

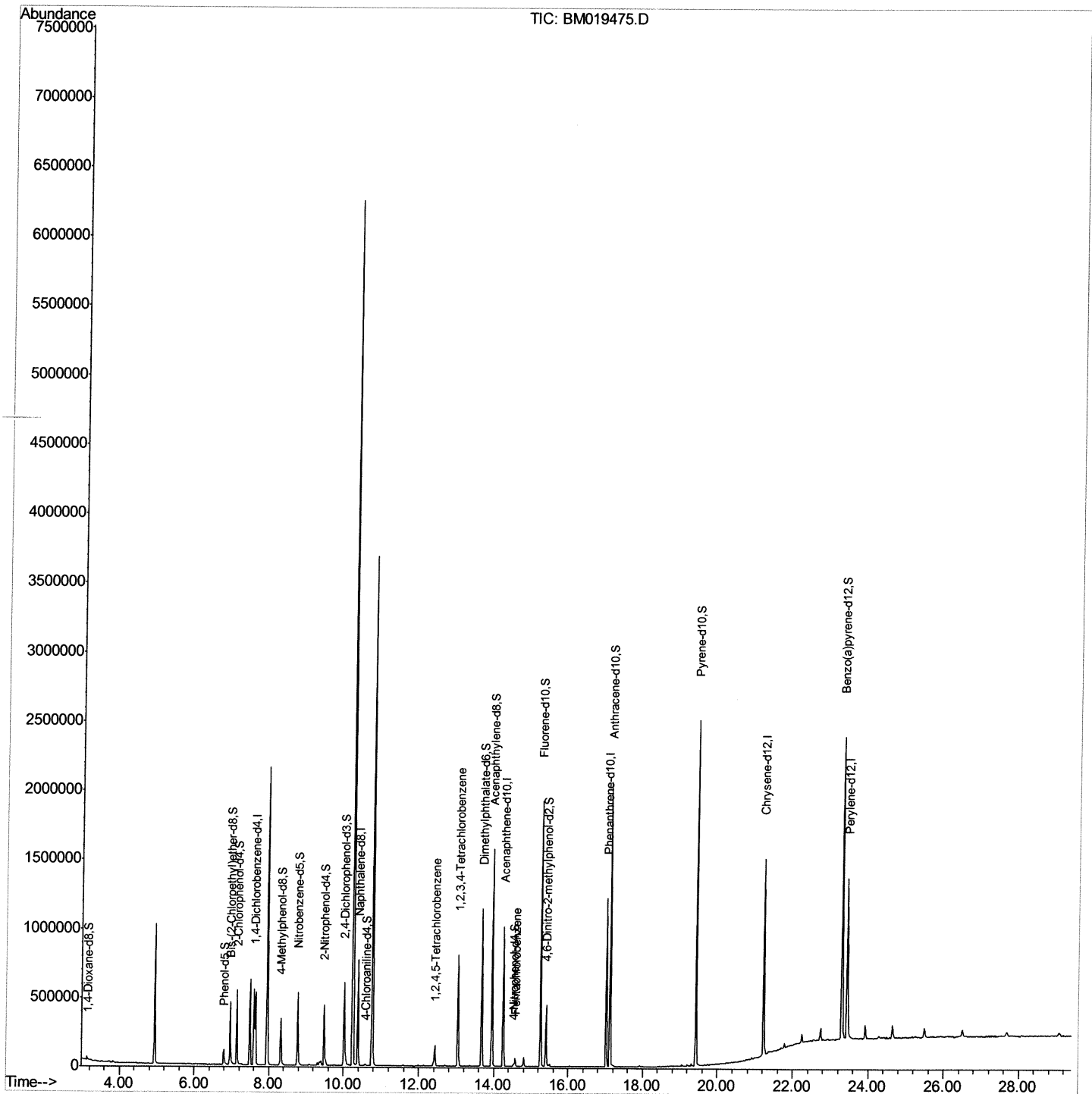
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019475.D
Acq On : 26 Mar 2019 18:03
Operator : JU/SJ
Sample : K2069-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09D6

Manual Integrations
APPROVED

Sohil
3/27/2019 5:46:29 PM

Quant Time: Mar 27 04:01:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 27 03:30:35 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

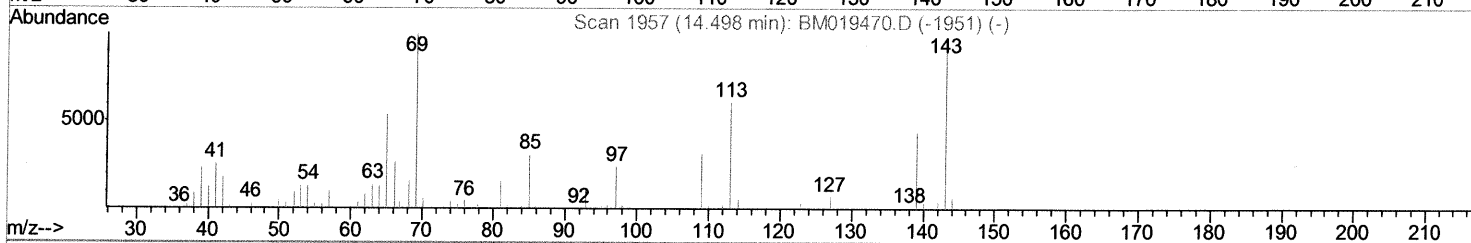
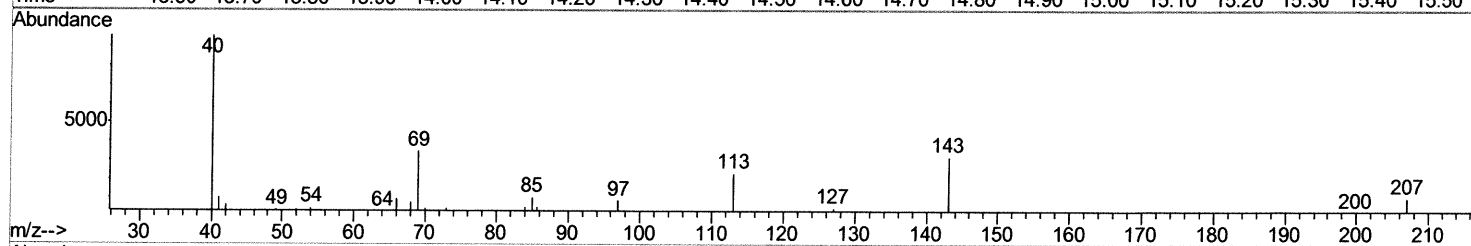
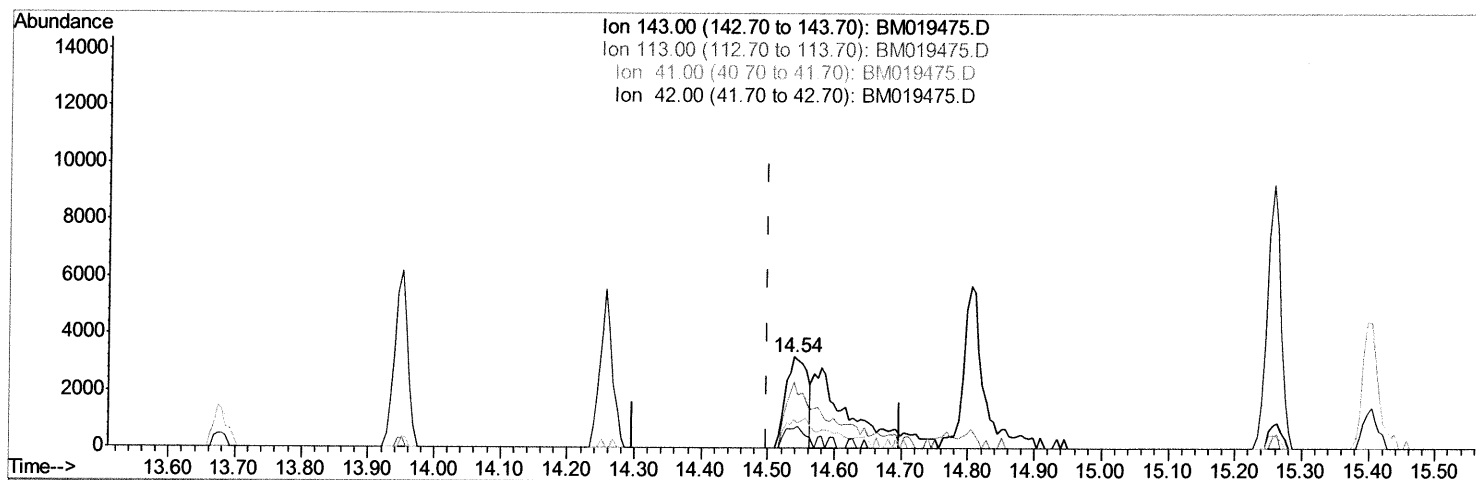
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TIC: BM019475.D

(52) 4-Nitrophenol-d4 (S)

14.539min (+0.041) 1.78ng/ul

response 7315

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	72.09
41.00	32.60	31.44
42.00	22.30	20.59

Quantitation Report (Qedit)

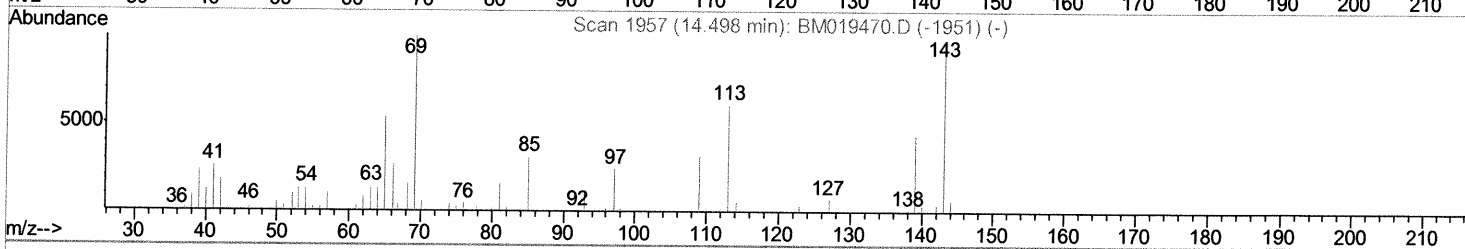
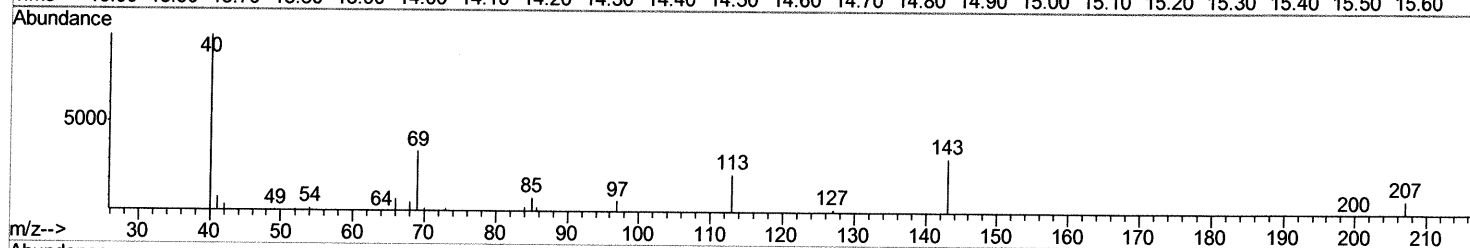
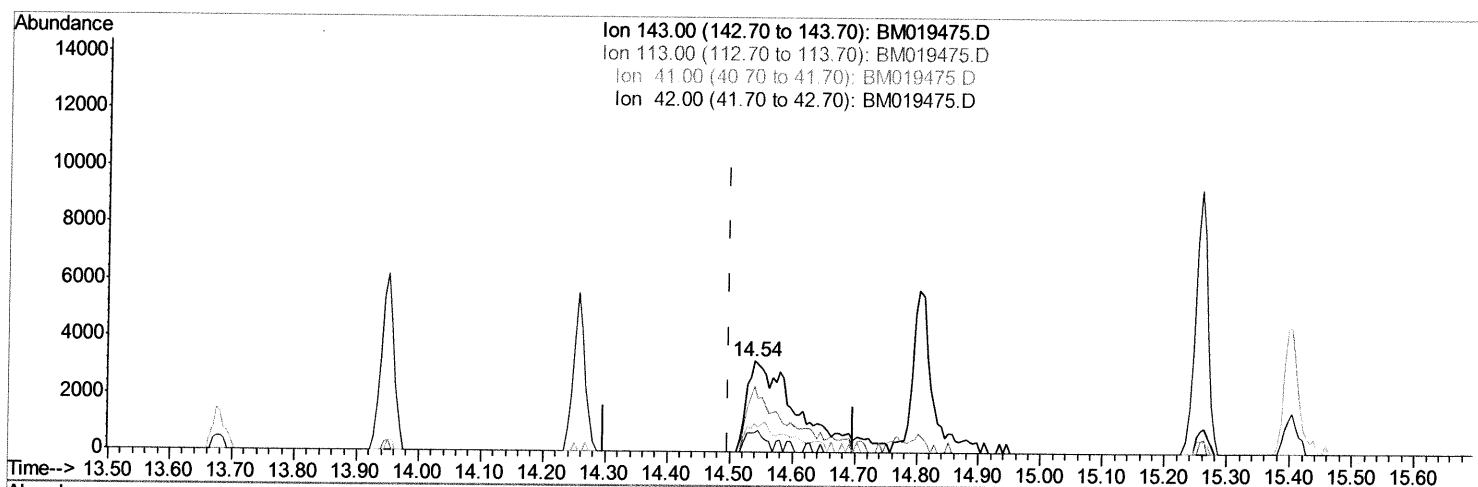
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Manual Integrations
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 Response via : Initial Calibration



TIC: BM019475.D

(52) 4-Nitrophenol-d4 (S)

14.539min (+0.041) 4.24ng/ul m 50 4/2/19

response 17369

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	72.09
41.00	32.60	31.44
42.00	22.30	20.59

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
 Data File : BM019475.D
 Acq On : 26 Mar 2019 18:03
 Operator : JU/SJ
 Sample : K2069-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 C09D6

Manual Integrations
 APPROVED

Sohil
 3/27/2019 5:46:29 PM

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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	153476	20.00	ng/ul	0.00
18) Naphthalene-d8	10.38	136	608011	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	324090	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	679901	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	712604	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	820428	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.15	96	17193	4.39	ng/uL	0.00
5) Phenol-d5	6.79	99	87322	6.39	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	237632	25.97	ng/ul	0.00
9) 2-Chlorophenol-d4	7.13	132	255500	24.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	155662	15.17	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	139075	32.07	ng/ul	0.00
22) 2-Nitrophenol-d4	9.48	143	157721	32.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.01	165	256994	28.22	ng/ul	0.00
29) 4-Chloroaniline-d4	10.58	131	14438	1.32	ng/ul	0.04
44) Dimethylphthalate-d6	13.68	166	843311	32.25	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	1078240	32.97	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	17369m)	4.24	ng/ul	0.04
58) Fluorene-d10	15.26	176	746558	33.14	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	116882	26.00	ng/ul	0.00
71) Anthracene-d10	17.12	188	1150195	35.25	ng/ul	0.00
79) Pyrene-d10	19.42	212	1210329	36.77	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	1398895	32.16	ng/ul	0.00
Target Compounds						
37) 1,2,4,5-Tetrachlorobenzene	12.43	216	56031	5.433	ng/ul	97
73) 1,2,3,4-Tetrachlorobenzene	13.05	216	262697	26.650	ng/uL	97
74) Pentachlorobenzene	14.57	250	15948	1.547	ng/uL	92

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

SJ 4/4/19