

Quantitation Report (QT/LSC Reviewed)

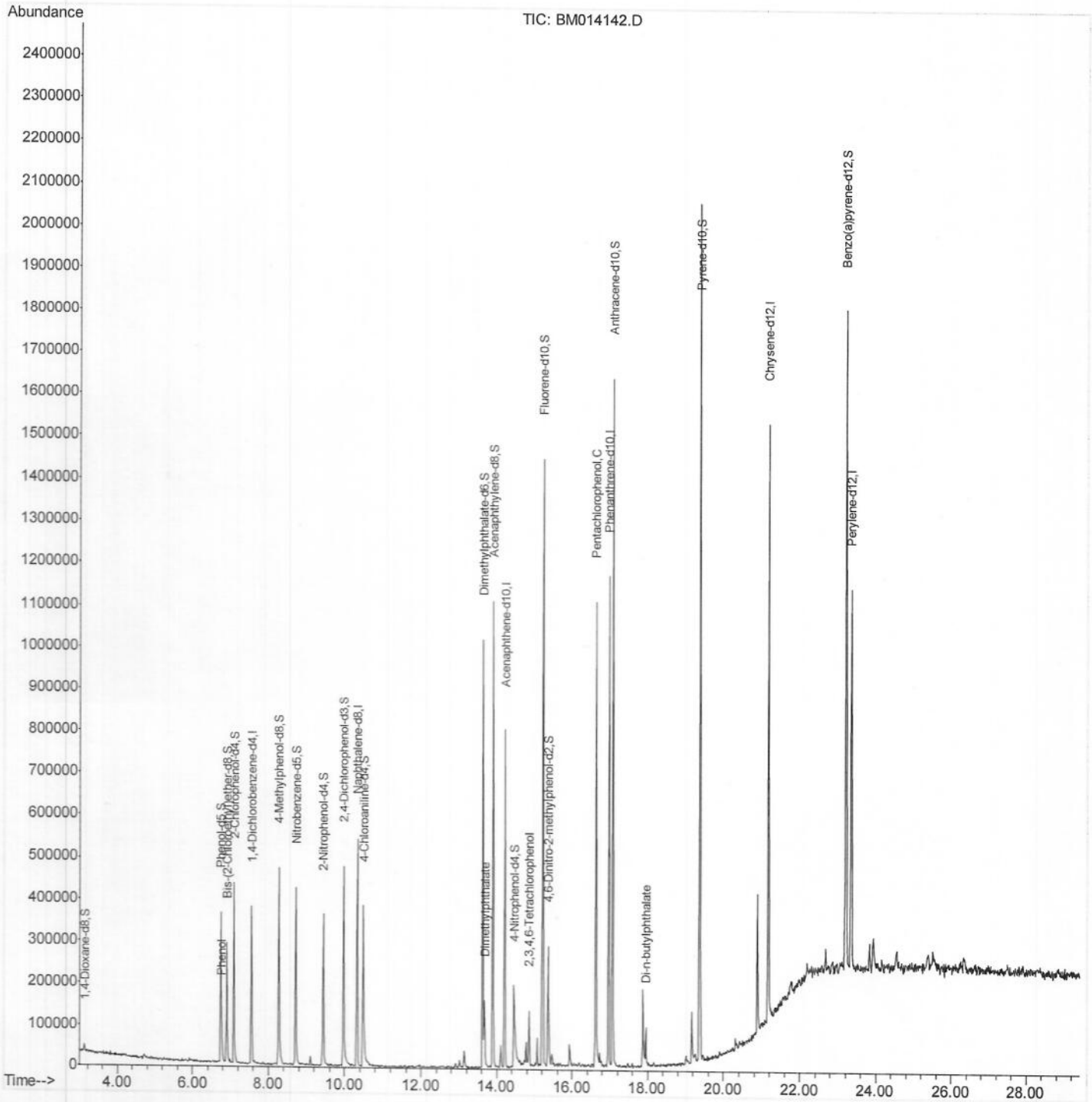
Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_M\Data\BM020218\
 Data File : BM014142.D
 Acq On : 03 Feb 2018 08:17
 Operator : SJ/JU
 Sample : J1339-13
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6D22

Manual Integrations
 APPROVED

Sohil
 2/5/2018 3:57:53 PM

Quant Time: Feb 05 13:05:35 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM013118.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Feb 03 04:17:00 2018
 Response via : Initial Calibration



Quantitation Report (Qedit)

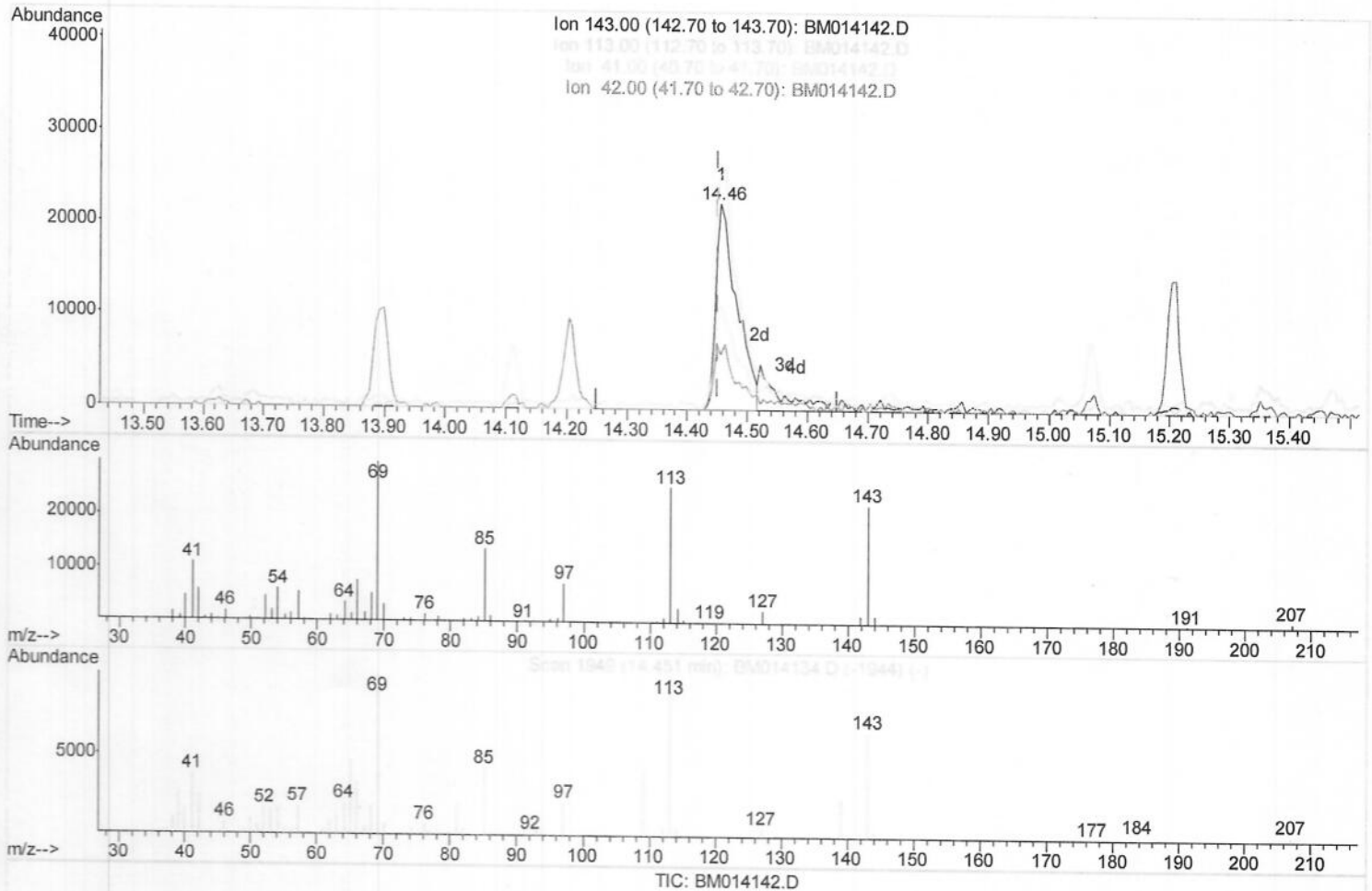
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(51) 4-Nitrophenol-d4 (S)

14.457min (+0.006) 19.56ng/ul

response 53639

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	113.83
41.00	34.80	49.08#
42.00	20.50	26.88#

Quantitation Report (Qedit)

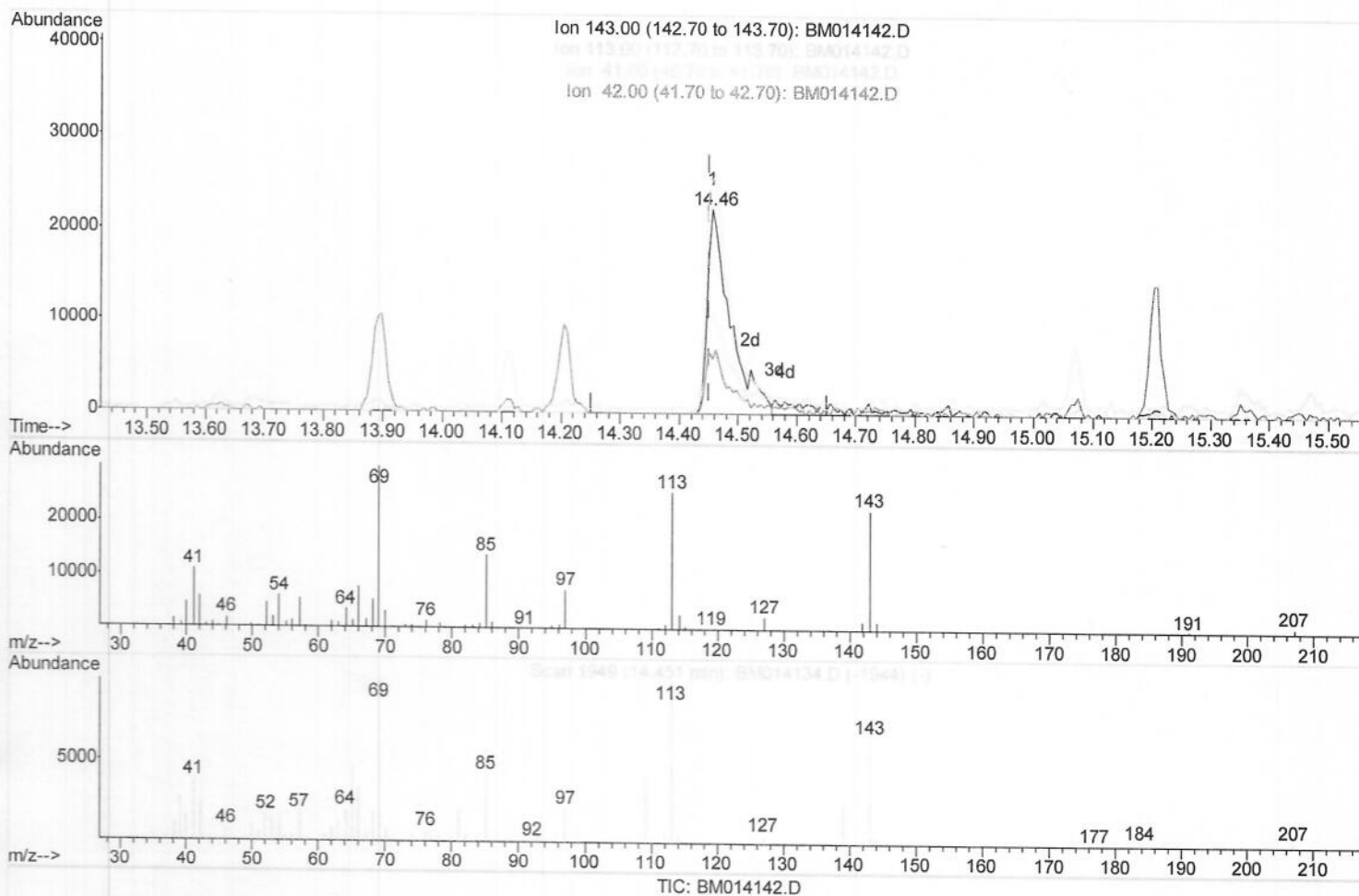
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(51) 4-Nitrophenol-d4 (S)

14.457min (+0.006) 21.97ng/ul m

response 60257

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	113.83
41.00	34.80	49.08#
42.00	20.50	26.88#

JU
02/03/18

Quantitation Report (Qedit)

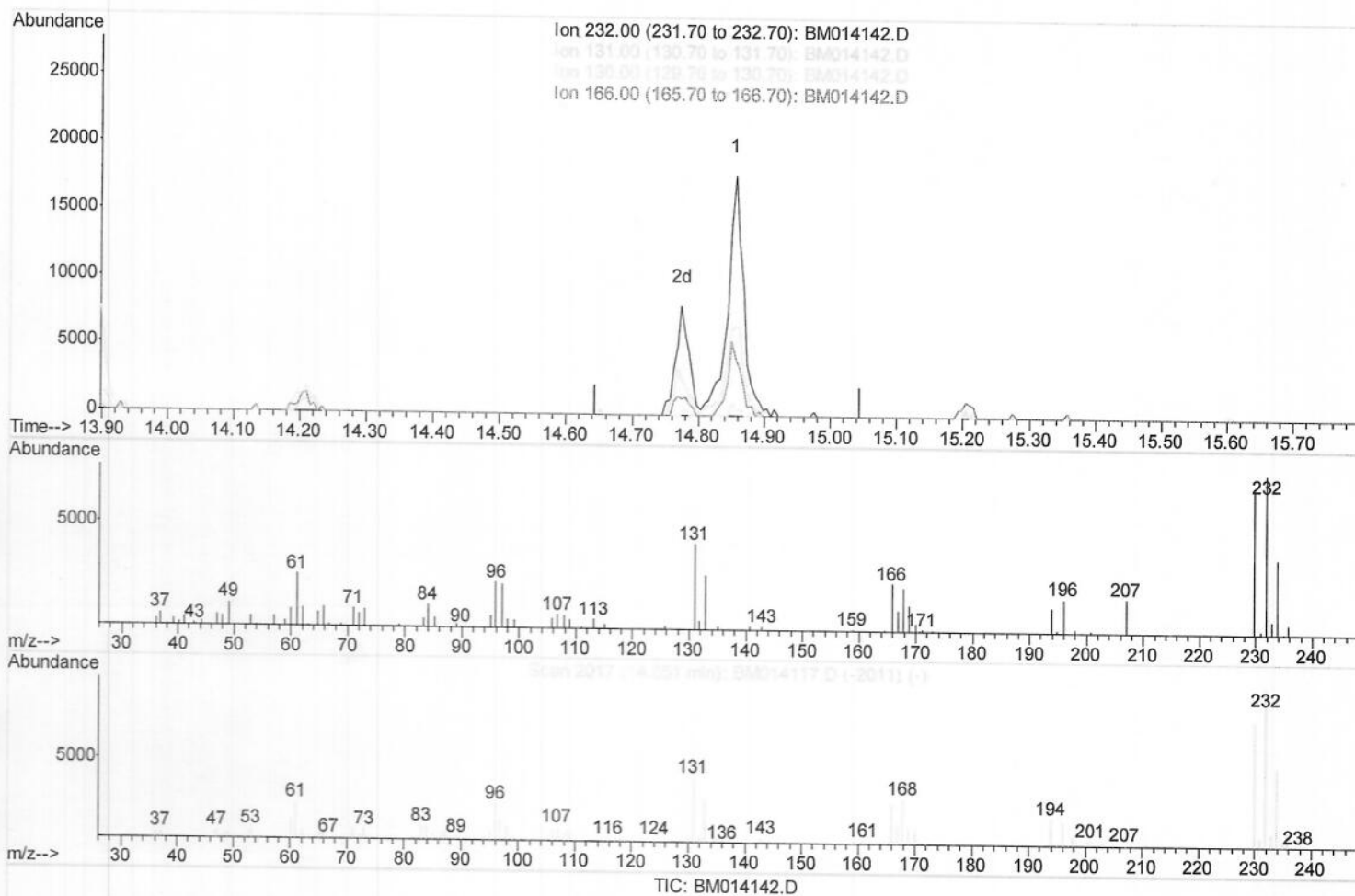
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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 F6D22

Manual Integrations
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Sohil
 2/5/2018 3:57:53 PM

Quant Time: Feb 05 06:11:43 2018
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(55) 2,3,4,6-Tetrachlorophenol

14.845min (-14.845) 0.00ng/ul d

response 0

Ion	Exp%	Act%
232.00	100	0.00
131.00	36.20	0.00
130.00	2.50	0.00
166.00	26.80	0.00

Quantitation Report (Qedit)

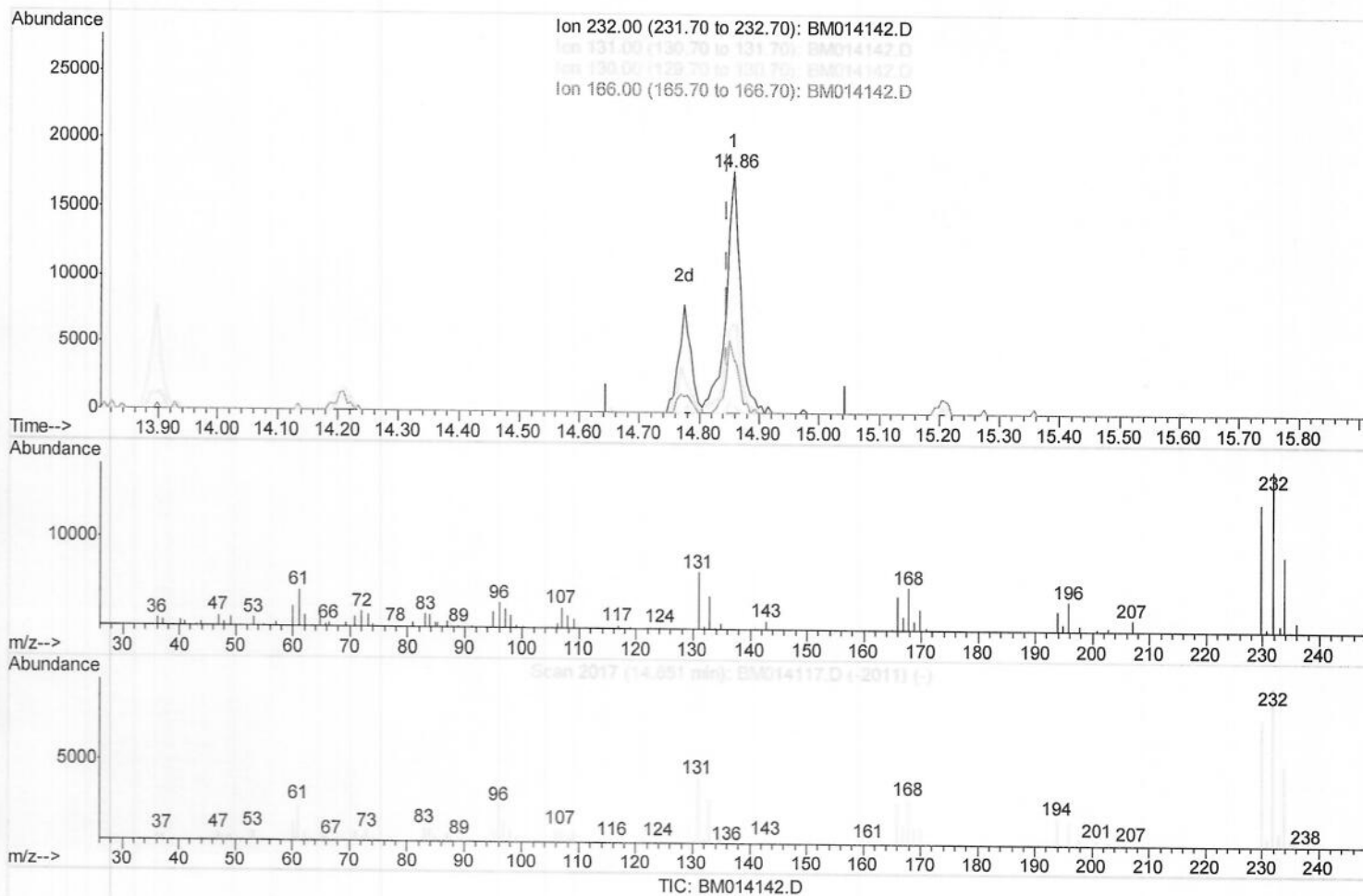
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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

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(55) 2,3,4,6-Tetrachlorophenol

14.857min (+0.012) 5.64ng/ul m

response 29784

Ion	Exp%	Act%
232.00	100	100
131.00	36.20	36.73
130.00	2.50	2.23
166.00	26.80	22.76

JV
 02/03/18

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	93400	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	384527	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.20	164	239831	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.96	188	610746	20.00	ng/ul	0.00
75) Chrysene-d12	21.17	240	696097	20.00	ng/ul	0.00
83) Perylene-d12	23.34	264	686310	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.12	96	10041	4.93	ng/uL	0.00
5) Phenol-d5	6.74	99	187429	26.70	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.90	67	112185	26.49	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	157579	28.35	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	161233	27.43	ng/ul	0.00
19) Nitrobenzene-d5	8.71	128	83911	30.23	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	90185	29.36	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	165306	25.58	ng/ul	0.00
29) 4-Chloroaniline-d4	10.48	131	196671	35.11	ng/ul	0.00
43) Dimethylphthalate-d6	13.63	166	621365	31.22	ng/ul	0.00
46) Acenaphthylene-d8	13.89	160	761864	31.55	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	60257m	21.97	ng/ul	0.00
57) Fluorene-d10	15.21	176	542297	30.44	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.36	200	70007	18.67	ng/ul	0.00
70) Anthracene-d10	17.06	188	862858	29.10	ng/ul	0.00
76) Pyrene-d10	19.37	212	1019287	30.41	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.21	264	1007173	29.90	ng/ul	0.00
Target Compounds						
6) Phenol	6.77	94	13678	1.948	ng/ul	93
44) Dimethylphthalate	13.67	163	94165	4.908	ng/ul#	95
55) 2,3,4,6-Tetrachlorophenol	14.86	232	29784m	5.639	ng/ul	
68) Pentachlorophenol	16.62	266	190015	47.515	ng/ul	99
73) Di-n-butylphthalate	17.94	149	60945	1.712	ng/ul	99

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