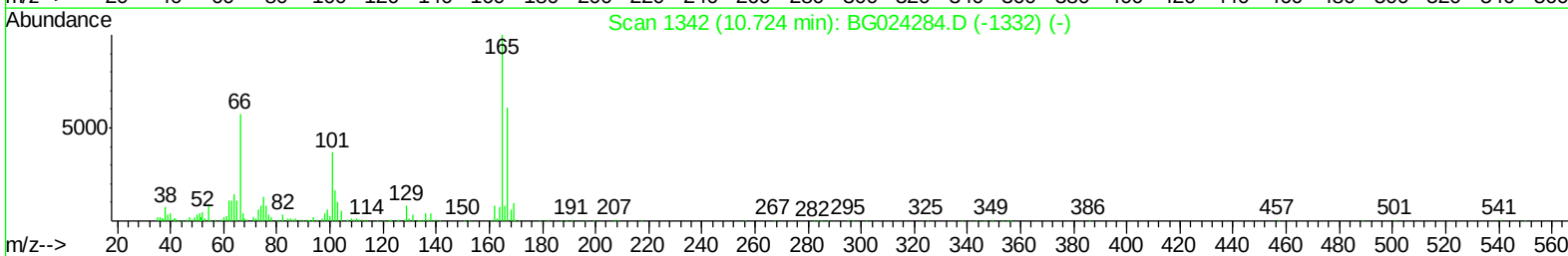
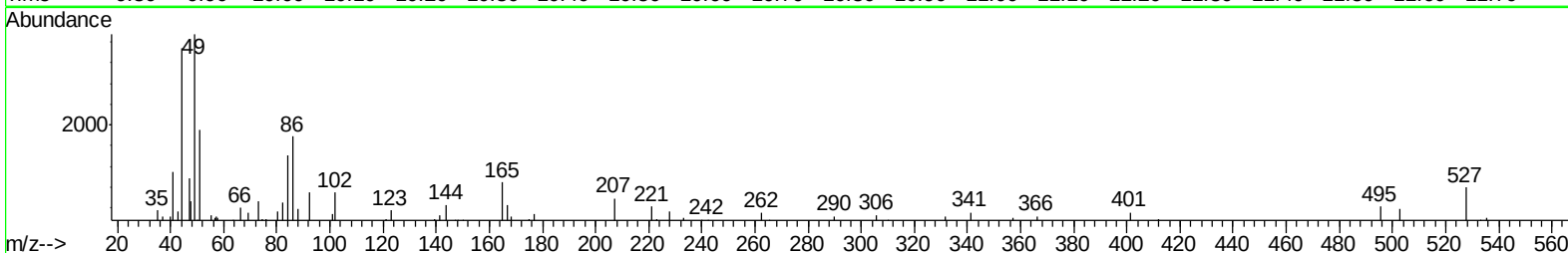
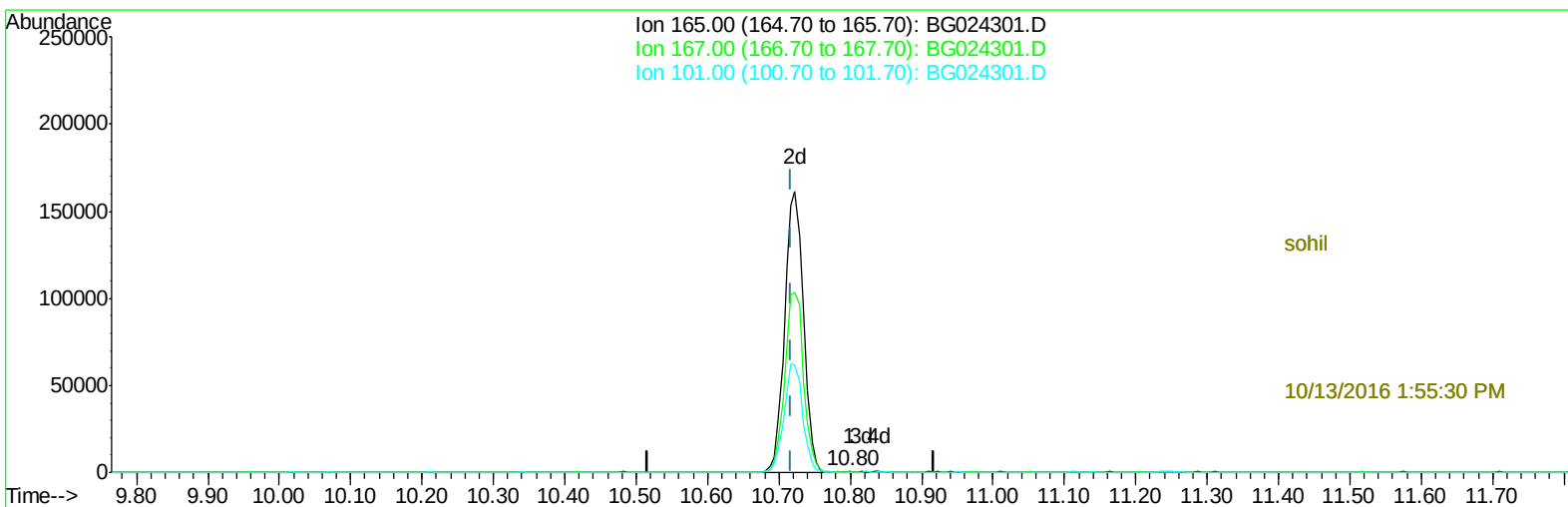


Data Path : Z:\HPCHEM1\BNA_G\Data\BG101216\
 Data File : BG024301.D
 Acq On : 12 Oct 2016 14:58
 Operator : UM/SJ
 Sample : H5145-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDN94

Manual Integrations
 APPROVED

Quant Time: Oct 12 15:35:24 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 12:34:51 2016
 Response via : Initial Calibration



TIC: BG024301.D

(26) 2,4-Dichlorophenol-d3 (S)

10.799min (+0.082) 0.06ng/ul

response 693

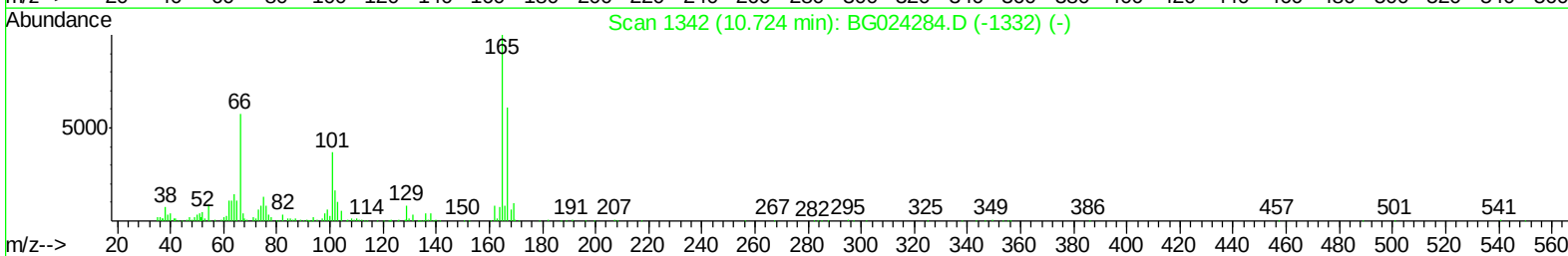
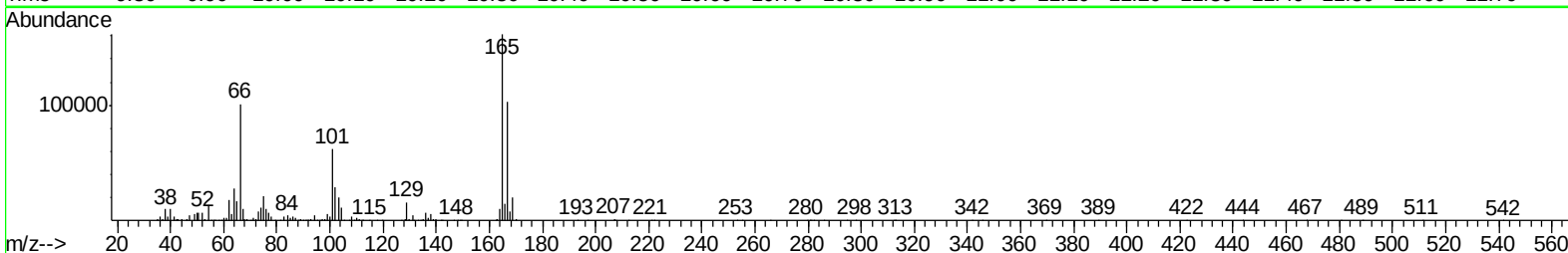
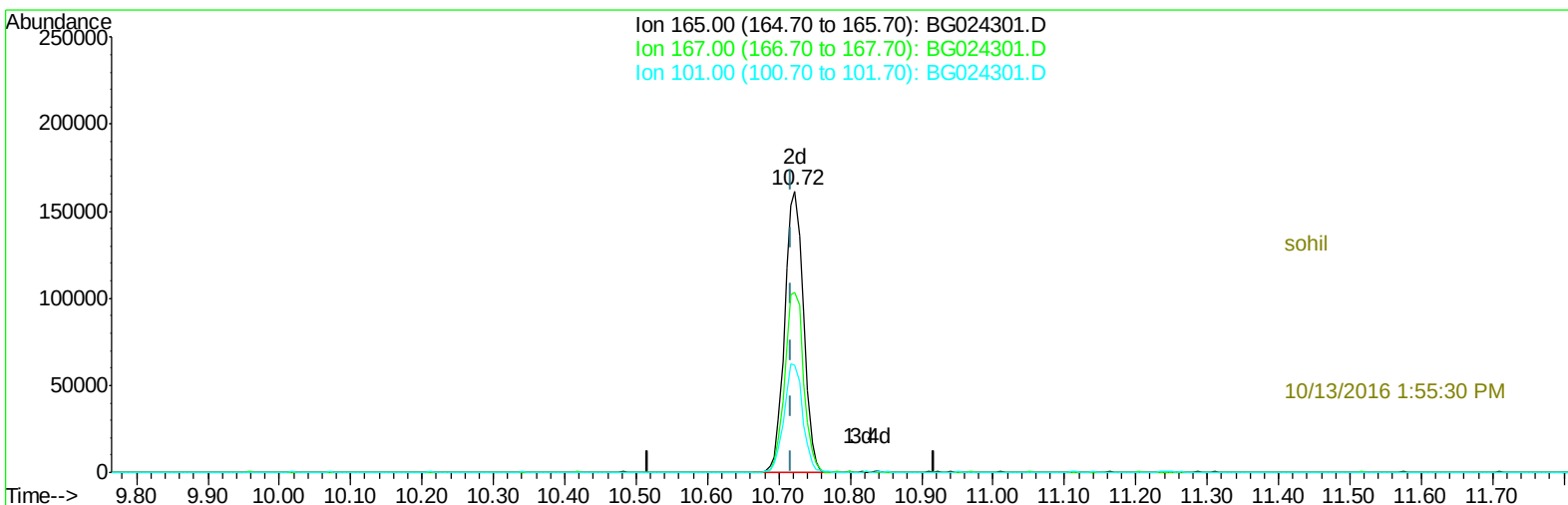
Ion	Exp%	Act%
165.00	100	100
167.00	58.70	50.23
101.00	29.80	31.04
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG101216\
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Sample : H5145-06
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Instrument :
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Manual Integrations
APPROVED

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TIC: BG024301.D

(26) 2,4-Dichlorophenol-d3 (S)

10.723min (+0.006) 27.52ng/ul m

response 295882

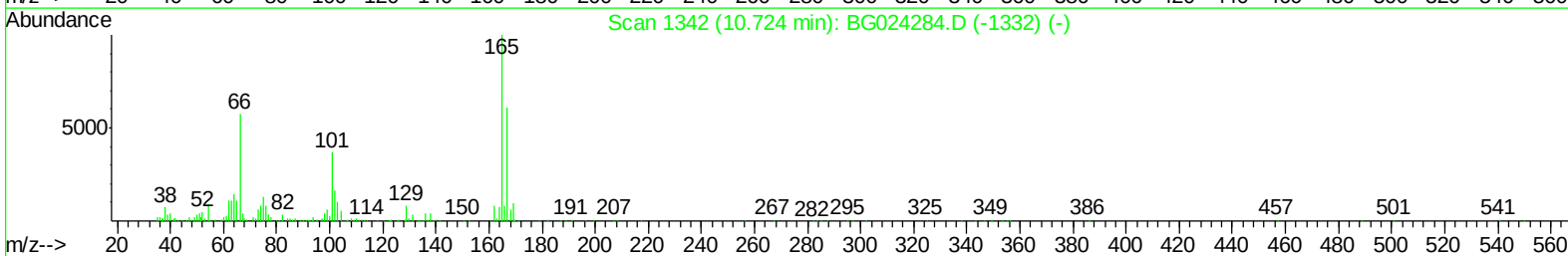
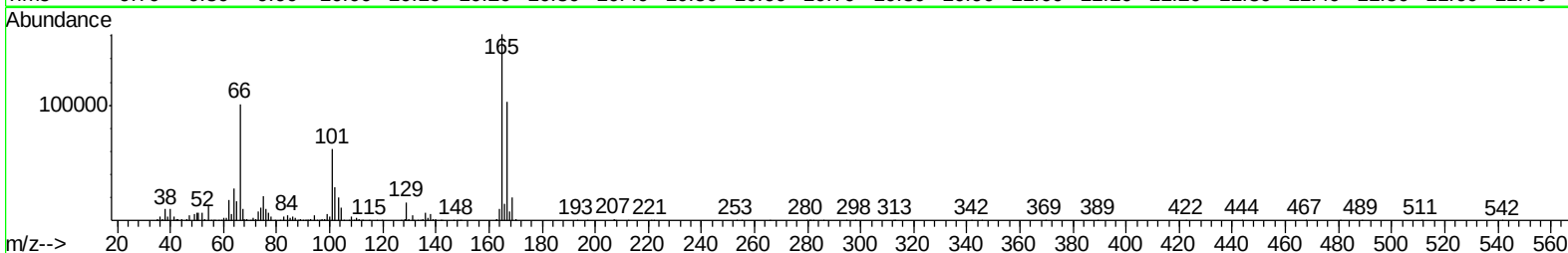
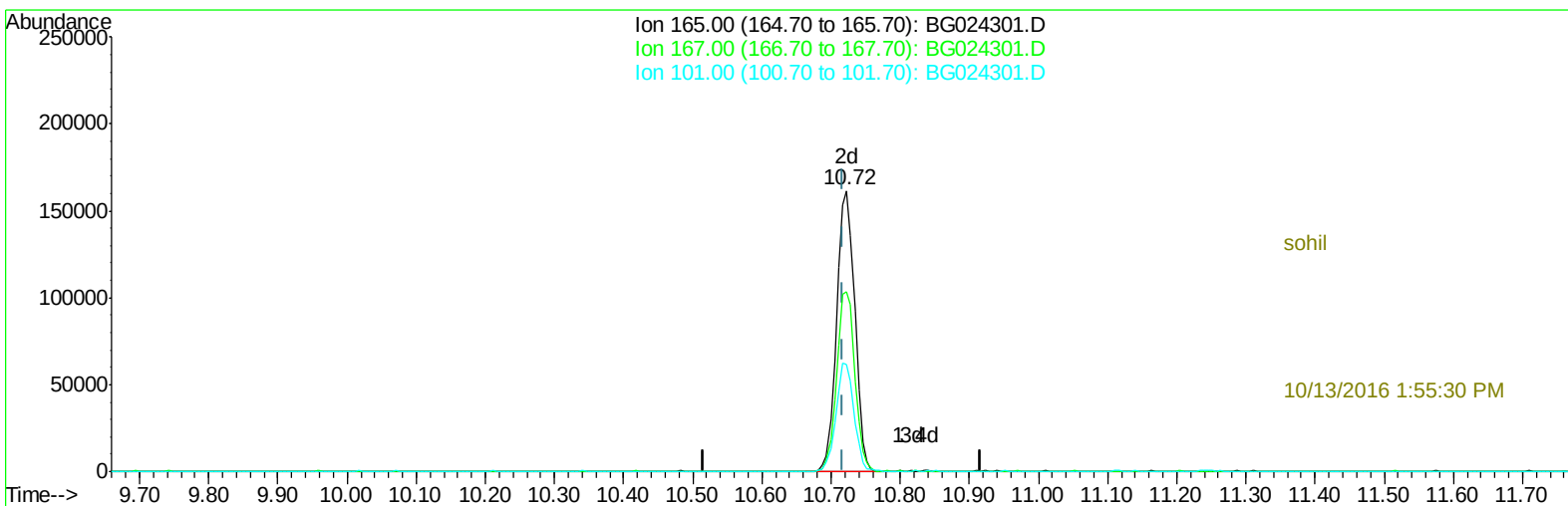
Ion	Exp%	Act%
165.00	100	100
167.00	58.70	63.94
101.00	29.80	38.12#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG101216\
Data File : BG024301.D
Acq On : 12 Oct 2016 14:58
Operator : UM/SJ
Sample : H5145-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Manual Integrations
APPROVED

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TIC: BG024301.D

(26) 2,4-Dichlorophenol-d3 (S)

10.723min (+0.006) 27.52ng/ul m

response 295882

Ion	Exp%	Act%
165.00	100	100
167.00	58.70	63.94
101.00	29.80	38.12#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG101216\
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Manual Integrations
 APPROVED

Quant Time: Oct 12 15:36:27 2016
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	131484	20.00	ng/ul	0.00
18) Naphthalene-d8	11.12	136	621812	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.91	164	515505	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1069183	20.00	ng/ul	0.00
75) Chrysene-d12	21.94	240	1188118	20.00	ng/ul	0.00
83) Perylene-d12	25.39	264	1223296	20.00	ng/ul	0.00

sohil

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.64	96	13005	4.70	ng/uL	0.00
5) Phenol-d5	7.41	99	335577	27.35	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	236573	28.53	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	238397	28.22	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	262947	26.51	ng/ul	0.00
19) Nitrobenzene-d5	9.47	128	131442	29.79	ng/ul	0.00
22) 2-Nitrophenol-d4	10.19	143	156095	31.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	295882m	27.52	ng/ul	0.00
29) 4-Chloroaniline-d4	11.25	131	381787	33.79	ng/ul	0.00
43) Dimethylphthalate-d6	14.29	166	1102258	30.62	ng/ul	0.00
46) Acenaphthylene-d8	14.60	160	1260433	30.57	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	172432	27.02	ng/ul	0.00
57) Fluorene-d10	15.89	176	1035456	30.43	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	196727	30.62	ng/ul	0.00
70) Anthracene-d10	17.74	188	1544754	31.99	ng/ul	0.00
76) Pyrene-d10	20.01	212	1737538	31.33	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.15	264	1769764	32.16	ng/ul	0.00

10/13/2016 1:55:30 PM

Target Compounds					Qvalue
6) Phenol	7.44	94	26826	2.15 ng/ul	94
44) Dimethylphthalate	14.34	163	666107	18.87 ng/ul	97

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

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