

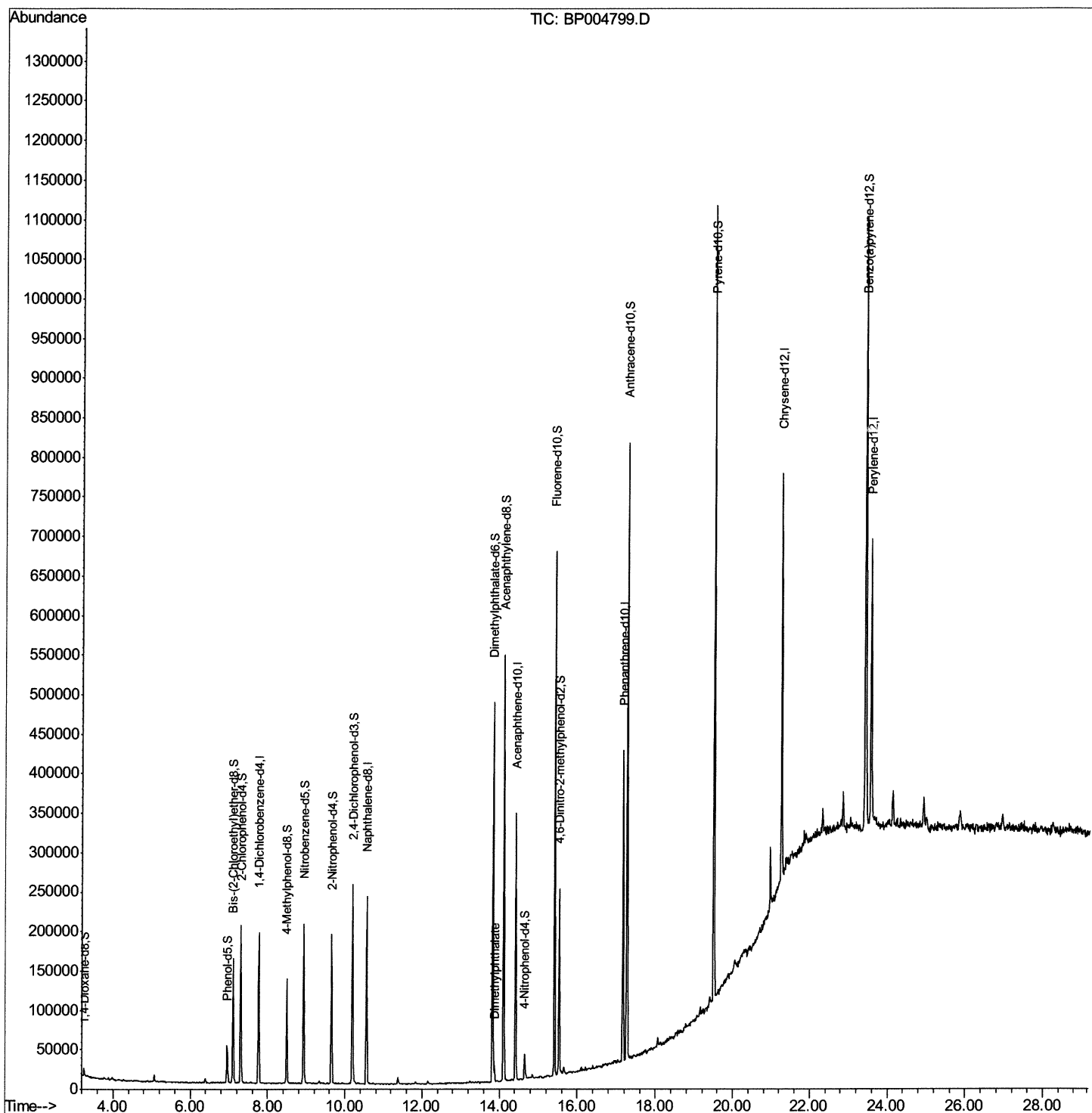
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP021721\
Data File : BP004799.D
Acq On : 18 Feb 2021 06:22
Operator : CG/JU
Sample : M1407-18
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBGM3

Manual Integrations
APPROVED

mohammad
2/18/2021 1:40:51 PM

Quant Time: Feb 18 10:01:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Feb 15 12:16:19 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

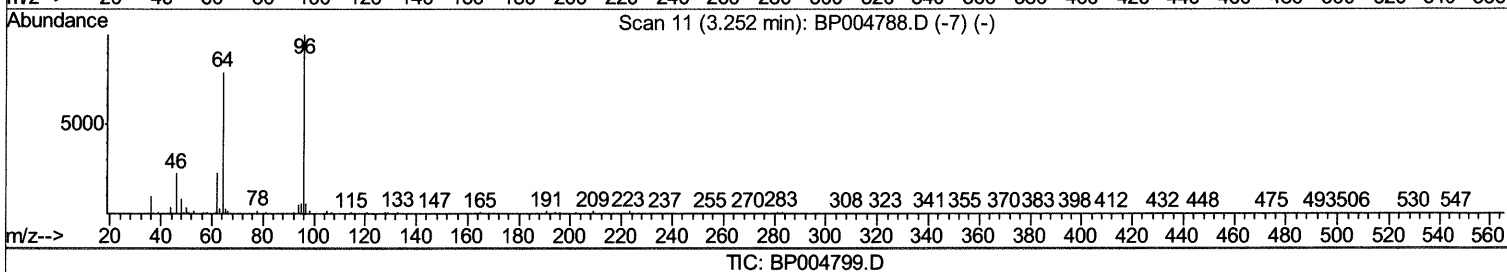
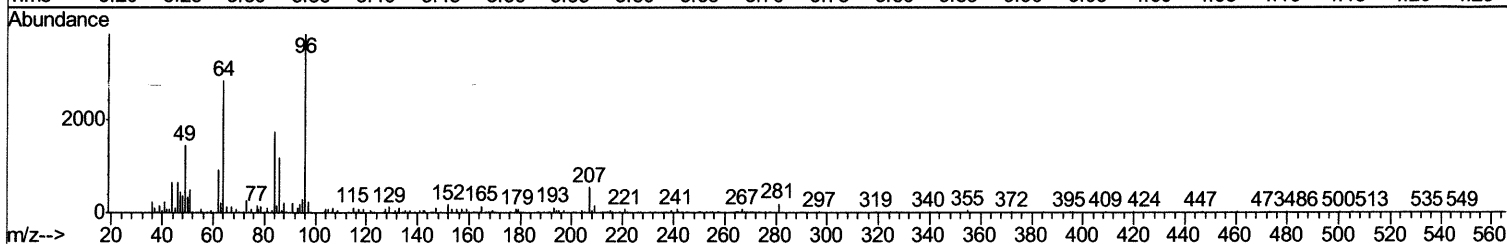
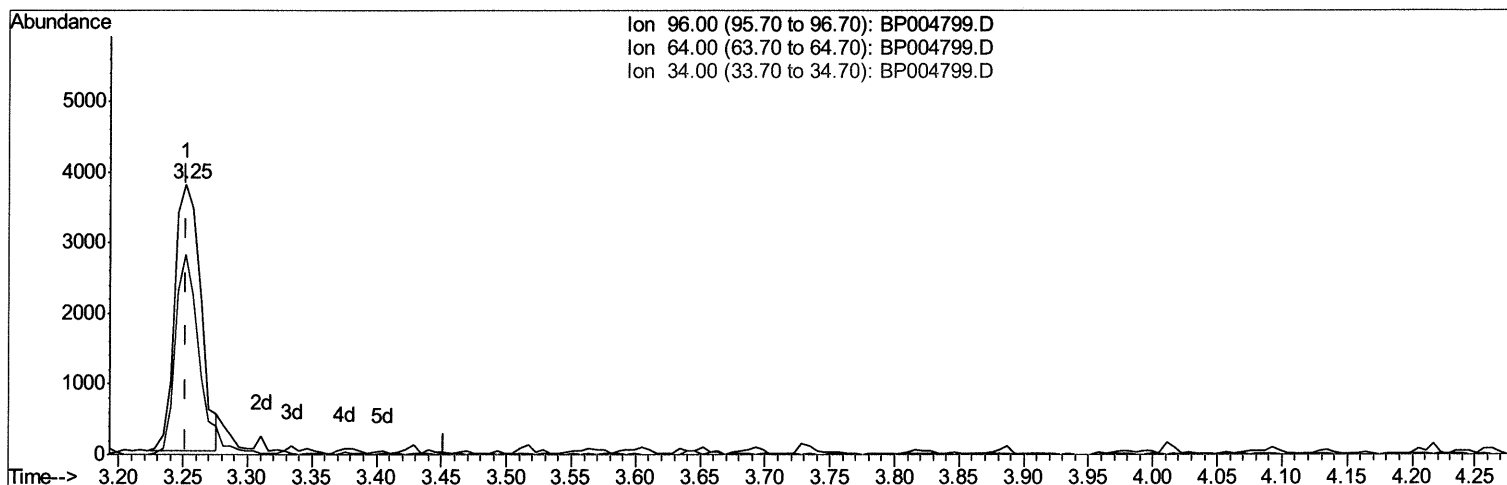
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(3) 1,4-Dioxane-d8 (S)

3.252min (-0.000) 4.21ng/uL

response 5308

Ion	Exp%	Act%
96.00	100	100
64.00	74.00	73.88
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

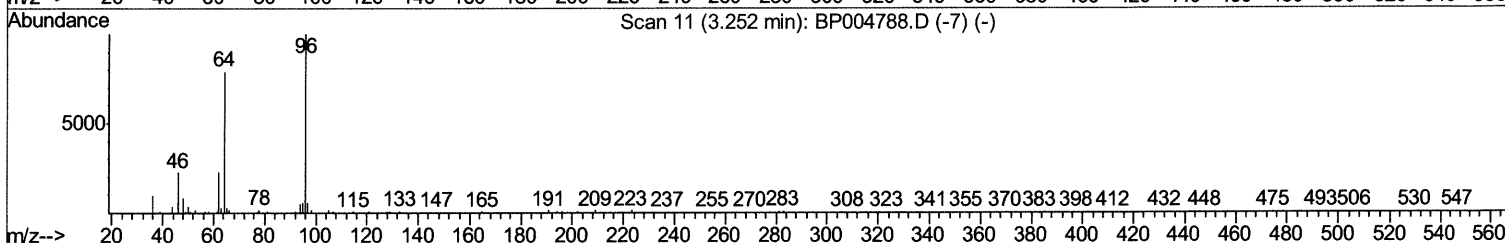
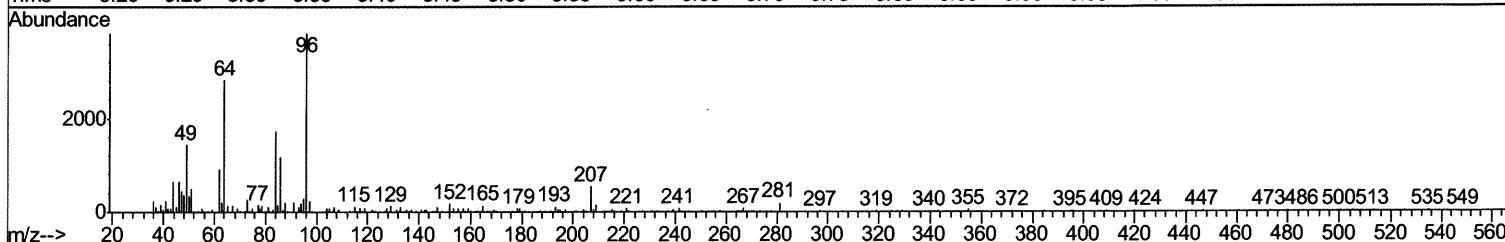
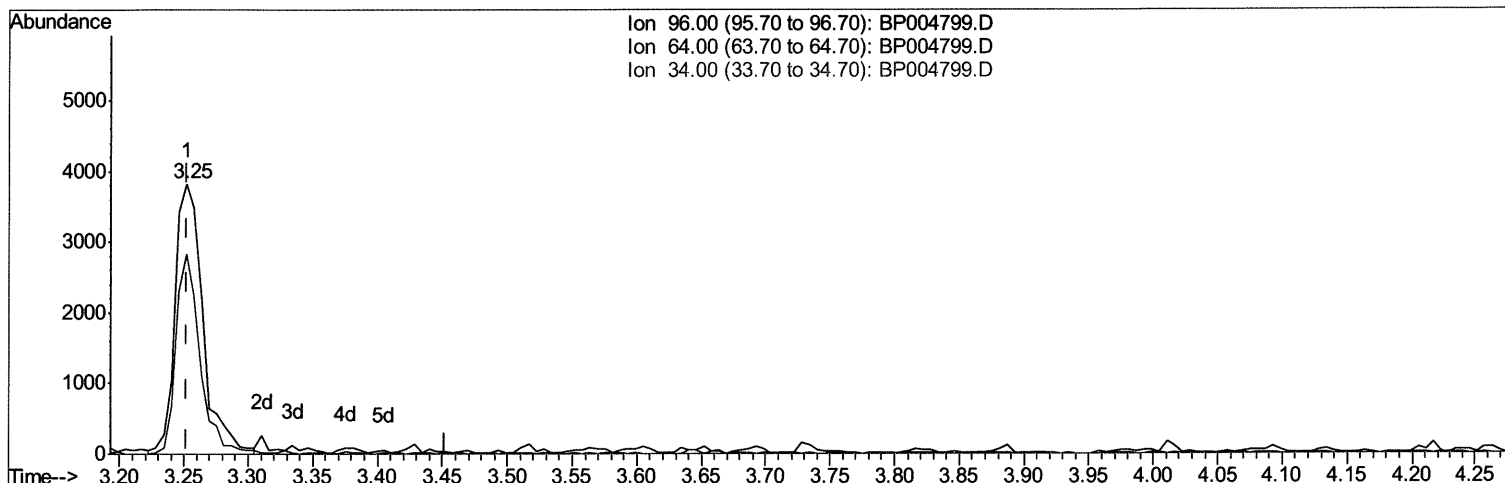
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Manual Integrations
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TIC: BP004799.D

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

3.252min (-0.000) 4.62ng/uL m 02/22/21 JU

response 5827

Ion	Exp%	Act%
96.00	100	100
64.00	74.00	73.88
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

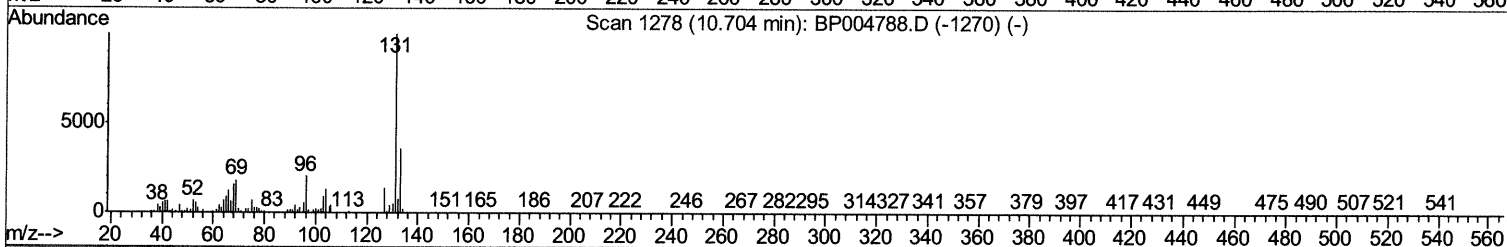
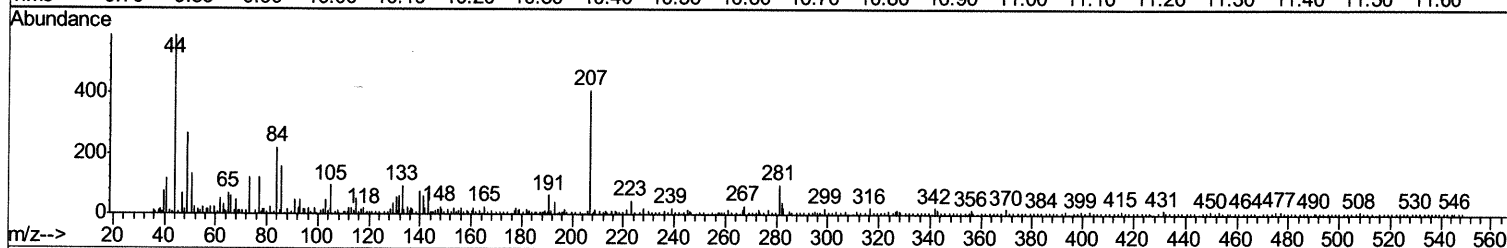
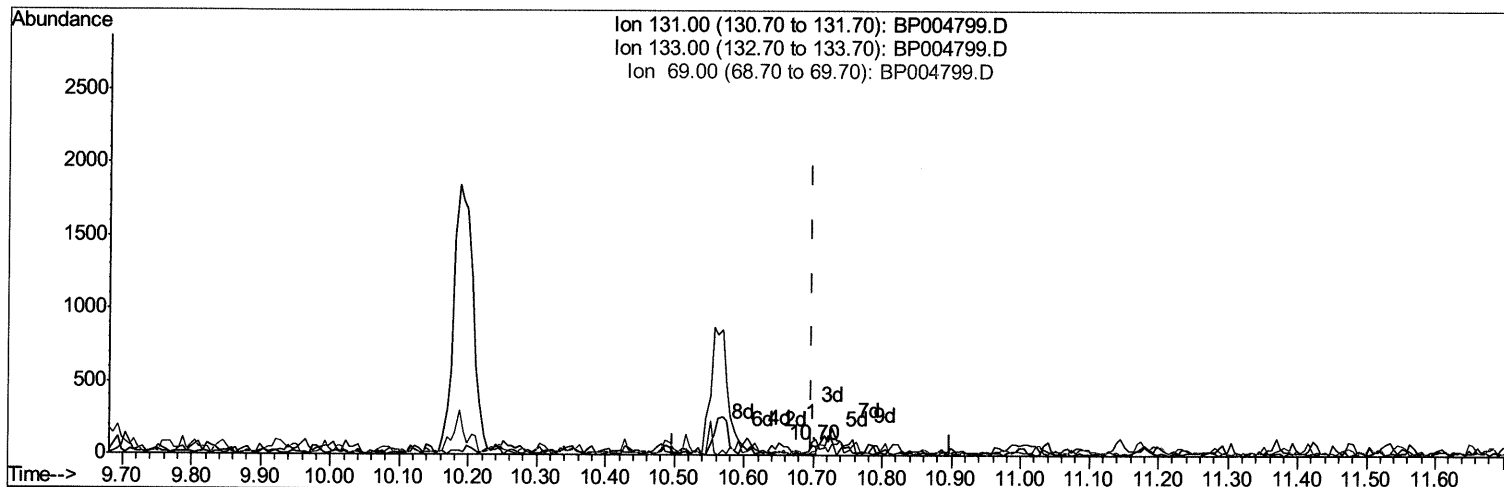
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 Sample : M1407-18
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBG3

Manual Integrations
 APPROVED

mohammad
 2/18/2021 1:40:51 PM

Quant Time: Feb 18 09:56:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
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TIC: BP004799.D

(29) 4-Chloroaniline-d4 (S)

10.699min (-0.000) 0.02ng/ul

response 52

Ion	Exp%	Act%
131.00	100	100
133.00	31.20	138.24#
69.00	20.50	39.71#
0.00	0.00	0.00

Quantitation Report (Qedit)

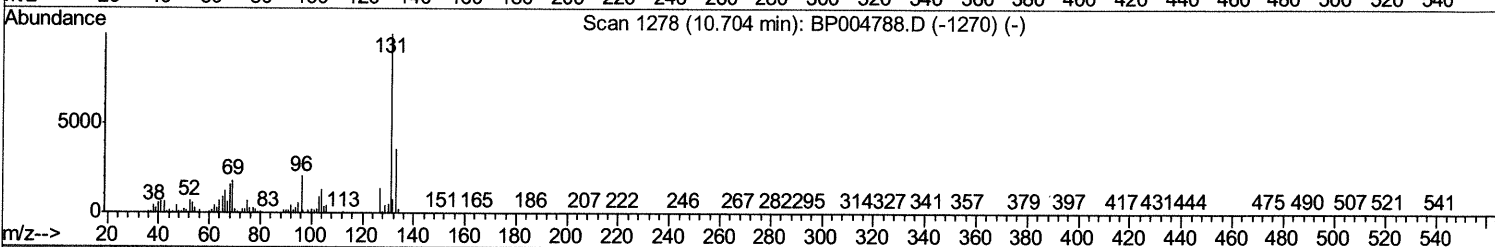
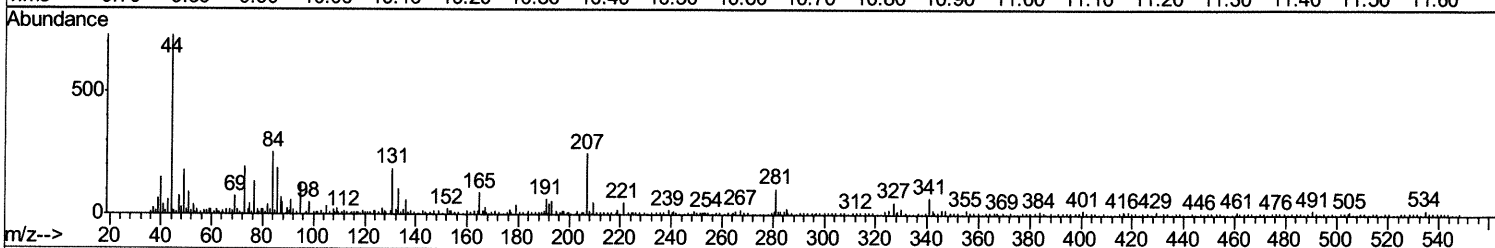
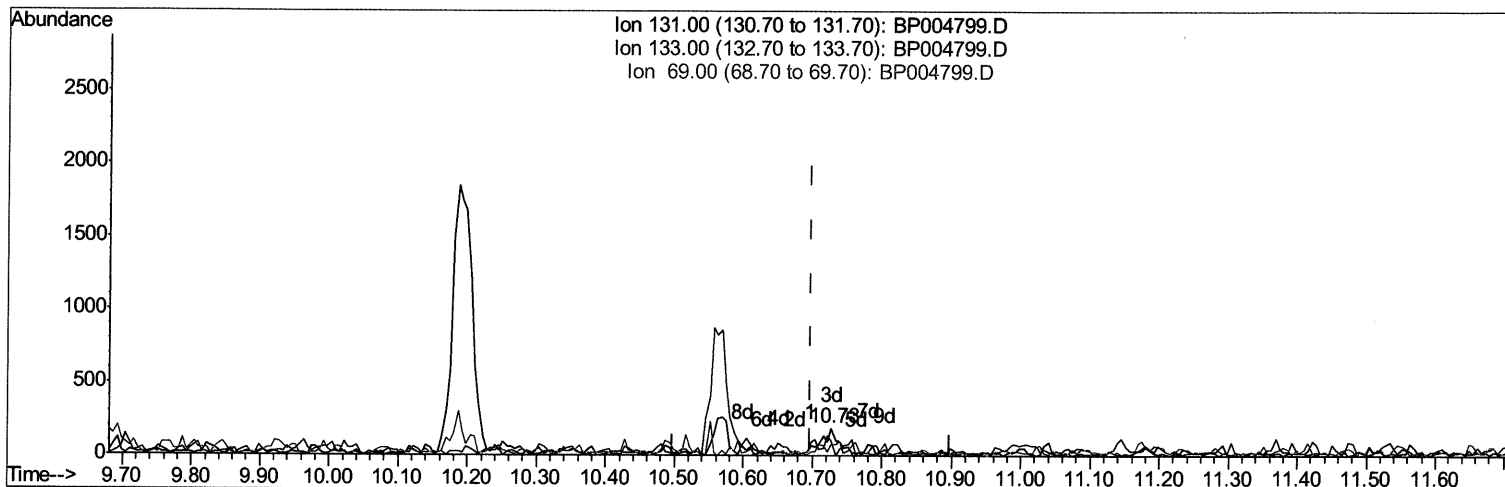
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 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
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Manual Integrations
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TIC: BP004799.D

(29) 4-Chloroaniline-d4 (S)

10.728min (+0.029) 0.11ng/ul m 02/22/21 JU

response 346

Ion	Exp%	Act%
131.00	100	100
133.00	31.20	55.38#
69.00	20.50	39.25#
0.00	0.00	0.00

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 ALS Vial : 35 Sample Multiplier: 1

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	48743	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	181825	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	106433	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	221661	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	245229	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	261224	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	5827m>	4.62	ng/uL >	0.00 02/22/21 JU
5) Phenol-d5	6.96	99	26167	6.69	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.10	67	62342	30.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	82881	27.96	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	46917	15.24	ng/ul	0.02
19) Nitrobenzene-d5	8.93	128	47674	37.98	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	54982	39.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	89764	32.37	ng/ul	0.01
29) 4-Chloroaniline-d4	10.73	131	346m>	0.11	ng/ul >	0.03 02/22/21 JU
44) Dimethylphthalate-d6	13.83	166	296376	38.36	ng/ul	0.00
47) Acenaphthylene-d8	14.11	160	349282	37.16	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	8542	6.33	ng/ul	0.04
58) Fluorene-d10	15.41	176	249650	37.71	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	47929	35.77	ng/ul	0.00
71) Anthracene-d10	17.27	188	422188	43.19	ng/ul	0.00
79) Pyrene-d10	19.51	212	498179	43.40	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	566034	42.83	ng/ul	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.87	163	12212	1.533	ng/ul#	93

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