

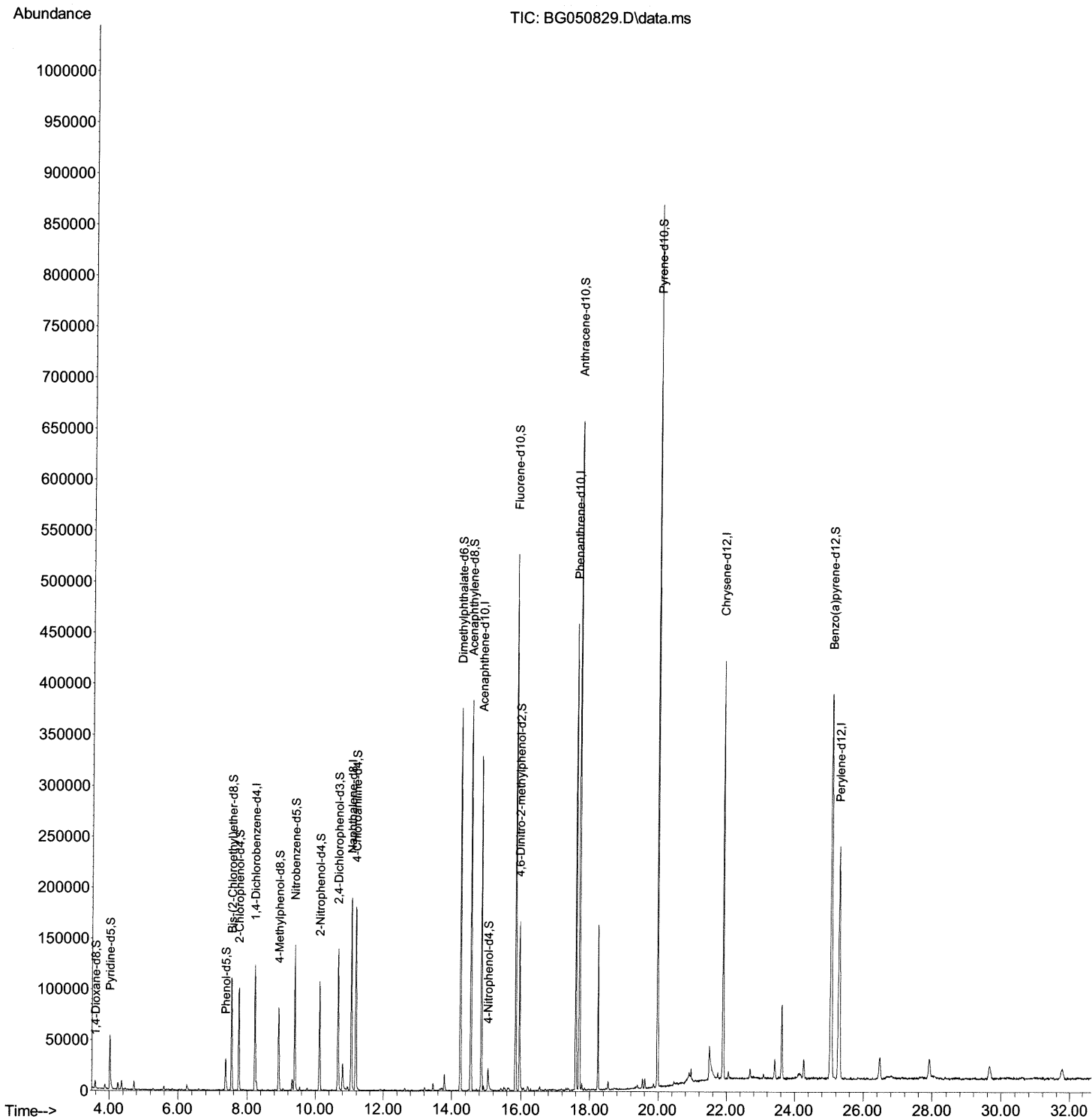
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
Data File : BG050829.D
Acq On : 2 Nov 2021 23:52
Operator : CG/JU
Sample : M4364-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BG340

Manual IntegrationsAPPROVED

Quant Time: Nov 03 01:34:46 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2021
Supervised By :mohammad ahmed 11/08/2021



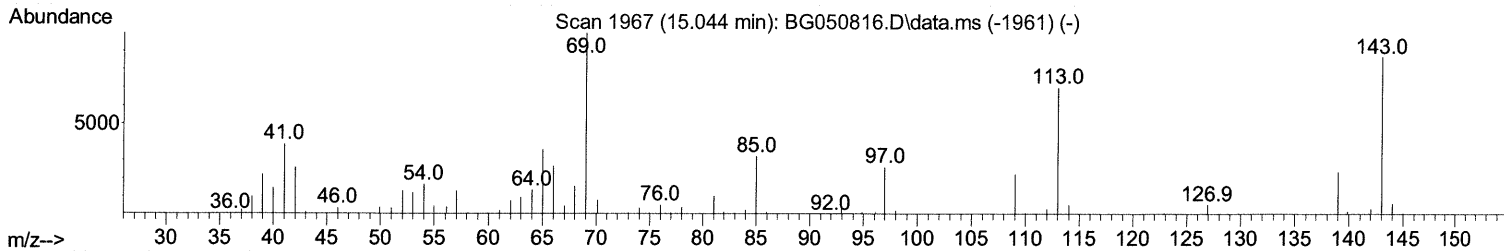
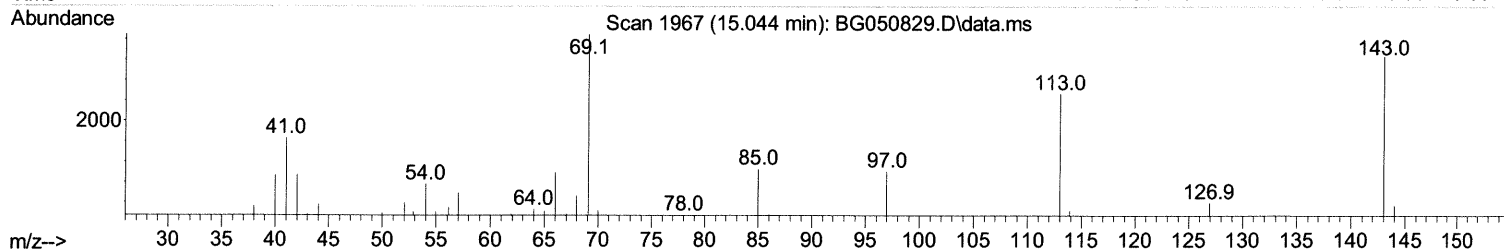
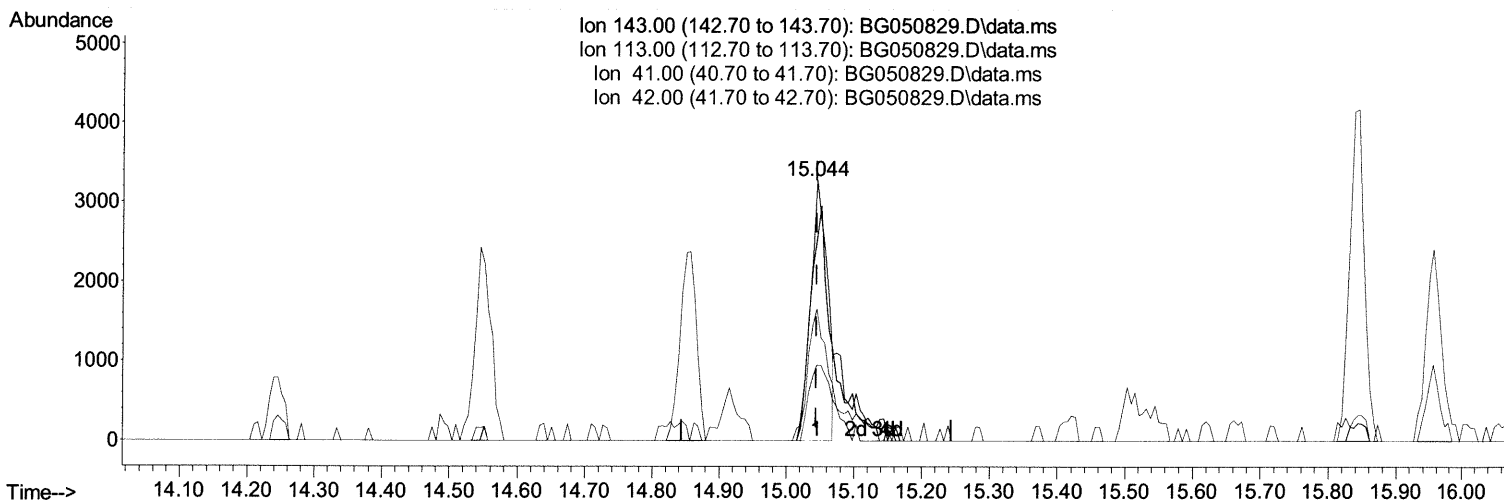
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TIC: BG050829.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.044min (+ 0.000) 3.50 ng/ul

response 5714

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	77.46
41.00	44.40	50.66
42.00	29.70	28.82

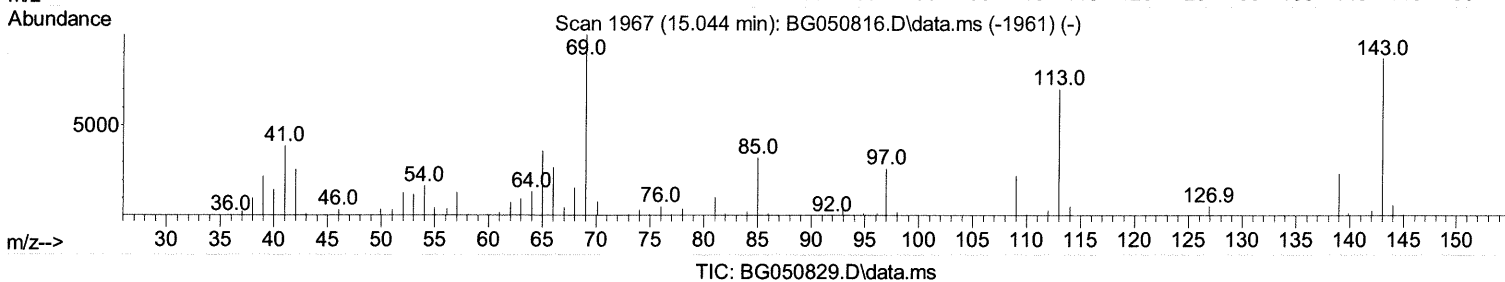
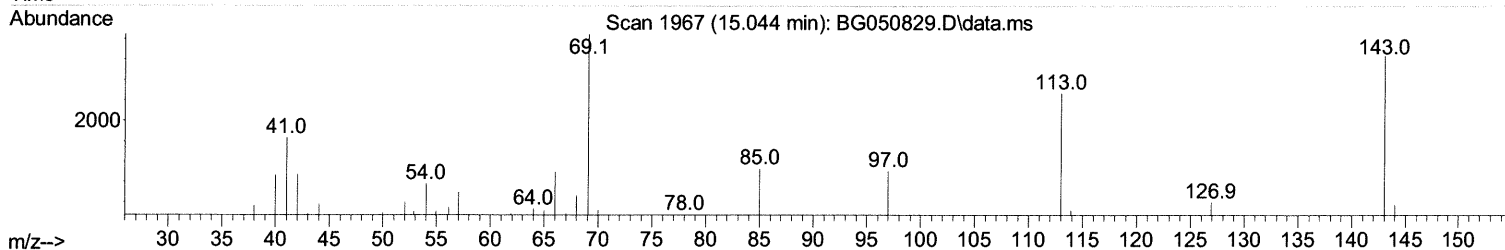
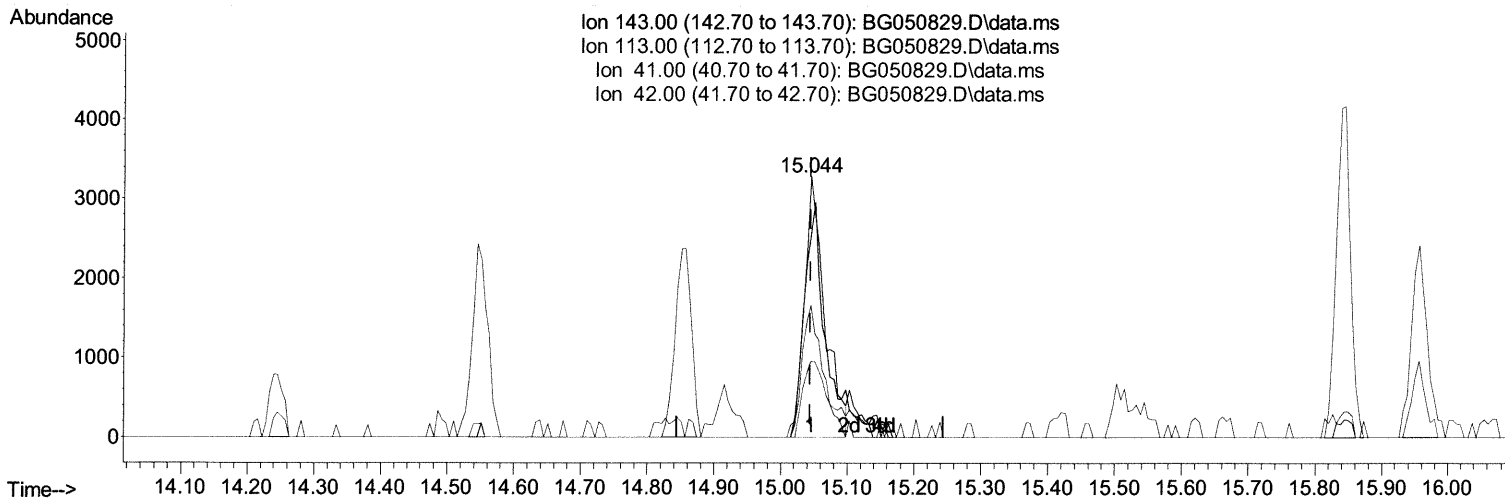
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(54) 4-Nitrophenol-d4 (S)

15.044min (+ 0.000) 4.27 ng/ul m 11041231

response 6971

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	77.46
41.00	44.40	50.66
42.00	29.70	28.82

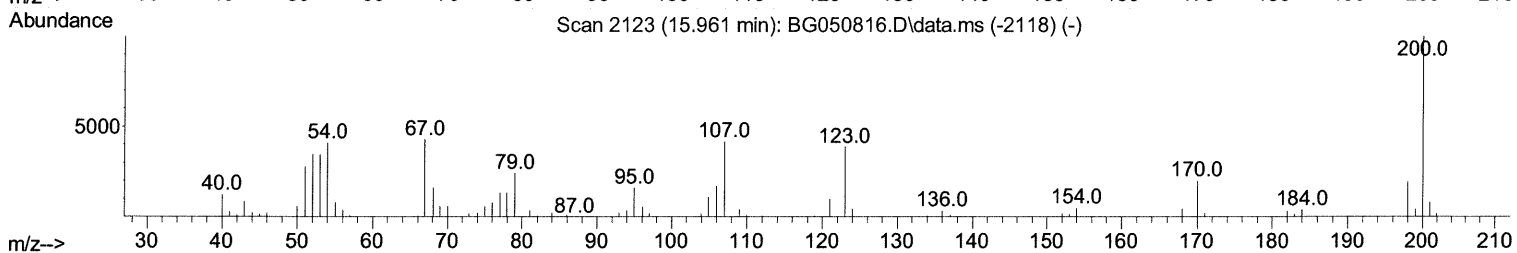
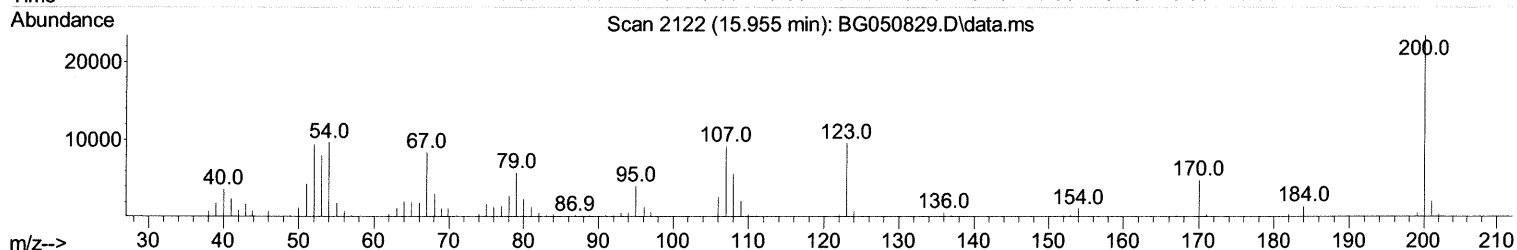
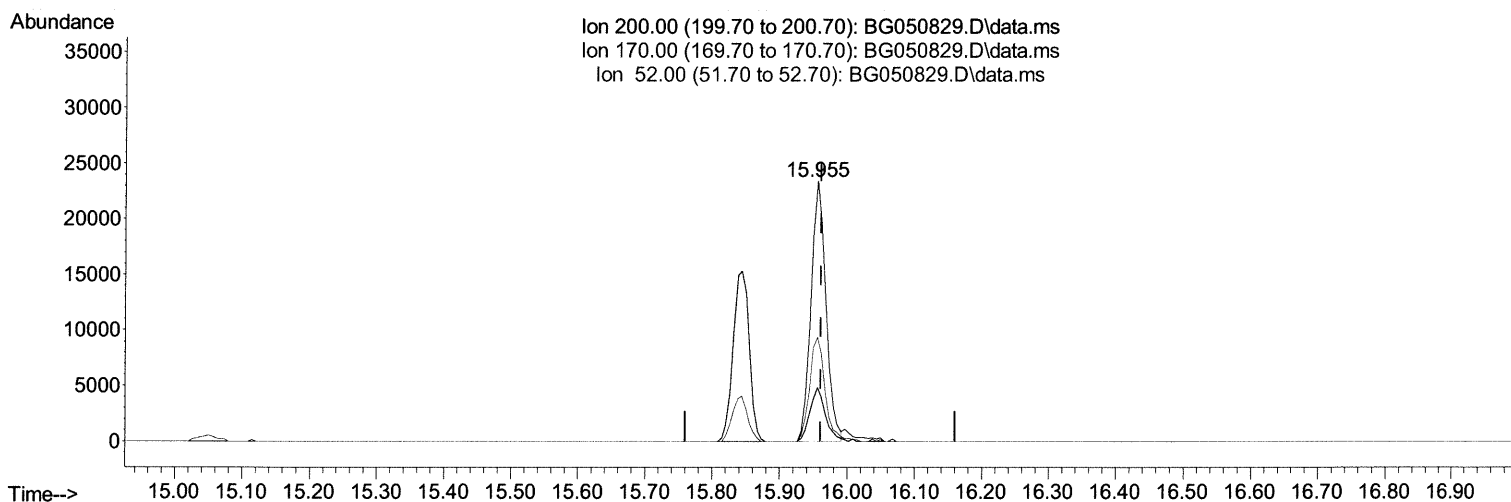
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TIC: BG050829.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.955min (-0.005) 21.15 ng/u1

response 35746

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	20.60
52.00	47.40	39.99
0.00	0.00	0.00

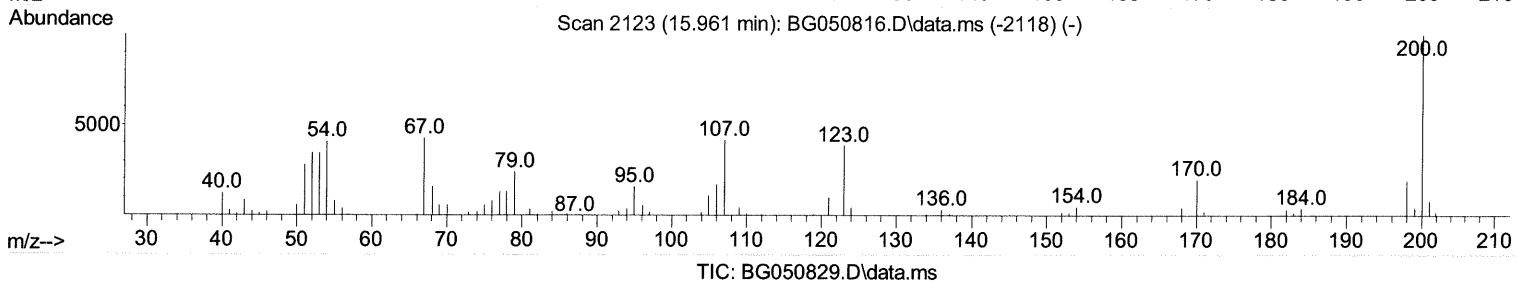
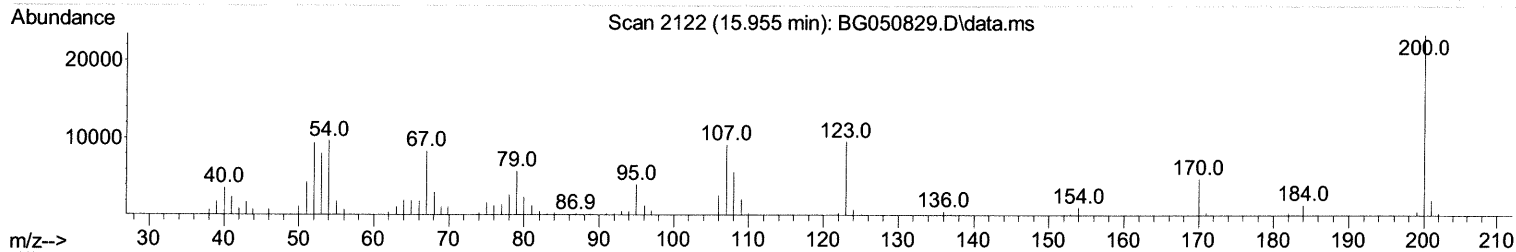
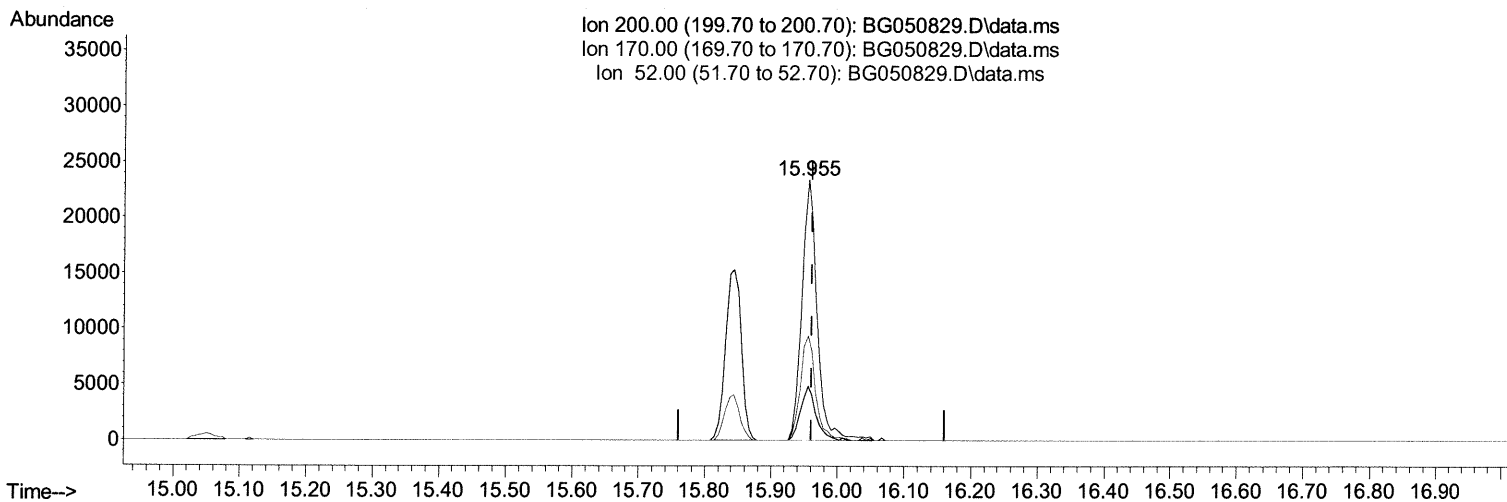
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Sample : M4364-08
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.955min (-0.005) 21.70 ng/ul m 11/04/21 JU

response 36674

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	20.60
52.00	47.40	39.99
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : M4364-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	33084	20.000	ng/ul	0.00
20) Naphthalene-d8	11.055	136	160327	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.856	164	117636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.600	188	278746	20.000	ng/ul	0.00
79) Chrysene-d12	21.895	240	246275	20.000	ng/ul	-0.01
88) Perylene-d12	25.297	264	228689	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.587	96	4247	4.143	ng/uL	0.00
4) Pyridine-d5	4.016	84	35467	11.566	ng/ul	0.00
7) Phenol-d5	7.371	99	17840	5.055	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	58900	25.834	ng/ul	0.00
11) 2-Chlorophenol-d4	7.759	132	45547	18.621	ng/ul	0.00
15) 4-Methylphenol-d8	8.928	113	32447	11.678	ng/ul	0.00
21) Nitrobenzene-d5	9.404	128	32502	23.855	ng/ul	0.00
24) 2-Nitrophenol-d4	10.127	143	33804	22.314	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	51478	20.172	ng/ul	0.00
31) 4-Chloroaniline-d4	11.184	131	102499	26.523	ng/ul	0.00
46) Dimethylphthalate-d6	14.245	166	242444	26.939	ng/ul	0.00
49) Acenaphthylene-d8	14.551	160	276530	24.662	ng/ul	0.00
54) 4-Nitrophenol-d4	15.044	143	6971m >	4.272	ng/ul >	0.00 11041214
60) Fluorene-d10	15.844	176	211719	26.556	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	36674m >	21.701	ng/ul >	0.00 11041214
73) Anthracene-d10	17.700	188	388260	29.461	ng/ul	0.00
81) Pyrene-d10	19.974	212	464923	29.228	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.062	264	386231	30.551	ng/ul	-0.01

Target Compounds Qvalue

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