

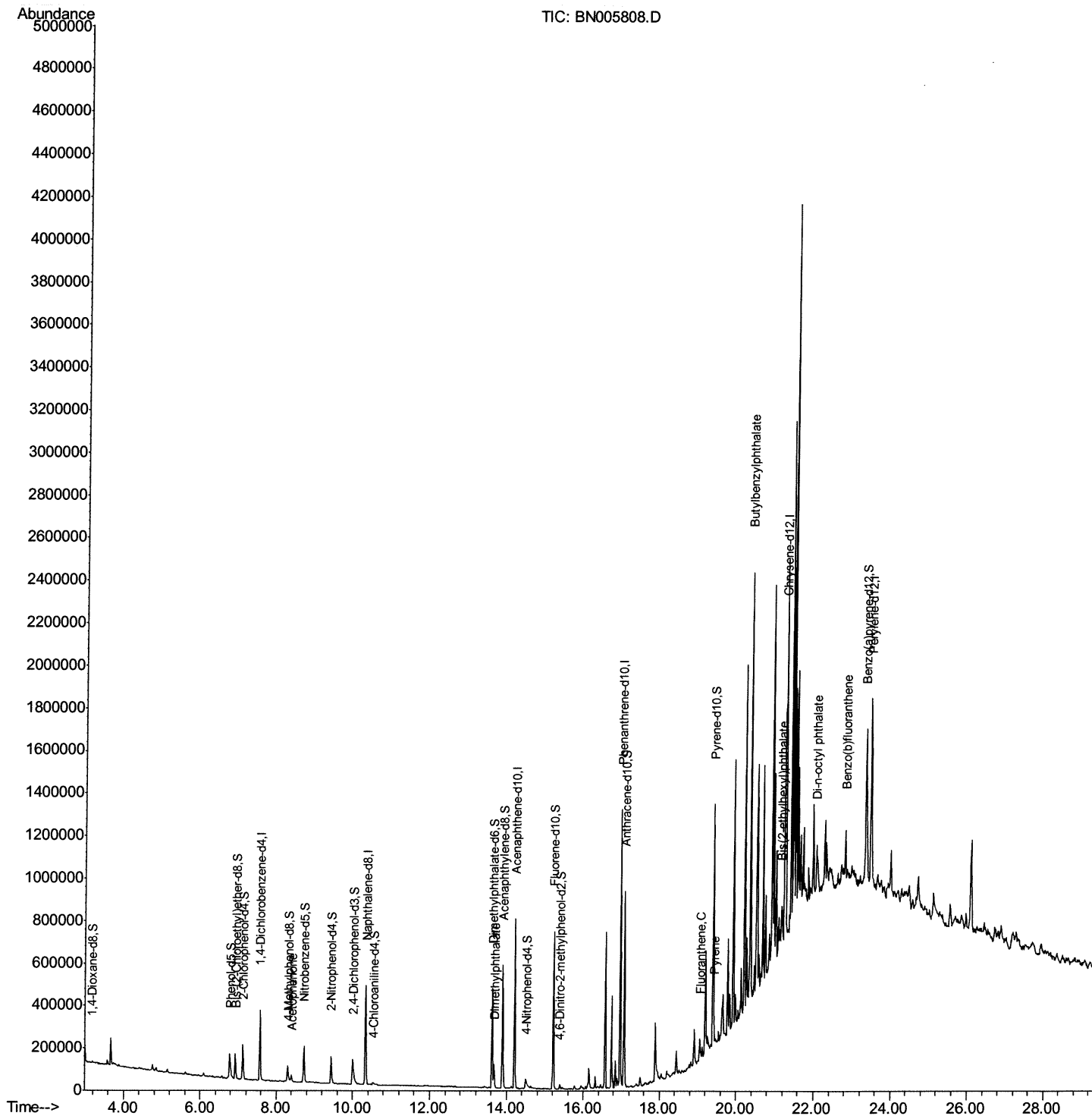
Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052019\
Data File : BN005808.D
Acq On : 20 May 2019 20:18
Operator : JU/SJ
Sample : K2693-10
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BFHL1

Manual Integrations
APPROVED

mohammad
5/21/2019 3:27:36 PM

Quant Time: May 21 04:43:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 21 04:05:47 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

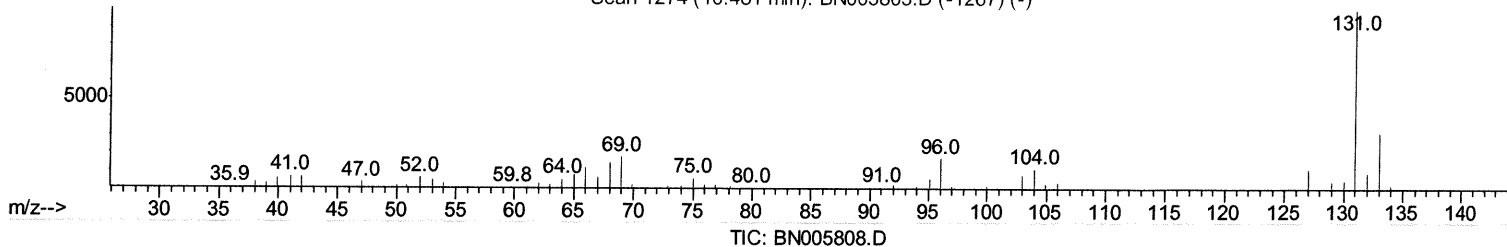
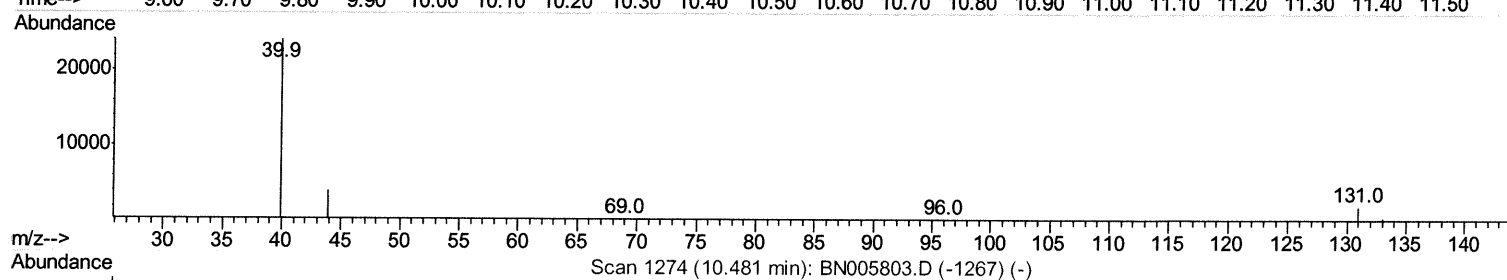
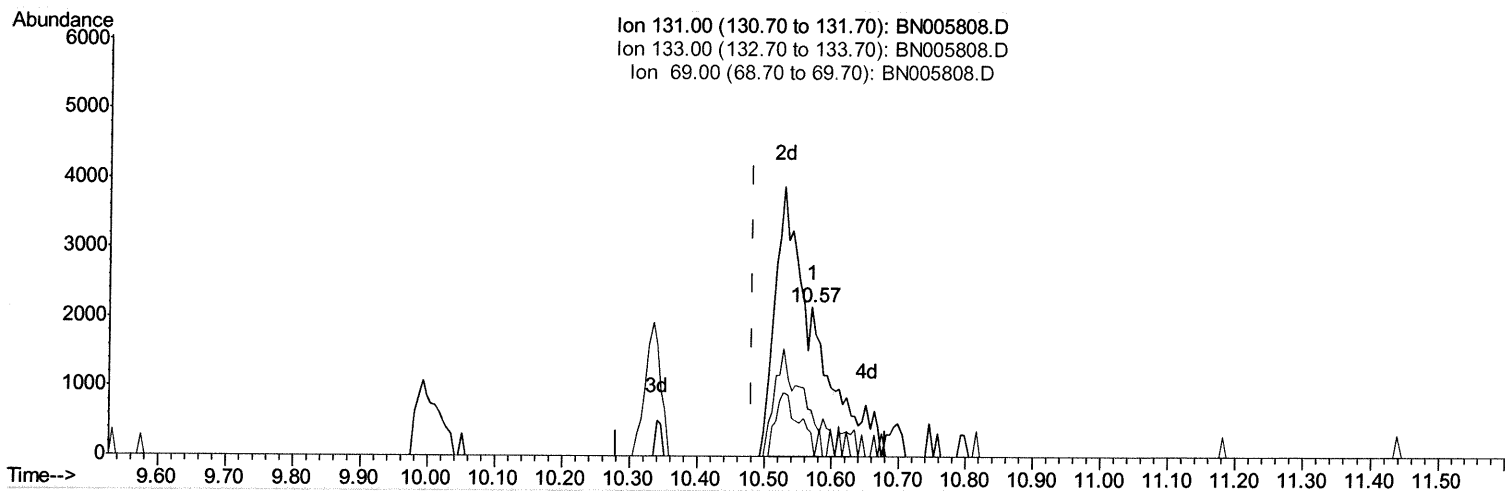
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(29) 4-Chloroaniline-d4 (S)

10.569min (+0.088) 0.22ng/ul

response 1565

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	31.14
69.00	18.90	16.47
0.00	0.00	0.00

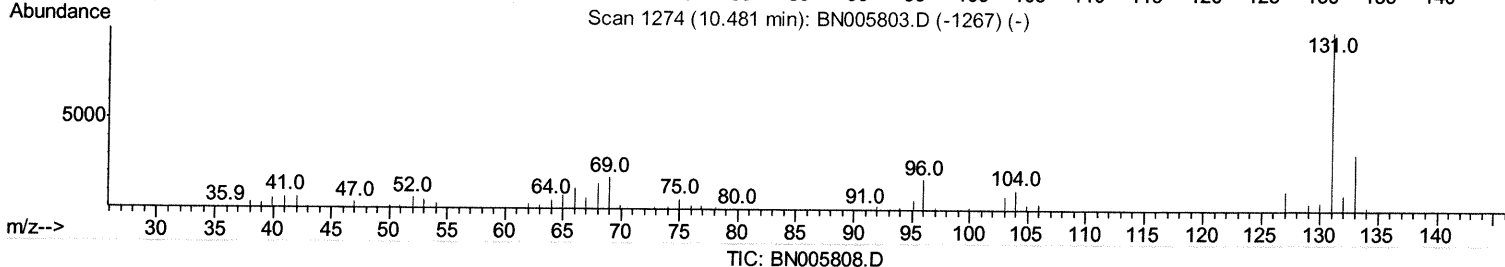
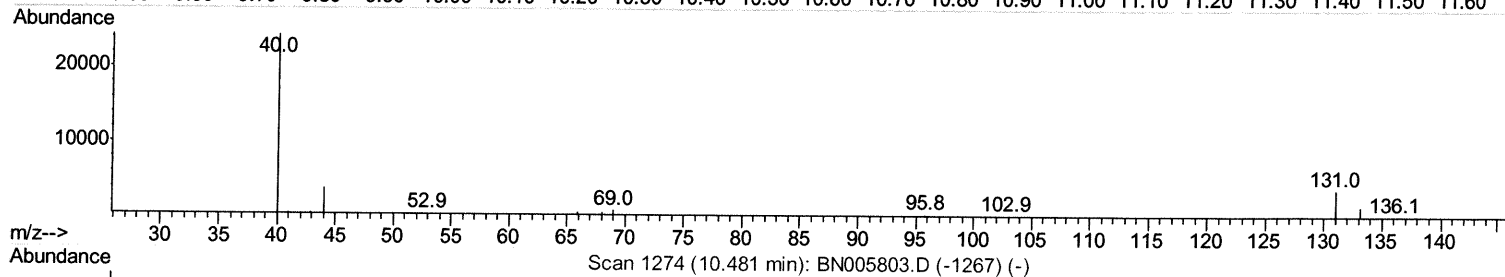
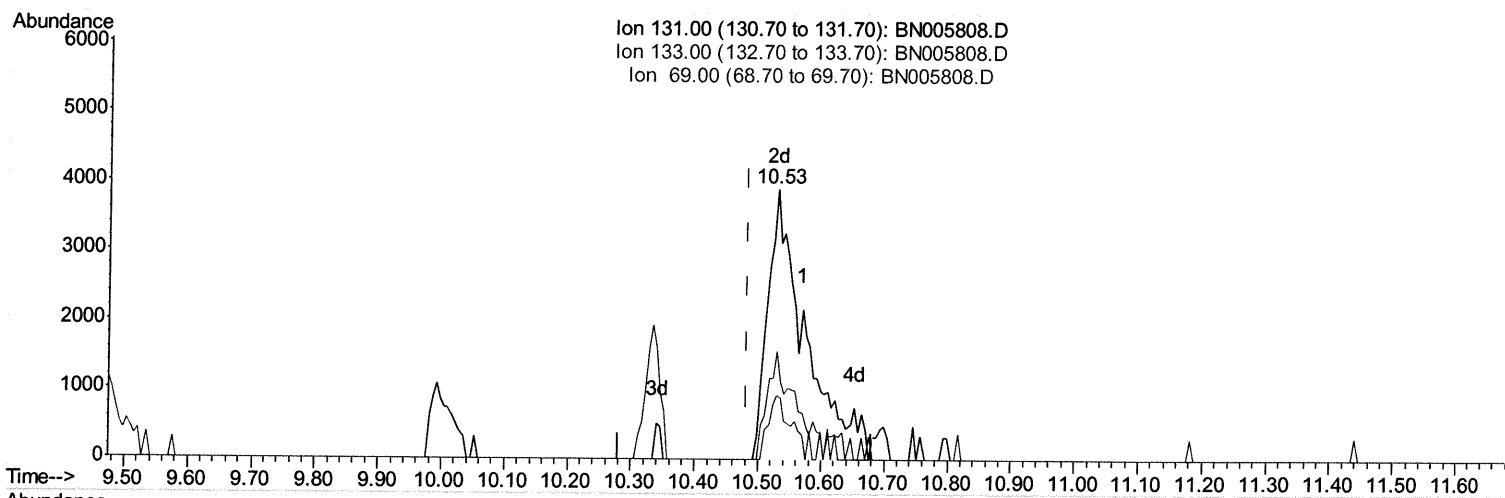
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(29) 4-Chloroaniline-d4 (S)

10.528min (+0.047) 2.25ng/ul m *JU 05/22/19*

response 16178

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	39.94
69.00	18.90	23.70#
0.00	0.00	0.00

Quantitation Report (Qedit)

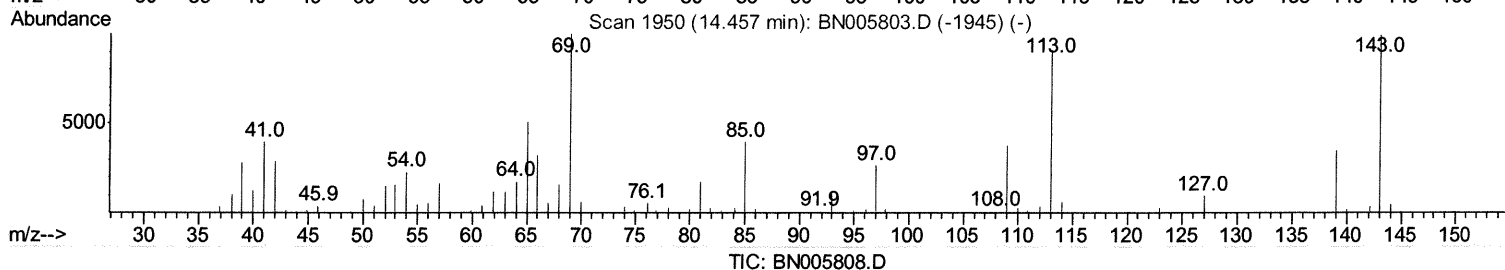
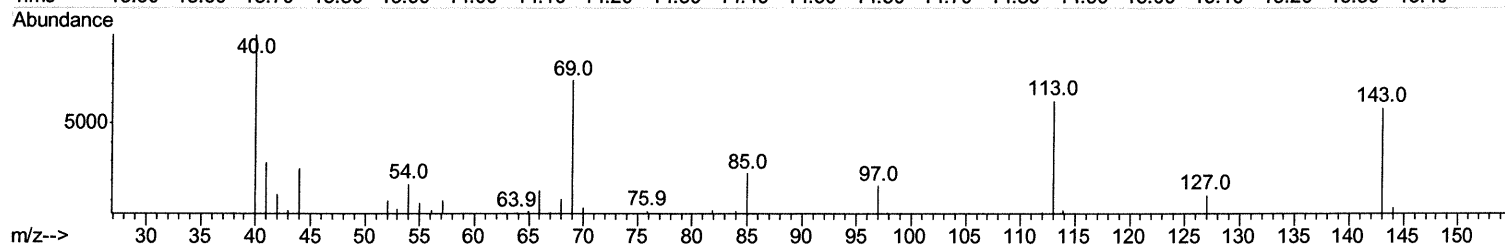
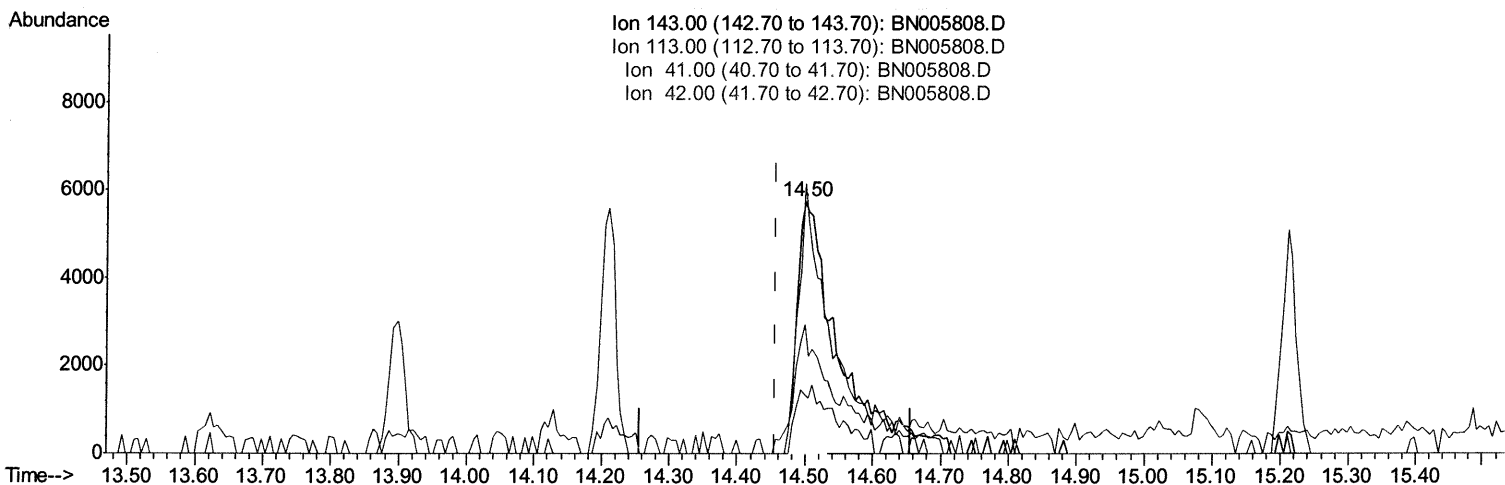
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Instrument :
 BNA_N
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Manual Integrations
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(51) 4-Nitrophenol-d4 (S)

14.498min (+0.041) 4.99ng/ul

response 14837

Ion	Exp%	Act%
143.00	100	100
113.00	82.50	107.09#
41.00	47.90	51.26
42.00	31.70	23.34#

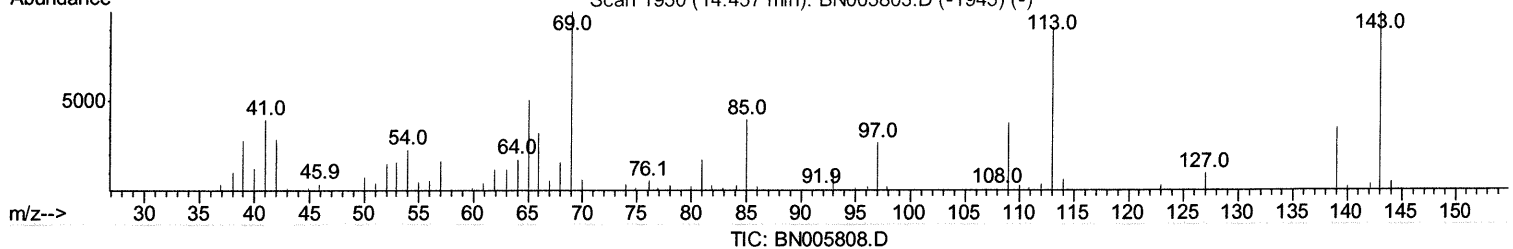
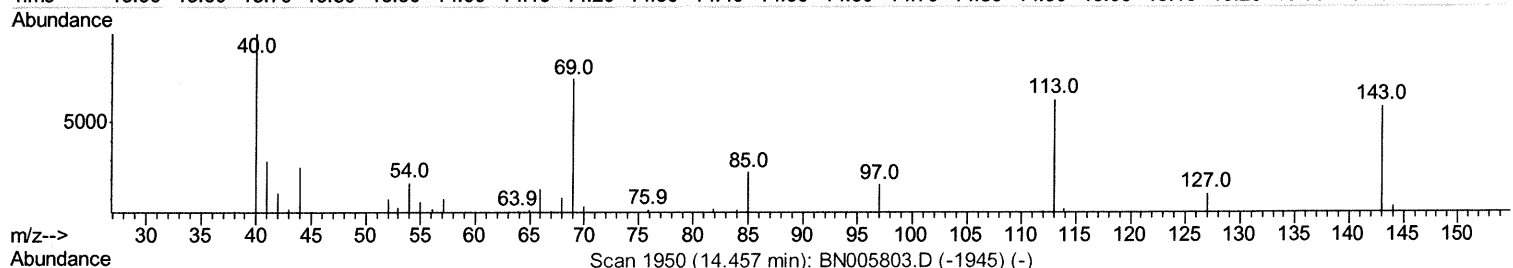
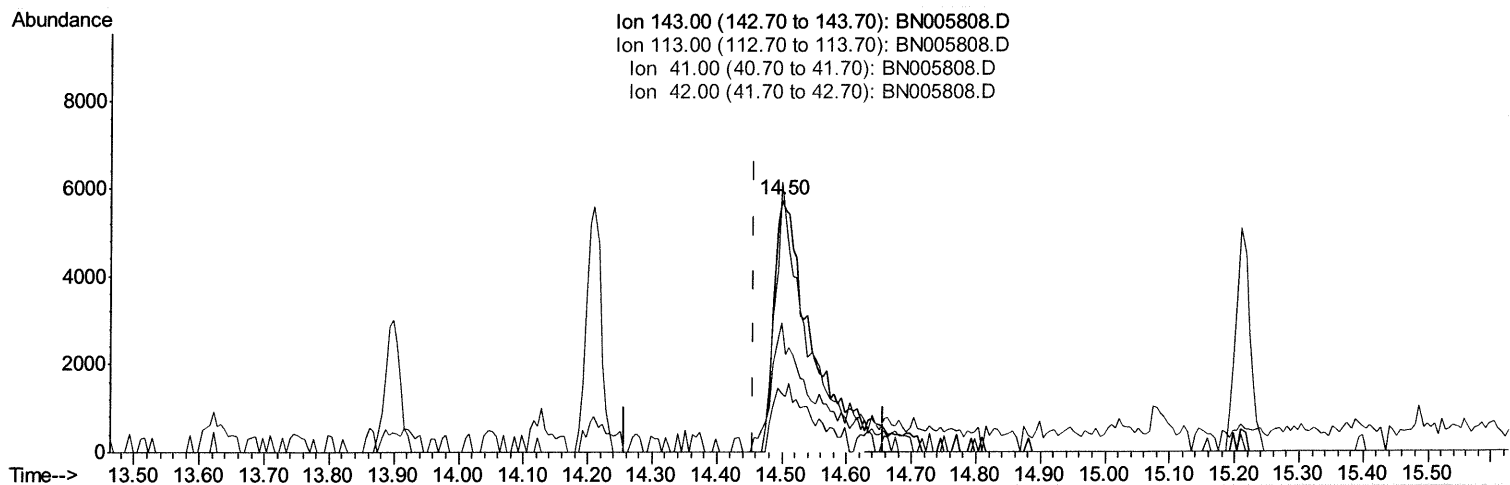
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual Integrations
APPROVED

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Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.498min (+0.041) 7.66ng/ul m ^{JU} 05/22/19

response 22783

Ion	Exp%	Act%
143.00	100	100
113.00	82.50	107.09#
41.00	47.90	51.26
42.00	31.70	23.34#

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	98118	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	404084	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.21	164	274526	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	723423	20.00	ng/ul	0.00
77) Chrysene-d12	21.23	240	653791	20.00	ng/ul	0.02
85) Perylene-d12	23.46	264	646933	20.00	ng/ul	0.04

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	4256	2.43	ng/uL	0.00
5) Phenol-d5	6.78	99	87674	11.56	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.92	67	56604	14.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.12	132	83723	13.73	ng/ul	0.00
13) 4-Methylphenol-d8	8.30	113	39200	6.05	ng/ul	0.02
19) Nitrobenzene-d5	8.73	128	44037	15.01	ng/ul	0.01
22) 2-Nitrophenol-d4	9.43	143	46488	13.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.99	165	86804	13.43	ng/ul	0.02
29) 4-Chloroaniline-d4	10.53	131	16178m>	2.25	ng/ul	0.05
43) Dimethylphthalate-d6	13.63	166	335389	15.88	ng/ul	0.00
46) Acenaphthylene-d8	13.90	160	420894	16.40	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	22783m>	7.66	ng/ul	0.04
57) Fluorene-d10	15.22	176	325376	17.94	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.39	200	5583	1.28	ng/ul	0.02
70) Anthracene-d10	17.07	188	536600	17.17	ng/ul	0.00
78) Pyrene-d10	19.39	212	644887	20.02	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.33	264	495896	17.09	ng/ul	0.04

JU 05/22/19

JU 05/22/19

Target Compounds

					Ovalue
14) Acetophenone	8.40	105	19630	1.891	ng/ul 96
44) Dimethylphthalate	13.67	163	76240	3.661	ng/ul 99
76) Fluoranthene	19.05	202	55856	1.260	ng/ul# 88
79) Pyrene	19.42	202	60477	1.506	ng/ul# 89
80) Butylbenzylphthalate	20.35	149	555283	33.760	ng/ul 92
83) Bis(2-ethylhexyl)phthalate	21.15	149	89059	3.816	ng/ul# 94
86) Di-n-octyl phthalate	22.05	149	80135	2.322	ng/ul 100
87) Benzo(b)fluoranthene	22.81	252	57257	1.609	ng/ul# 77

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