

(QT Reviewed)

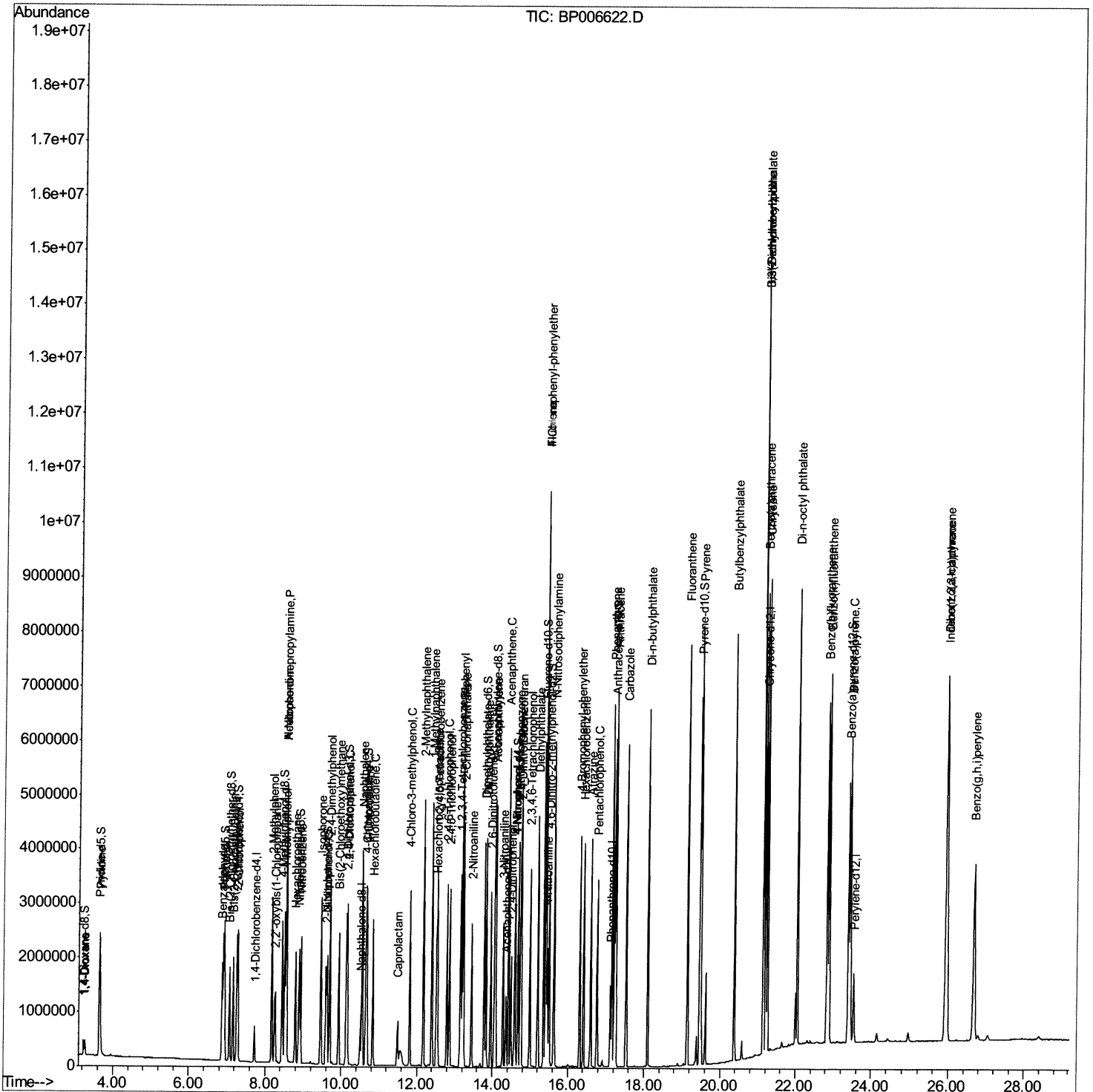
```
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP081021\  
Data File : BP006622.D  
Acq On    : 10 Aug 2021  14:00  
Operator  : CG/JU  
Sample    : SSTD08012  
Misc      :  
ALS Vial  : 7      Sample Multiplier: 1
```

Instrument :
BNA_P
ClientSampleId :
SSTD080612

Manual Integrations

mohammad
8/11/2021 9:27:46 PM

Quant Time: Aug 10 14:33:15 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 10 13:23:21 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

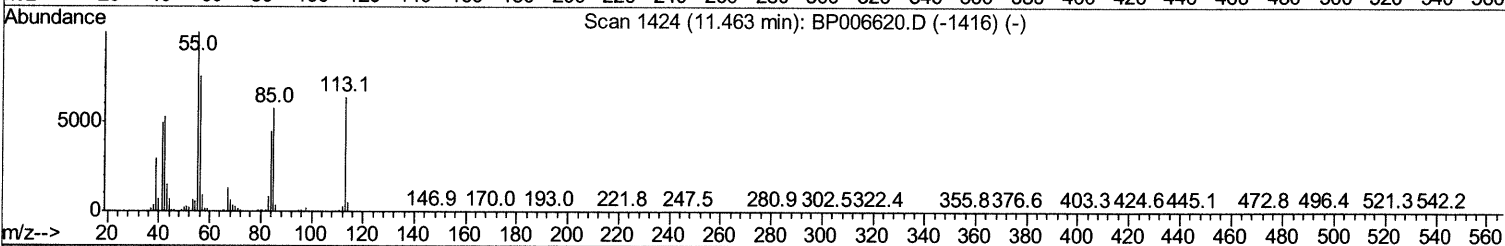
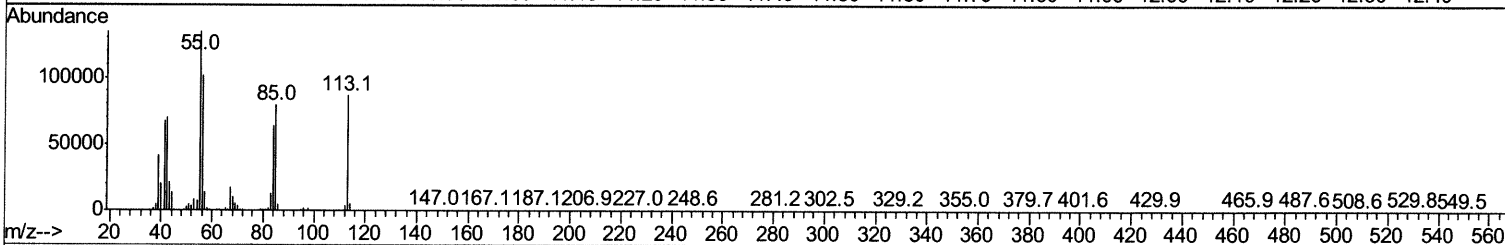
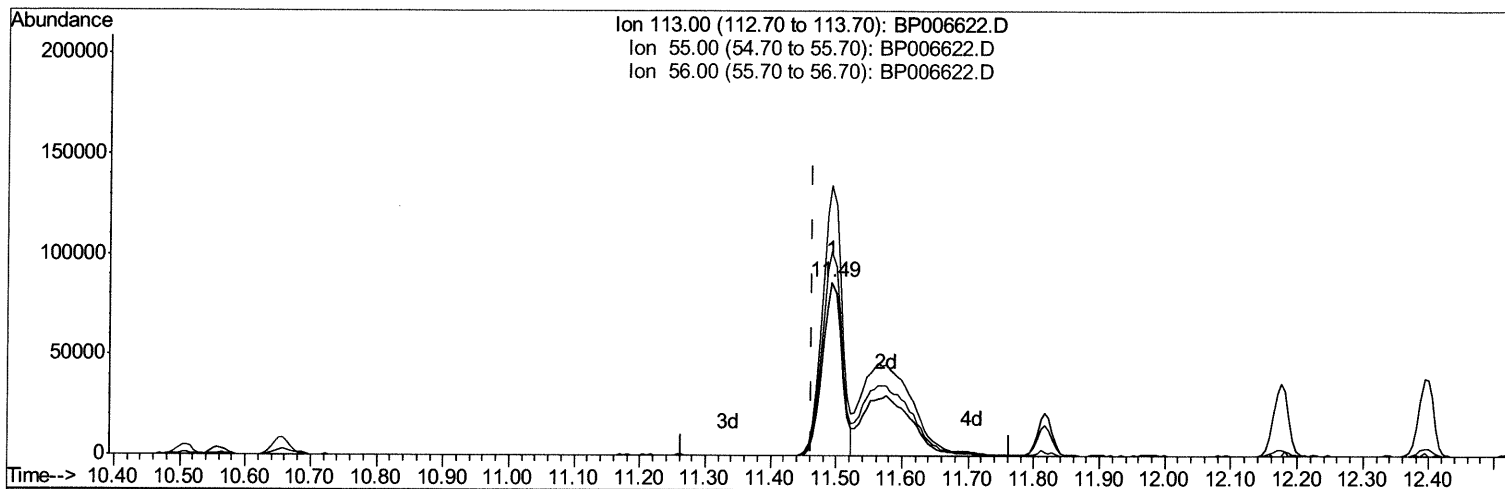
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP081021\
 Data File : BP006622.D
 Acq On : 10 Aug 2021 14:00
 Operator : CG/JU
 Sample : SST08012
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
 SST080612

Manual Integrations
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Quant Time: Aug 10 14:31:25 2021
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TIC: BP006622.D

(34) Caprolactam

11.493min (+0.029) 51.27ng/ul

response 176049

Ion	Exp%	Act%
113.00	100	100
55.00	158.10	155.88
56.00	118.70	117.76
0.00	0.00	0.00

Quantitation Report (Qedit)

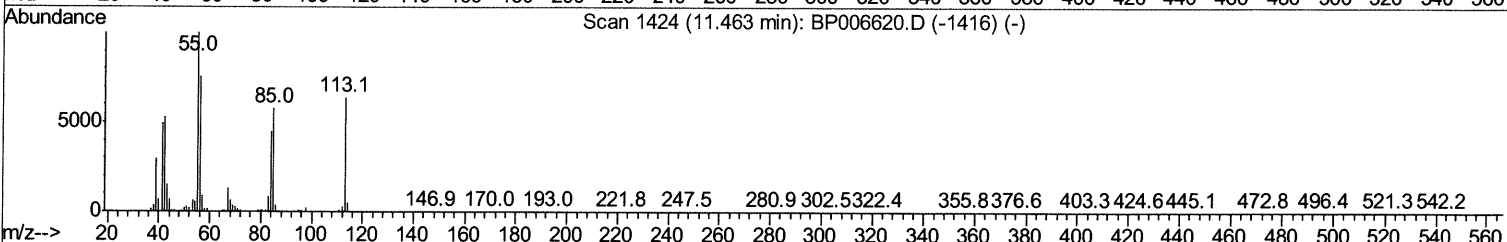
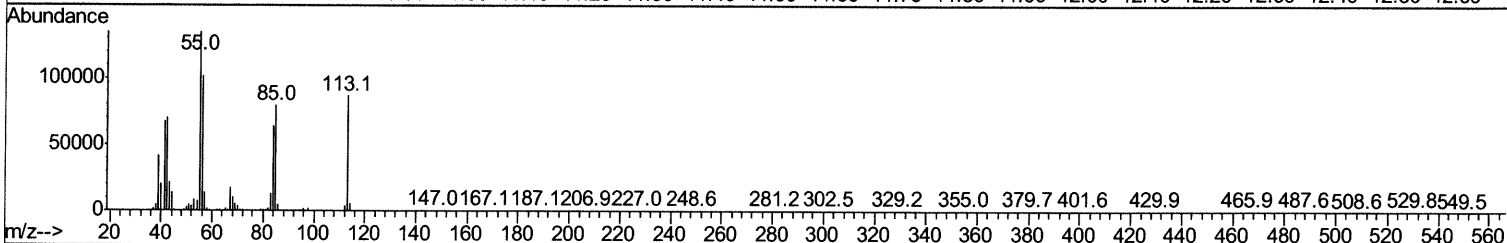
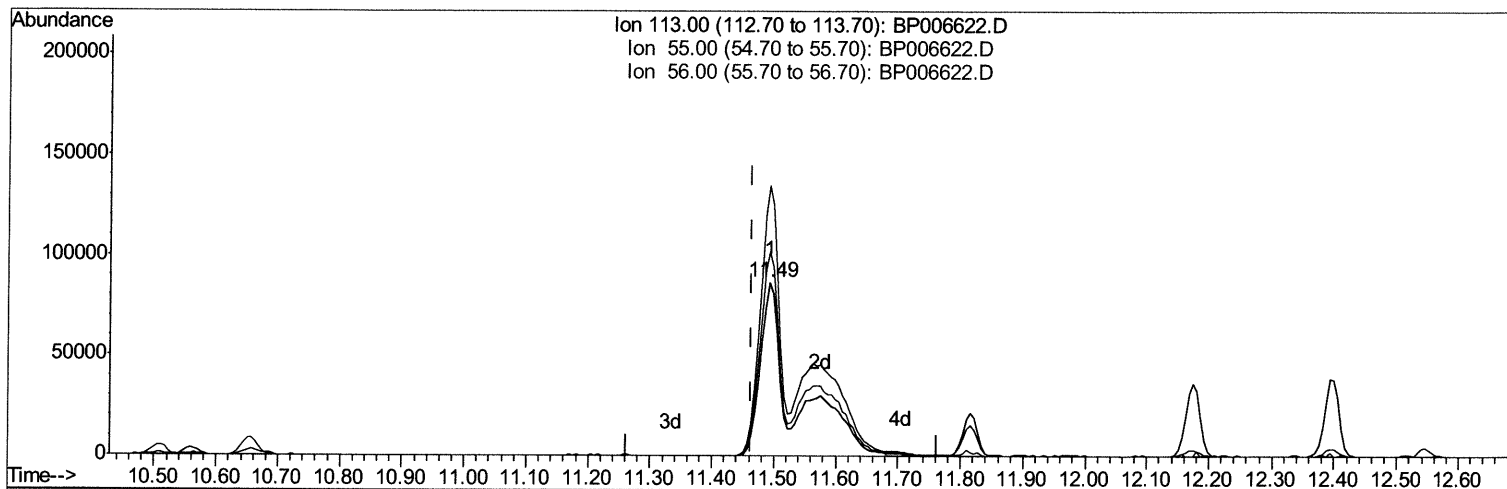
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TIC: BP006622.D

(34) Caprolactam

11.493min (+0.029) 95.95ng/ul m

response 329453

Ion	Exp%	Act%
113.00	100	100
55.00	158.10	155.88
56.00	118.70	117.76
0.00	0.00	0.00

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Quantitation Report (Qedit)

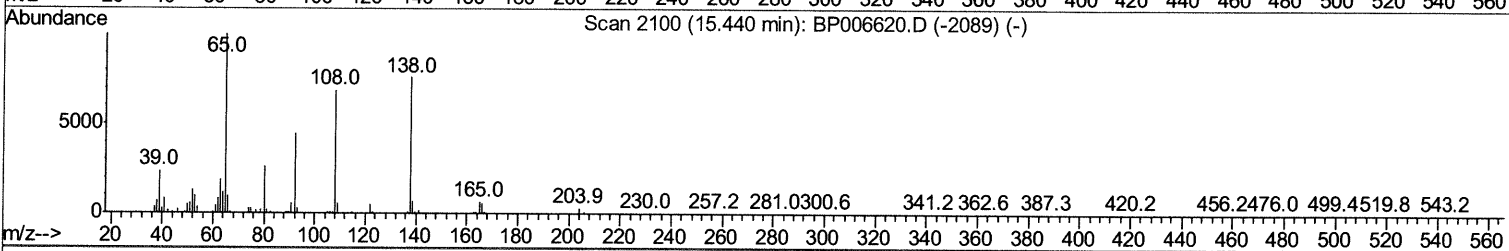
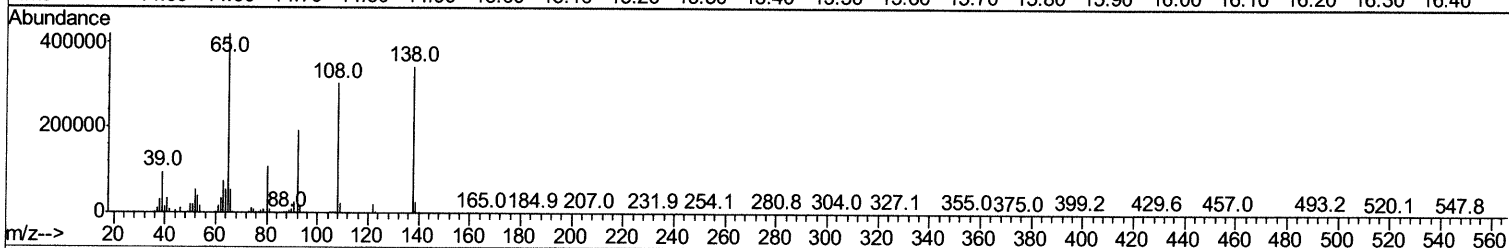
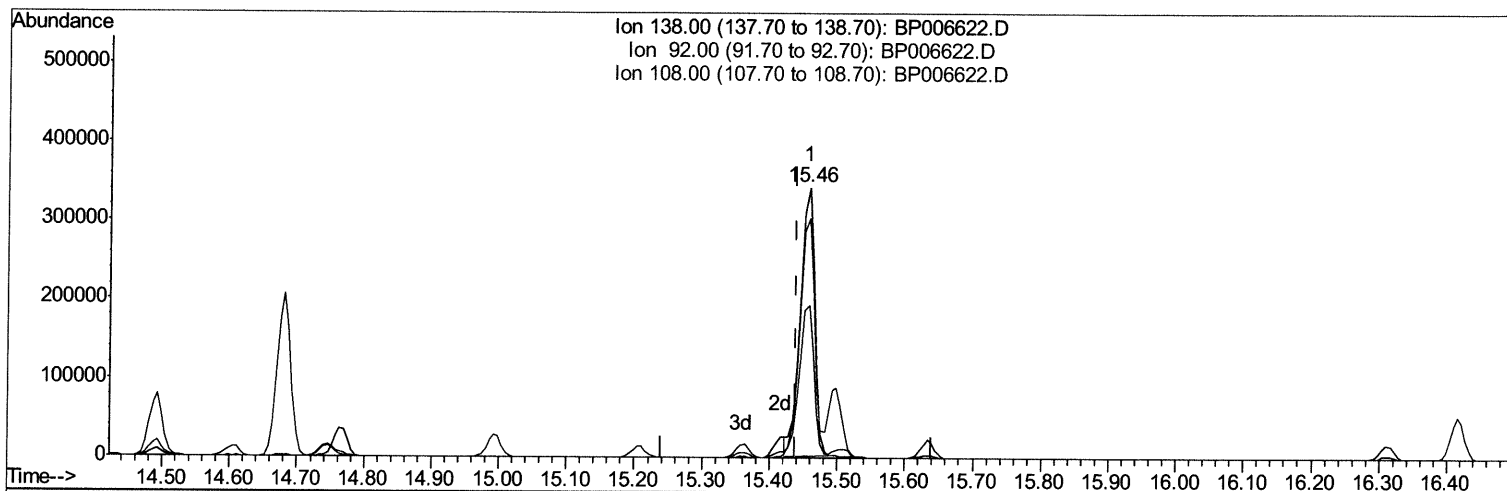
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TIC: BP006622.D

(63) 4-Nitroaniline

15.457min (+0.018) 82.30ng/ul

response 502715

Ion	Exp%	Act%
138.00	100	100
92.00	60.20	56.23
108.00	91.20	88.84
0.00	0.00	0.00

Quantitation Report (Qedit)

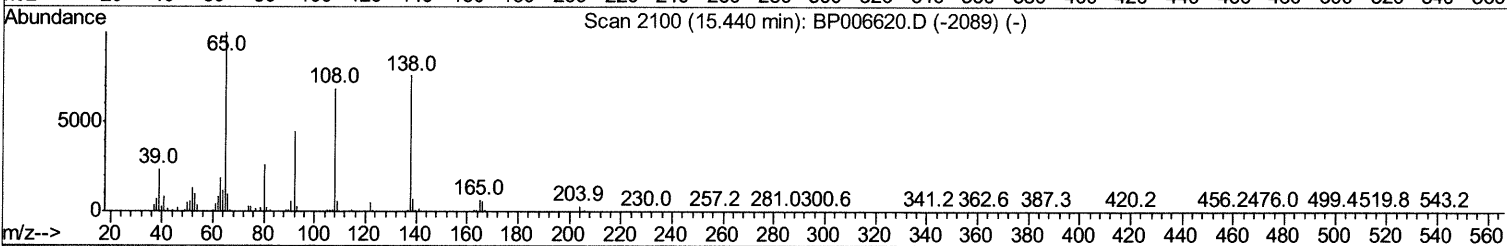
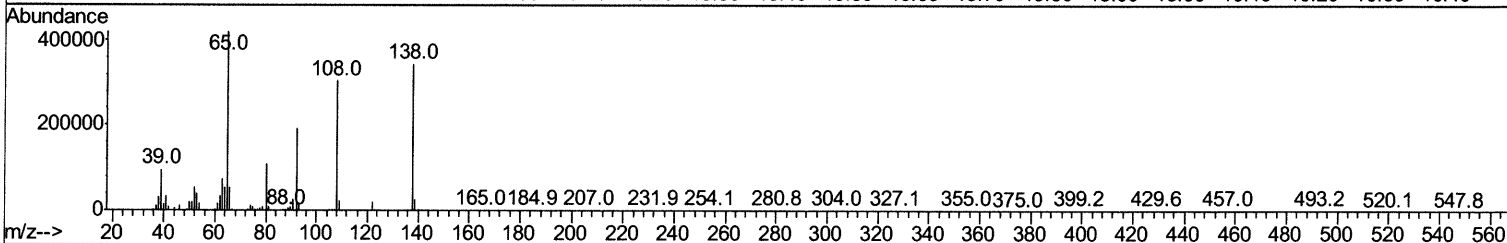
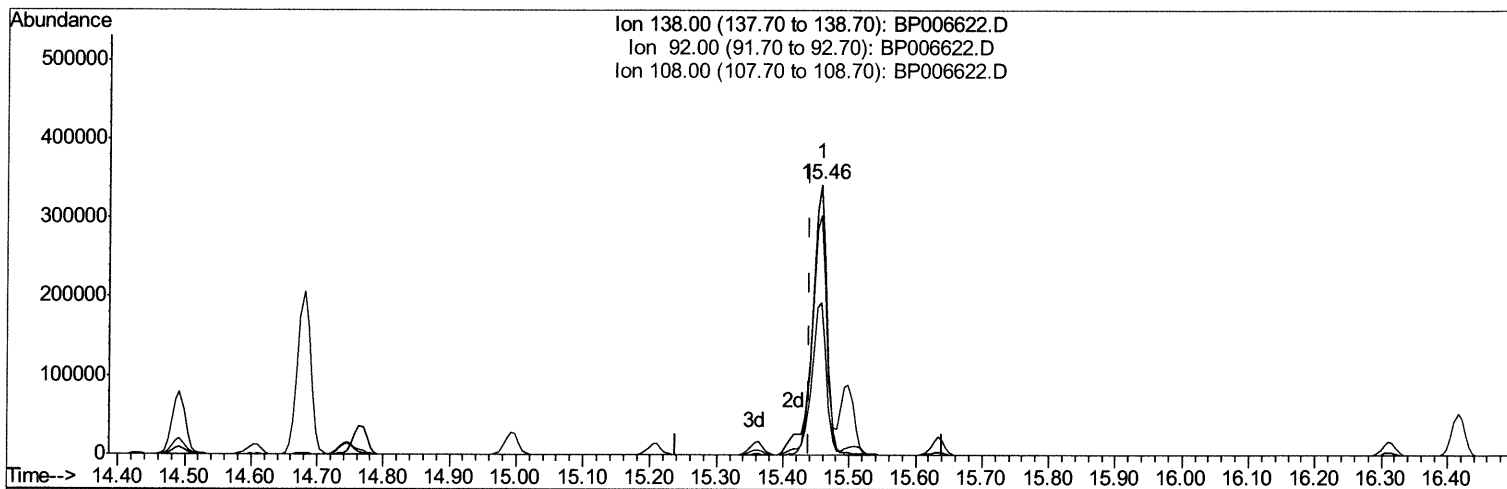
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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TIC: BP006622.D

(63) 4-Nitroaniline

15.457min (+0.018) 88.79ng/ul m

response 542339

Ion	Exp%	Act%
138.00	100	100
92.00	60.20	56.23
108.00	91.20	88.84
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	160508	20.00	ng/ul	0.00
20) Naphthalene-d8	10.51	136	727524	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.36	164	415602	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.12	188	818401	20.00	ng/ul	0.00
79) Chrysene-d12	21.22	240	775866	20.00	ng/ul	0.00
88) Perylene-d12	23.54	264	813968	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	145819	33.91	ng/uL	0.00
4) Pividine-d5	3.63	84	1077127	84.38	ng/ul	0.00
7) Phenol-d5	6.91	99	1285982	85.31	ng/ul	0.01
9) Bis-(2-Chloroethyl)ether-d	7.07	67	776605	82.80	ng/ul	0.00
11) 2-Chlorophenol-d4	7.26	132	998435	89.15	ng/ul	0.00
15) 4-Methylphenol-d8	8.45	113	1018709	85.93	ng/ul	0.00
21) Nitrobenzene-d5	8.89	128	507638	93.71	ng/ul	0.00
24) 2-Nitrophenol-d4	9.60	143	540934	104.58	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.14	165	900255	88.82	ng/ul	0.00
31) 4-Chloroaniline-d4	10.66	131	1365719	81.84	ng/ul	0.00
46) Dimethylphthalate-d6	13.79	166	2533542	85.65	ng/ul	0.00
49) Acenaphthylene-d8	14.06	160	3187520	87.71	ng/ul	0.00
54) 4-Nitrophenol-d4	14.59	143	534202	90.96	ng/ul	0.02
60) Fluorene-d10	15.36	176	2060769	82.87	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.50	200	450553	98.75	ng/ul	0.01
73) Anthracene-d10	17.22	188	3093663	83.74	ng/ul	0.00
81) Pyrene-d10	19.46	212	3299421	82.68	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.40	264	3521702	85.14	ng/ul	0.02

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.26	88	151957	33.471	ng/uL 98
5) Pividine	3.65	79	1115277	83.775	ng/ul 99
6) Benzaldehyde	6.88	77	624893	75.977	ng/ul 98
8) Phenol	6.93	94	1301258	84.447	ng/ul 98
10) Bis(2-Chloroethyl)ether	7.16	93	990216	81.696	ng/ul 98
12) 2-Chlorophenol	7.29	128	994061	87.548	ng/ul 99
13) 2-Methylphenol	8.18	108	959743	85.352	ng/ul 98
14) 2,2'-oxybis(1-Chloropropan	8.26	45	1145631	84.604	ng/ul 99
16) Acetophenone	8.55	105	1577978	86.561	ng/ul 100
17) N-Nitroso-di-n-propylamine	8.55	70	876635	91.920	ng/ul 99
18) 4-Methylphenol	8.51	108	1032095	85.477	ng/ul 98
19) Hexachloroethane	8.79	117	419360	90.291	ng/ul 98
22) Nitrobenzene	8.93	77	1286552	91.537	ng/ul 99
23) Isophorone	9.46	82	2346611	93.393	ng/ul 98
25) 2-Nitrophenol	9.63	139	562828	98.675	ng/ul 97
26) 2,4-Dimethylphenol	9.70	107	1183547	87.495	ng/ul 99
27) Bis(2-Chloroethoxy)methane	9.94	93	1354934	83.973	ng/ul 99
29) 2,4-Dichlorophenol	10.16	162	858997	86.382	ng/ul 99
30) Naphthalene	10.56	128	3093830	81.964	ng/ul 100
32) 4-Chloroaniline	10.68	127	1345909	81.876	ng/ul 100
33) Hexachlorobutadiene	10.84	225	496165	81.786	ng/ul 99
34) Caprolactam	11.49	113	329453m	95.946	ng/ul 99

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35) 4-Chloro-3-methylphenol	11.82	107	1086308	91.508	ng/ul	97
36) 2-Methylnaphthalene	12.18	142	2125813	82.191	ng/ul	99
37) 1-Methylnaphthalene	12.40	142	2121273	81.481	ng/ul	99
39) 1,2,4,5-Tetrachlorobenzene	12.55	216	914803	81.679	ng/ul	97
40) Hexachlorocyclopentadiene	12.52	237	604118	82.915	ng/ul	99
41) 2,4,6-Trichlorophenol	12.79	196	640423	92.278	ng/ul	99
42) 2,4,5-Trichlorophenol	12.86	196	690196	90.894	ng/ul	99
43) 1,1'-Biphenyl	13.20	154	2566303	82.374	ng/ul	98
44) 2-Chloronaphthalene	13.23	162	1941099	82.522	ng/ul	99
45) 2-Nitroaniline	13.45	65	760975	109.517	ng/ul	99
47) Dimethylphthalate	13.84	163	2448466	84.330	ng/ul	100
48) 2,6-Dinitrotoluene	13.96	165	529004	99.663	ng/ul	99
50) Acenaphthylene	14.09	152	3019388	85.318	ng/ul	100
51) 3-Nitroaniline	14.28	138	586883	94.365	ng/ul	98
52) Acenaphthene	14.43	153	2147393	82.702	ng/ul	99
53) 2,4-Dinitrophenol	14.49	184	339002	107.685	ng/ul	99
55) 4-Nitrophenol	14.61	109	477921	100.603	ng/ul	97
56) Dibenzofuran	14.77	168	2833844	81.381	ng/ul	100
57) 2,4-Dinitrotoluene	14.75	165	761620	100.407	ng/ul	100
58) 2,3,4,6-Tetrachlorophenol	14.99	232	558260	94.275	ng/ul	97
59) Diethylphthalate	15.20	149	2658298	88.396	ng/ul	99
61) Fluorene	15.42	166	2378656	83.043	ng/ul	100
62) 4-Chlorophenyl-phenylether	15.42	204	1087825	81.462	ng/ul	98
63) 4-Nitroaniline	15.46	138	542339m	88.786	ng/ul	98
66) 4,6-Dinitro-2-methylphenol	15.51	198	433341	97.161	ng/ul#	98
67) N-Nitrosodiphenylamine	15.63	169	1983609	81.588	ng/ul	99
68) 4-Bromophenyl-phenylether	16.32	248	646365	98.619	ng/ul	96
69) Hexachlorobenzene	16.42	284	725498	82.539	ng/ul	100
70) Atrazine	16.60	200	680830	80.920	ng/ul	99
71) Pentachlorophenol	16.77	266	482192	87.262	ng/ul	99
72) Phenanthrene	17.16	178	3608253	82.107	ng/ul	100
74) Anthracene	17.26	178	3662103	82.185	ng/ul	98
75) 1,2,3,4-Tetrachlorobenzene	13.16	216	908595	80.993	ng/uL	99
76) Pentachlorobenzene	14.68	250	871857	79.721	ng/uL	99
77) Carbazole	17.53	167	3371857	82.722	ng/ul	100
78) Di-n-butylphthalate	18.10	149	4500300	92.378	ng/ul	100
80) Fluoranthene	19.14	202	4095816	83.294	ng/ul	100
82) Pyrene	19.49	202	4158023	81.489	ng/ul	97
83) Butylbenzylphthalate	20.38	149	2137616	102.101	ng/ul	96
84) 3,3'-Dichlorobenzidine	21.15	252	1383532	80.932	ng/ul	99
85) Benzo(a)anthracene	21.21	228	3952587	82.214	ng/ul	98
86) Bis(2-ethylhexyl)phthalate	21.15	149	3234313	98.668	ng/ul	98
87) Chrysene	21.26	228	3830072	83.113	ng/ul	99
89) Di-n-octyl phthalate	22.05	149	5485497	96.706	ng/ul	100
90) Benzo(b)fluoranthene	22.84	252	4263506	82.465	ng/ul	99
91) Benzo(k)fluoranthene	22.89	252	4053918	82.617	ng/ul	98
93) Benzo(a)pyrene	23.45	252	3780923	83.664	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	25.95	276	5002314	86.473	ng/ul	97
95) Dibenzo(a,h)anthracene	25.97	278	4189934	85.517	ng/ul	99
96) Benzo(g,h,i)perylene	26.69	276	4128967	86.388	ng/ul	98

2021/8/11

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed