

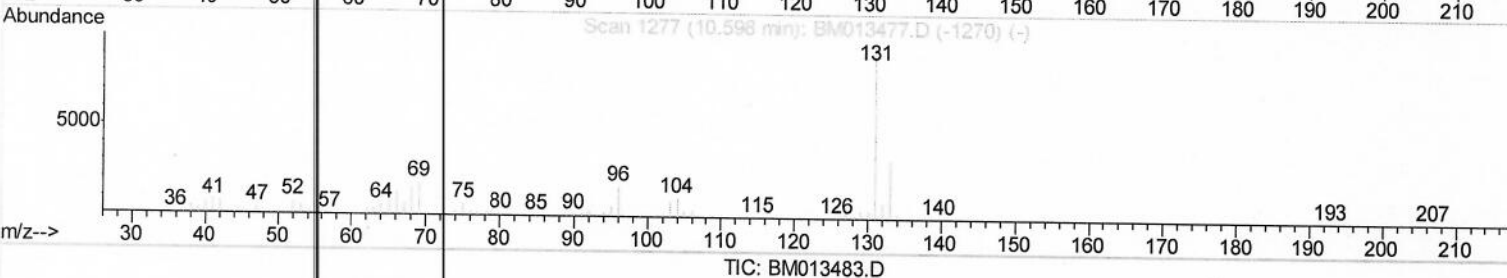
Quantitation Report (Qedit)

ALS Vial : 40	Sample Multiplier: 1
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Response via : Initial Calibration

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0.00	0.00	0.00
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Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\

Quantitation Report (Qedit)

Data File : BM013483.D

Acq On : 16 Dec 2017 18:58

Operator : SJ/JU

Sample : I6775-10DL 30X

Misc :

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 18 03:14:39 2017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M

Quant Title : SVDA CALIBRATION

QLast Update : Mon Dec 18 03:13:05 2017

Response via : Initial Calibration

Instrument :

BNA_M

Client Sampled :

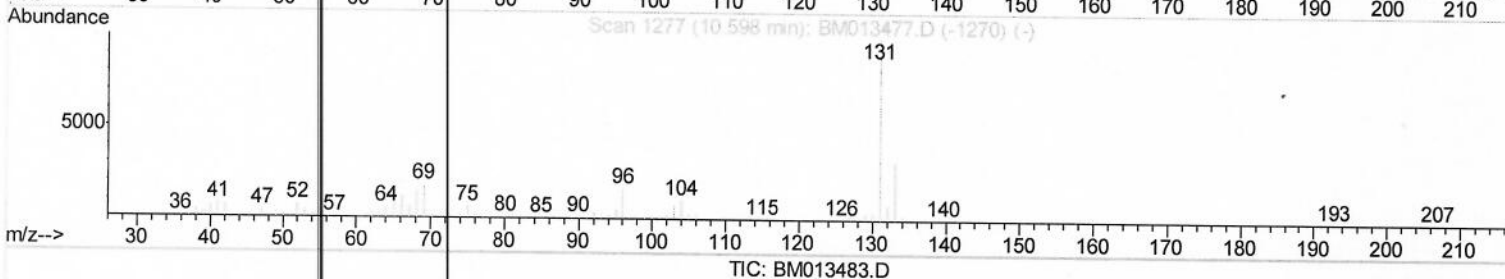
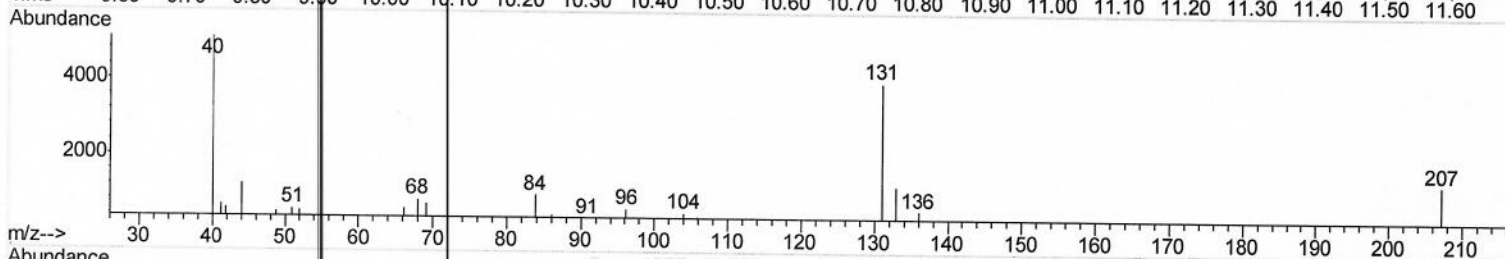
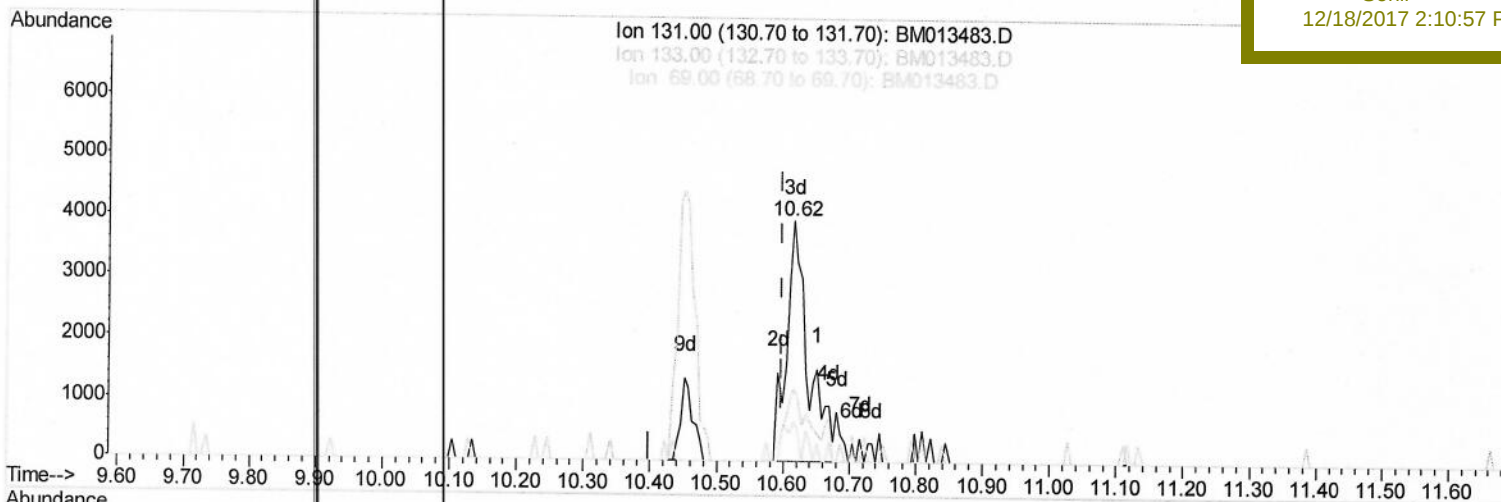
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Manual Integrations

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TIC: BM013483.D

(29) 4-Chloroaniline-d4 (S)

10.616min (+0.017) 0.54ng/ul m 15J12125117

response 9494

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	30.49
69.00	24.30	16.95#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\
 Quantitation Report (QT Reviewed)
 Data File : BM013483.D
 Acq On : 16 Dec 2017 18:58
 Operator : SJ/JU
 Sample : I6775-10DL 30X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9A22DL

Quant Time: Dec 18 03:41:44 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 18 03:13:05 2017
 Response via : Initial Calibration

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	244668	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	1102943	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	619047	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	1264227	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	903504	20.00	ng/ul	0.00
83) Perylene-d12	23.48	264	824392	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.86	99	11176	0.53	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.02	67	8442	0.70	ng/ul	0.00
9) 2-Chlorophenol-d4	7.22	132	9724	0.59	ng/ul	0.00
13) 4-Methylphenol-d8	8.40	113	8015	0.44	ng/ul	0.01
19) Nitrobenzene-d5	8.84	128	5644	0.65	ng/ul	0.01
22) 2-Nitrophenol-d4	9.55	143	5550	0.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.10	165	8238	0.43	ng/ul	0.02
29) 4-Chloroaniline-d4	10.62	131	9494m	0.54	ng/ul	0.02
43) Dimethylphthalate-d6	13.73	166	36140	0.66	ng/ul	0.00
46) Acenaphthylene-d8	14.01	160	42070	0.62	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.32	176	29848	0.63	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.17	188	44624	0.70	ng/ul	0.00
76) Pyrene-d10	19.46	212	38310	0.85	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.34	264	25964	0.62	ng/ul	0.00
Target Compounds						
68) Pentachlorophenol	16.72	266	348817	36.80	ng/ul	Qvalue 99

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

5J12125117

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BNA_M

ClientSampleId :

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Manual Integrations

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