

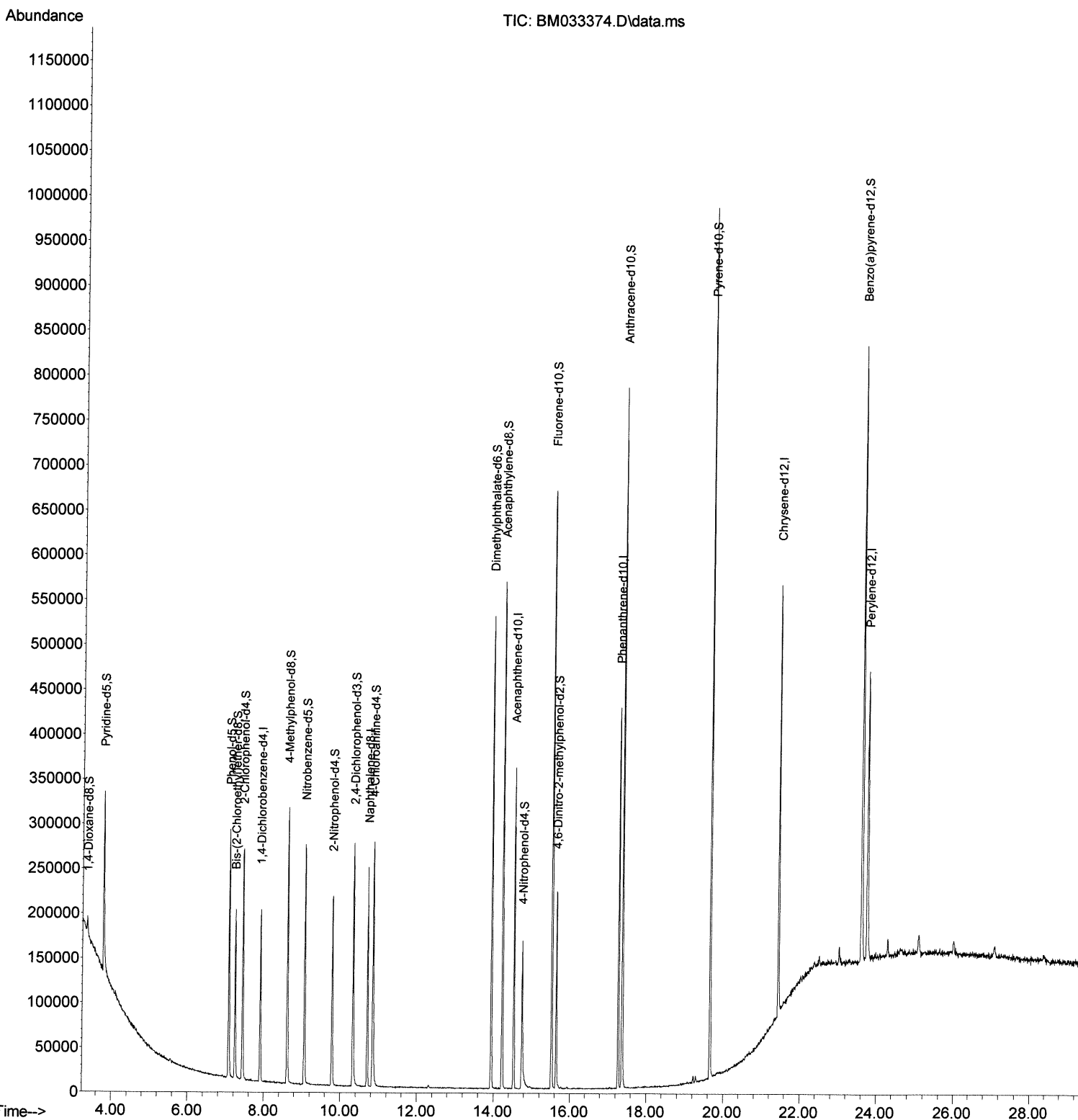
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\
Data File : BM033374.D
Acq On : 10 Dec 2021 01:47
Operator : CG/JU
Sample : PB141276BL
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK276

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021
Supervised By :mohammad ahmed 12/15/2021

Quant Time: Dec 10 04:54:51 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 13:25:37 2021
Response via : Initial Calibration



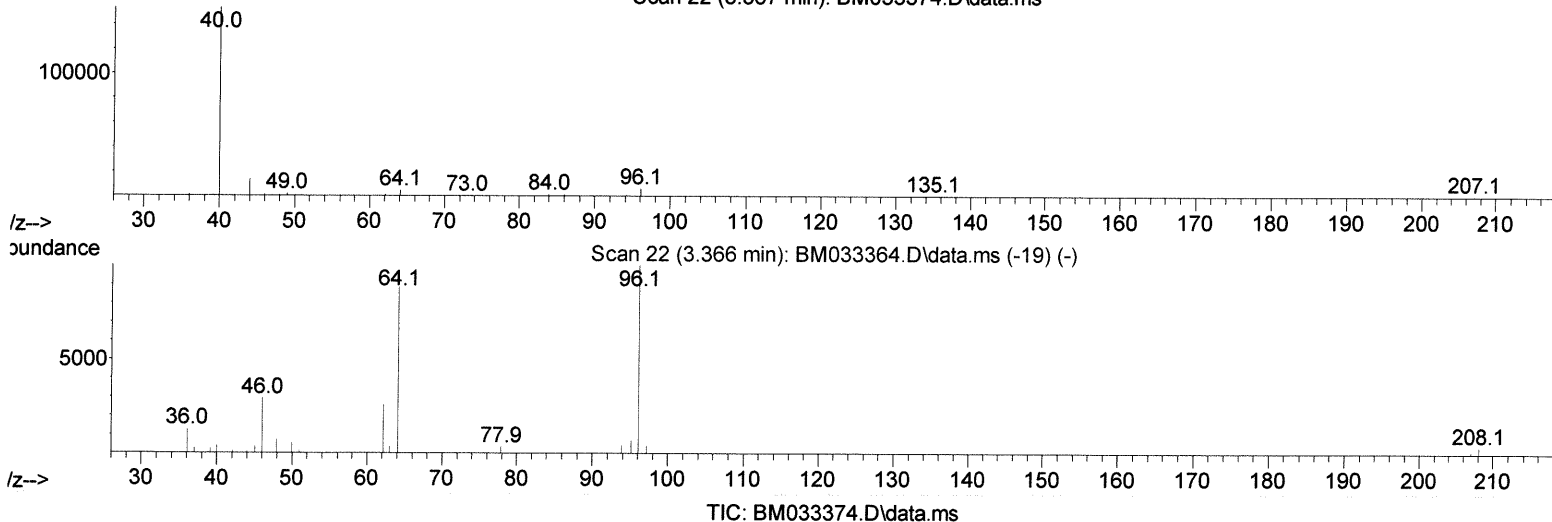
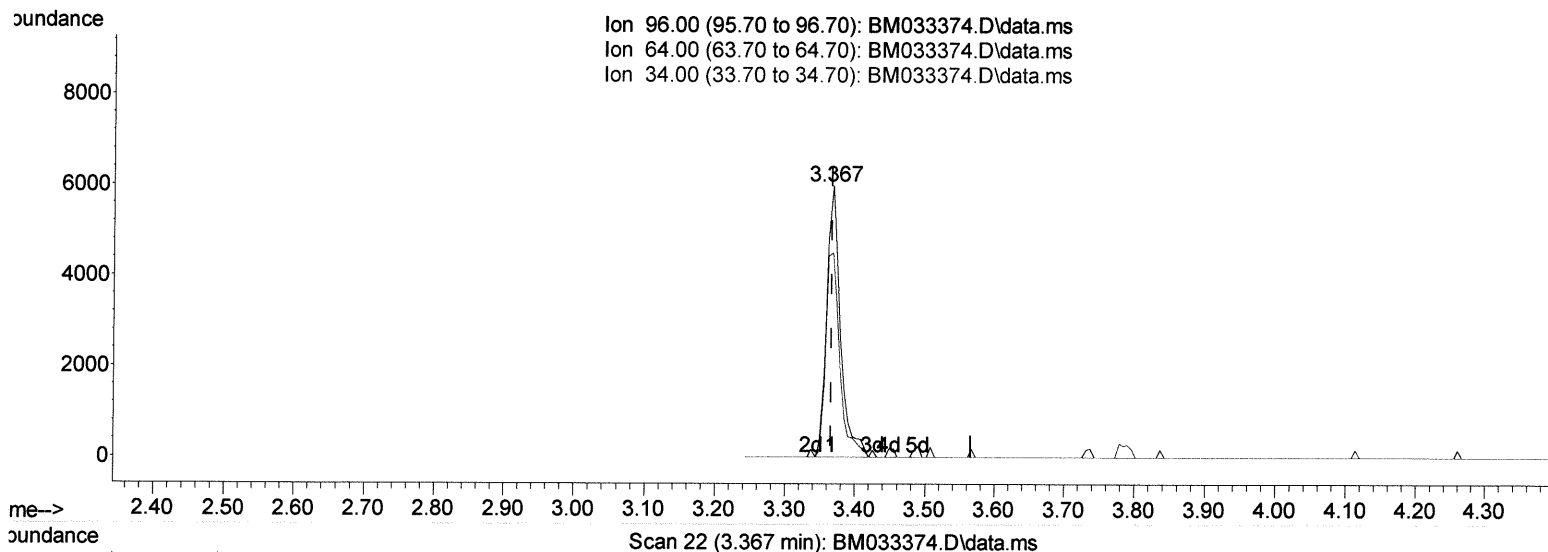
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(3) 1,4-Dioxane-d8 (S)

3.367min (+ 0.001) 6.69 ng/uL

response 8129

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	75.23
34.00	0.00	0.00
0.00	0.00	0.00

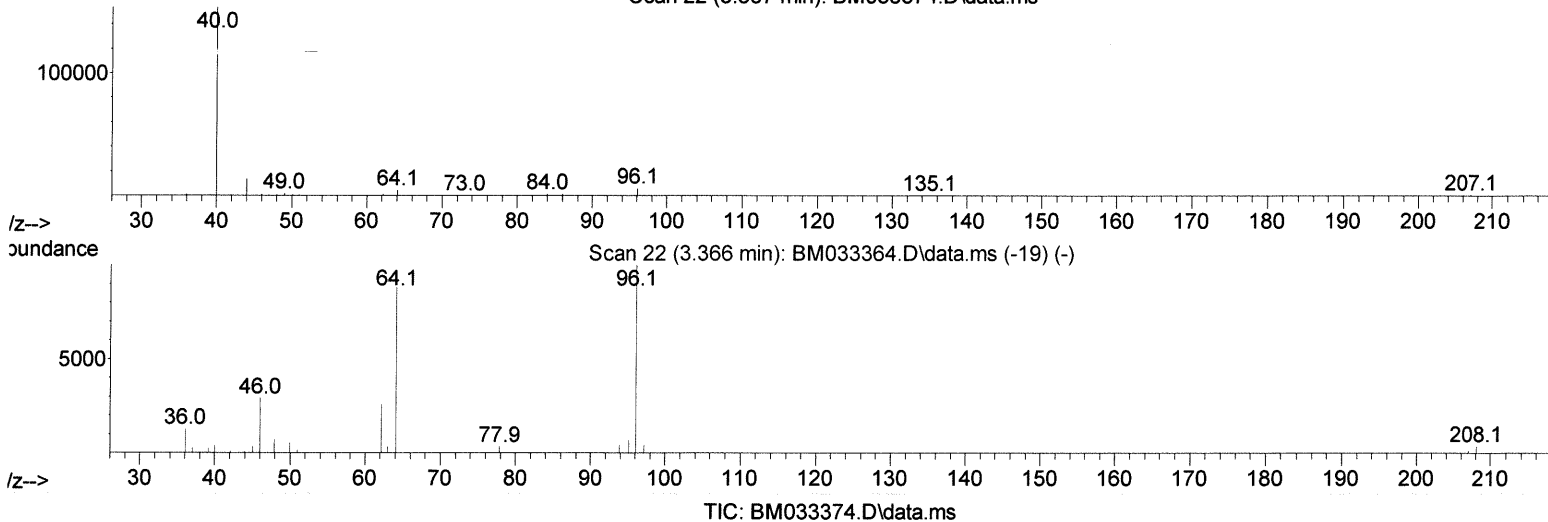
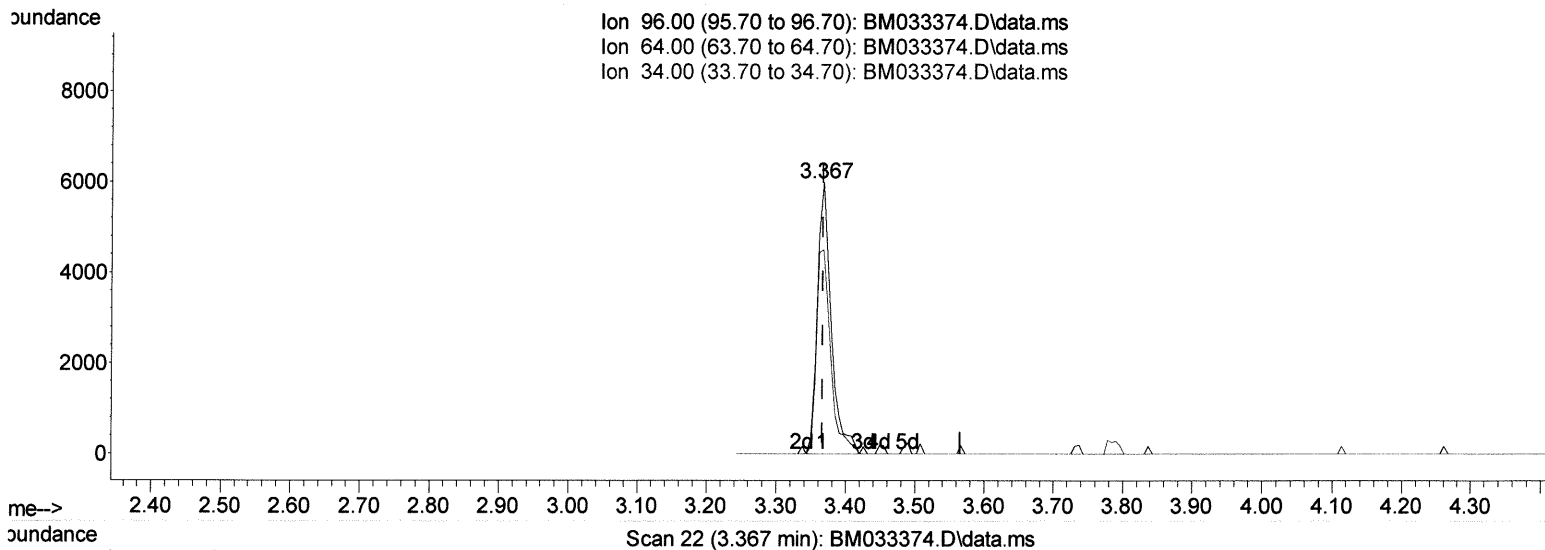
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(3) 1,4-Dioxane-d8 (S)

3.367min (+ 0.001) 6.87 ng/uL m

response 8339

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	75.23
34.00	0.00	0.00
0.00	0.00	0.00

Jyoti

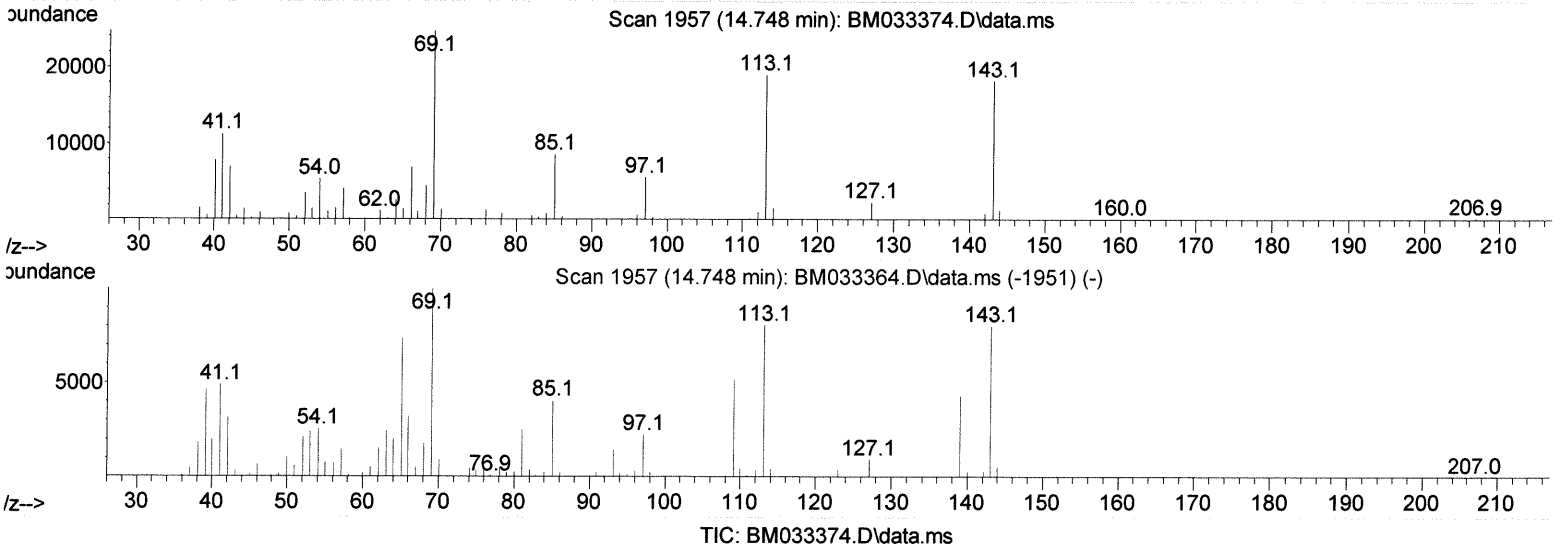
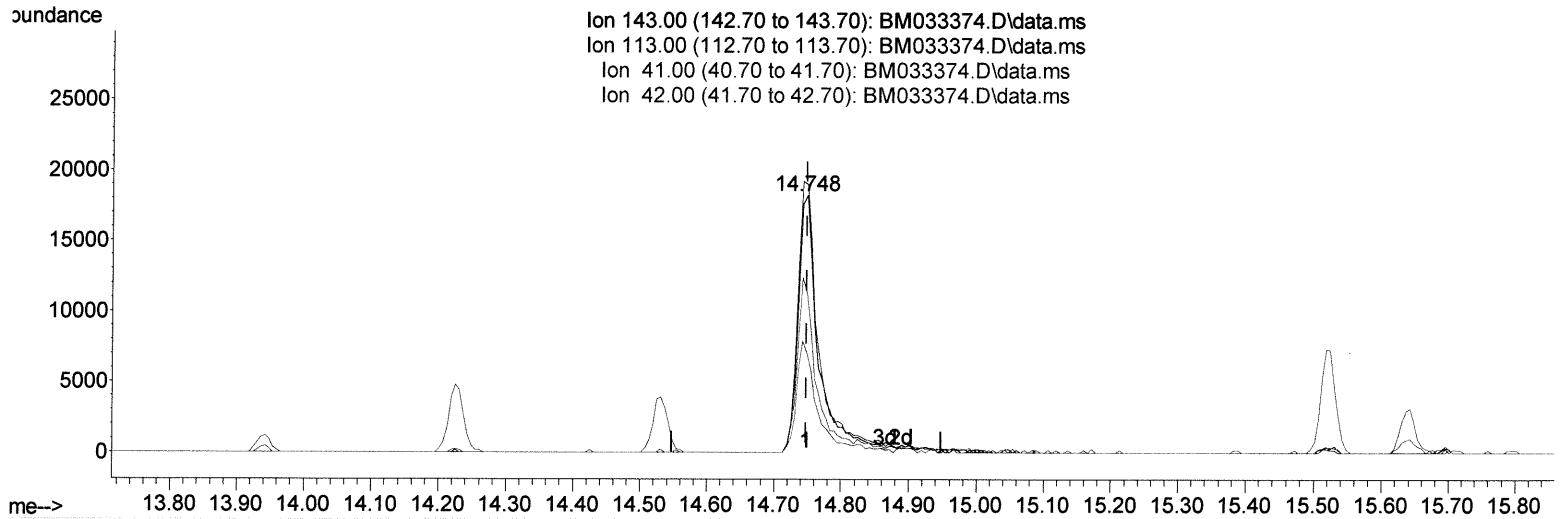
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(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.000) 25.58 ng/ul

response 38809

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	104.26
41.00	57.20	61.83
42.00	39.50	38.47

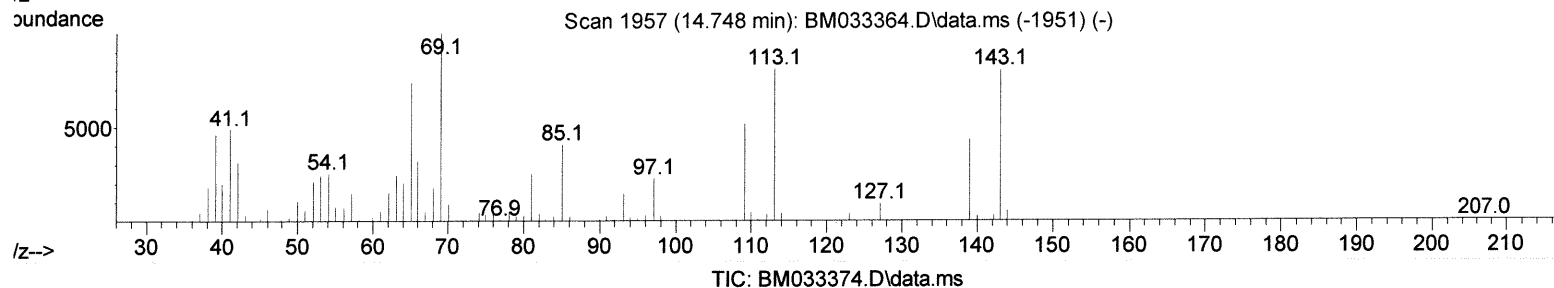
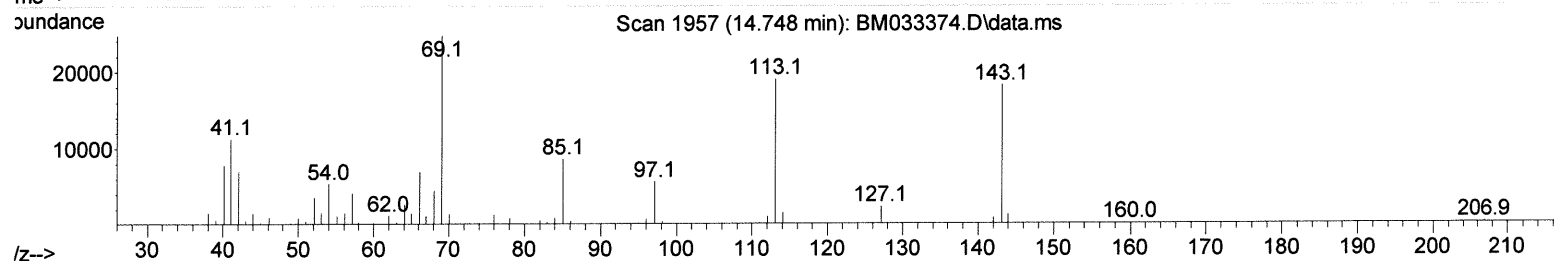
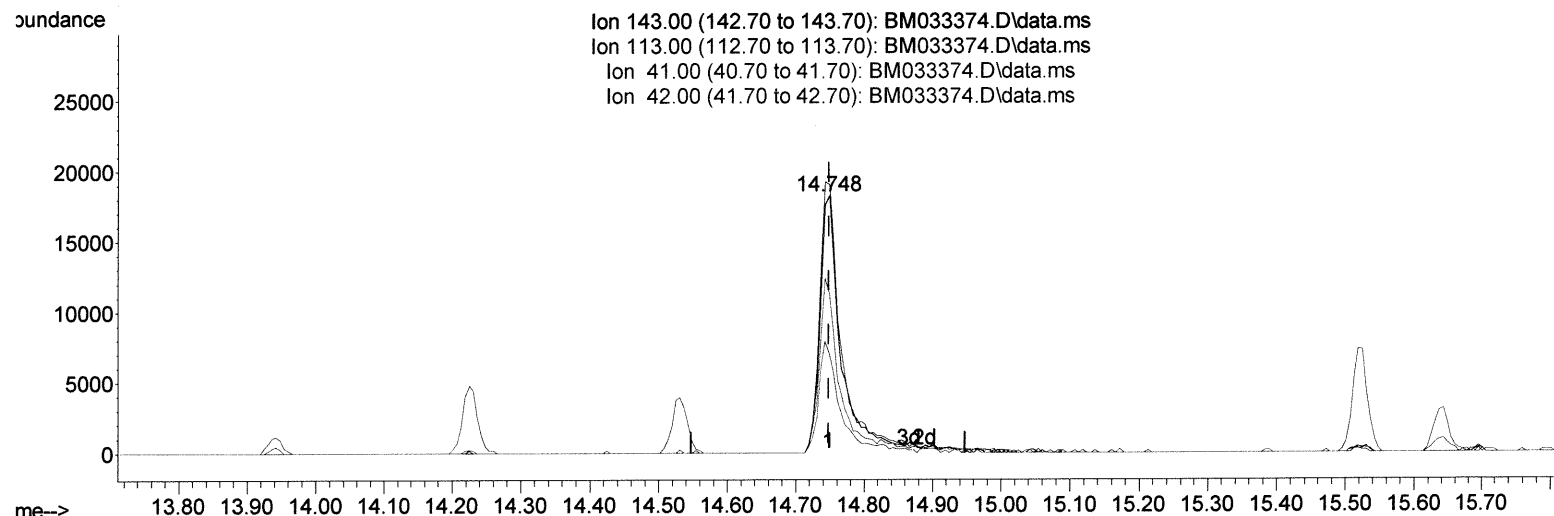
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 ALS Vial : 29 Sample Multiplier: 1

Instrument :
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 Supervised By :mohammad ahmed 12/15/2021



(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.000) 24.43 ng/ul m

response 37065

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	104.26
41.00	57.20	61.83
42.00	39.50	38.47

Ju 12/20/21

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.907	152	45643	20.000	ng/ul	0.00
20) Naphthalene-d8	10.701	136	178486	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.531	164	111844	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.272	188	235987	20.000	ng/ul	0.00
79) Chrysene-d12	21.430	240	224995	20.000	ng/ul	0.00
88) Perylene-d12	23.753	264	226768	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.367	96	8339m	6.866	ng/ul	0.00
4) Pyridine-d5	3.784	84	110208	31.295	ng/ul	0.00
7) Phenol-d5	7.078	99	132699	30.703	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.237	67	89821	31.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.443	132	98003	32.387	ng/ul	0.00
15) 4-Methylphenol-d8	8.613	113	100641	29.749	ng/ul	0.00
21) Nitrobenzene-d5	9.066	128	50186	34.653	ng/ul	0.00
24) 2-Nitrophenol-d4	9.790	143	51954	34.965	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.331	165	90311	32.162	ng/ul	0.00
31) 4-Chloroaniline-d4	10.843	131	131079	31.484	ng/ul	0.00
46) Dimethylphthalate-d6	13.942	166	298281	35.694	ng/ul	0.00
49) Acenaphthylene-d8	14.225	160	366634	35.391	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	37065m	24.428	ng/ul	0.00
60) Fluorene-d10	15.525	176	254755	34.081	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.642	200	37274	26.170	ng/ul	0.00
73) Anthracene-d10	17.372	188	411679	35.299	ng/ul	0.00
81) Pyrene-d10	19.654	212	484057	38.494	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.606	264	439910	35.790	ng/ul	0.00

Target Compounds Qvalue

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