4/2/19

Target Compounds				Ovalue	
45) Dimethylphthalate	13.70	163	41397	1.568 nq/ul#	98

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