

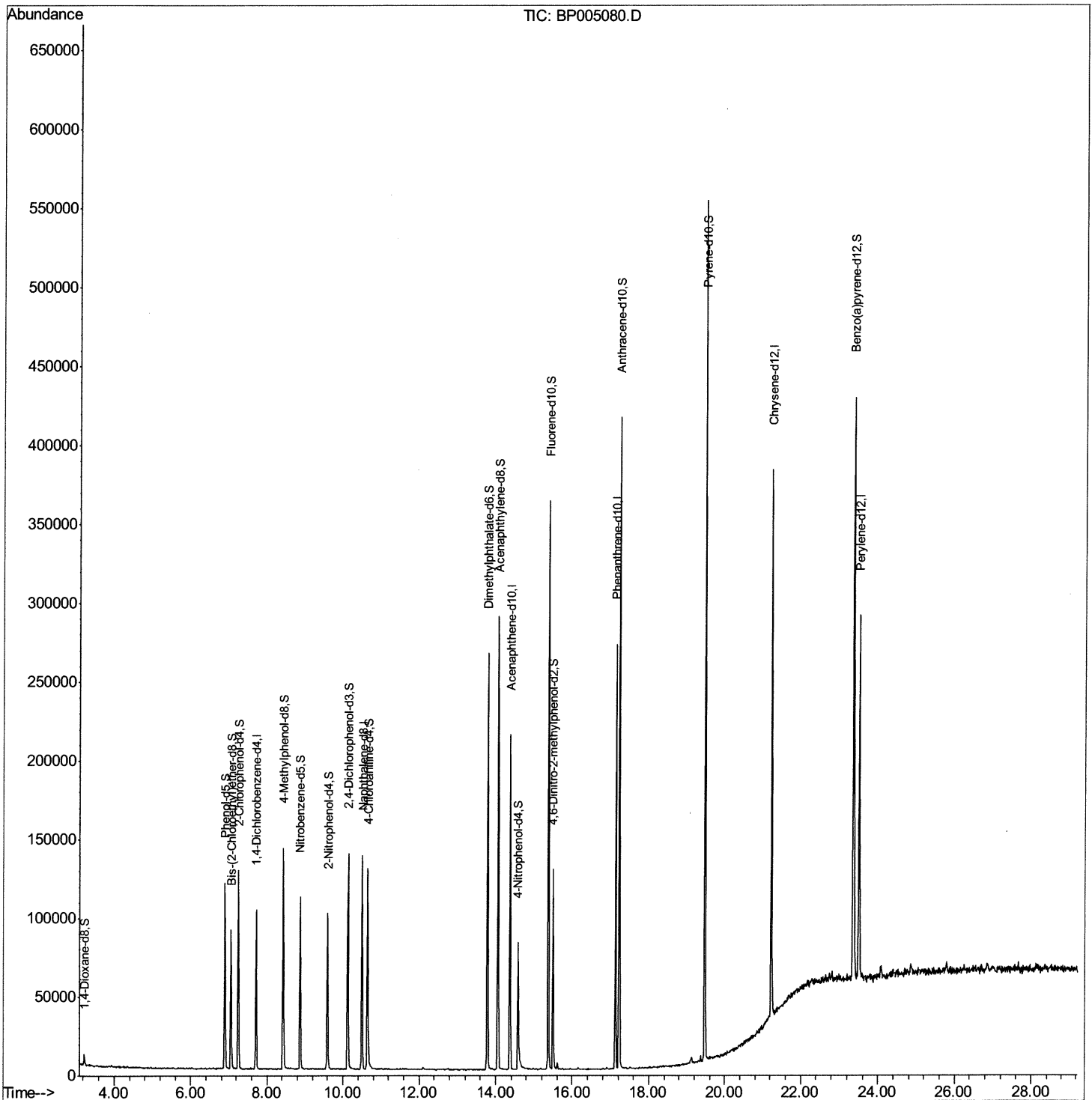
Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP032921\
Data File : BP005080.D
Acq On : 29 Mar 2021 18:31
Operator : CG/JU
Sample : PB135393BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK93

Manual Integrations
APPROVED

mohammad
3/30/2021 4:12:11 PM

Quant Time: Mar 29 19:02:34 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP032921MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 29 15:02:46 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

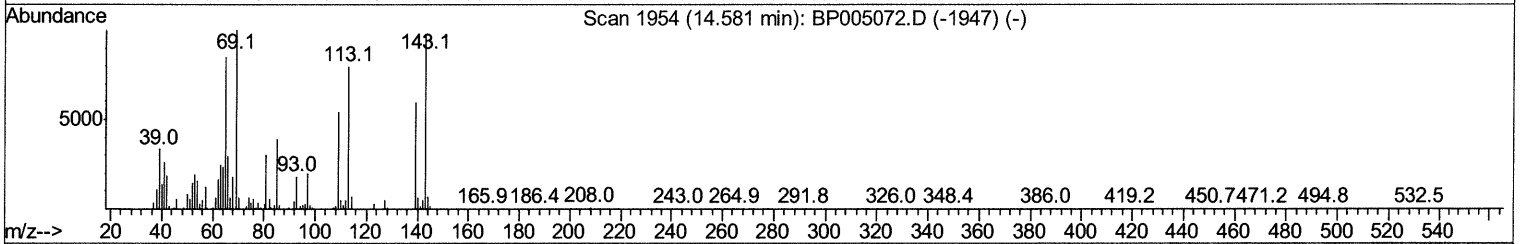
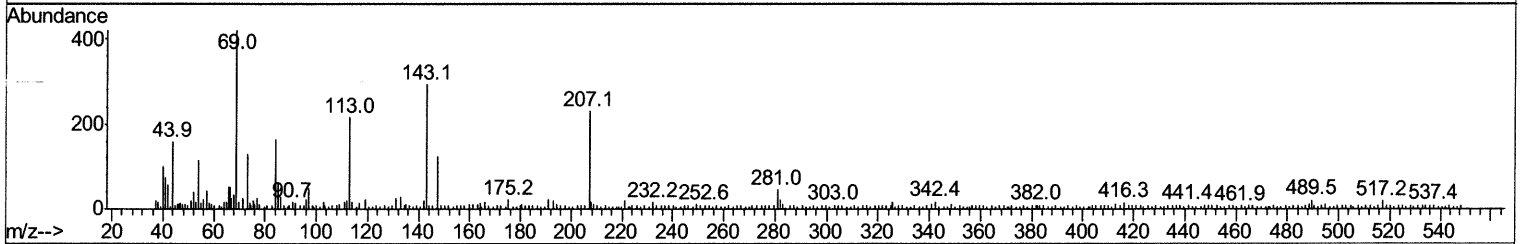
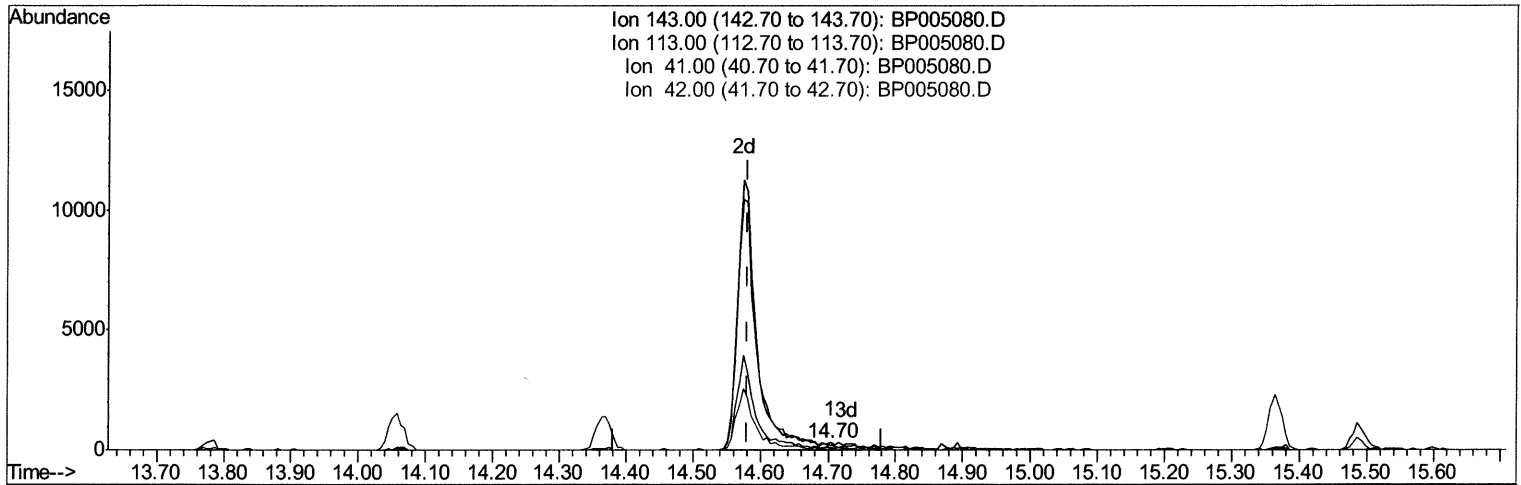
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TIC: BP005080.D

(52) 4-Nitrophenol-d4 (S)

14.704min (+0.124) 0.28ng/ul

response 227

Ion	Exp%	Act%
143.00	100	100
113.00	81.90	73.72
41.00	27.10	25.94
42.00	18.80	20.14

Quantitation Report (Qedit)

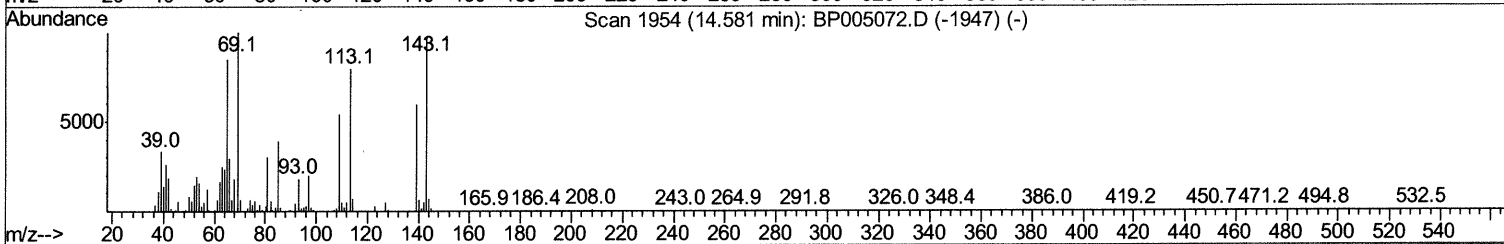
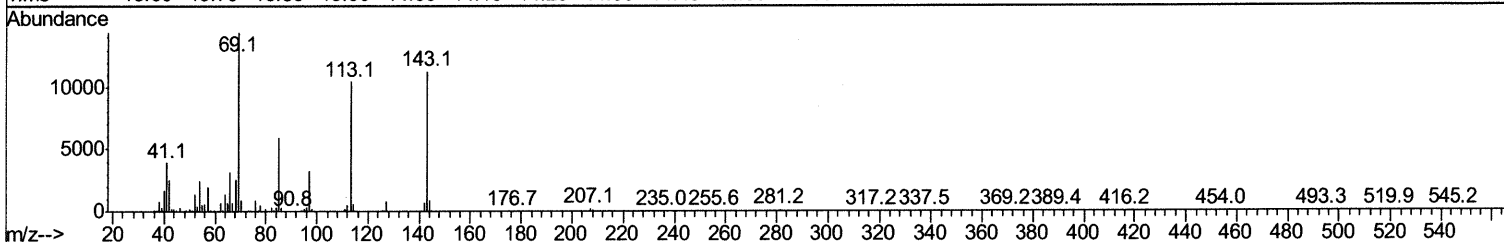
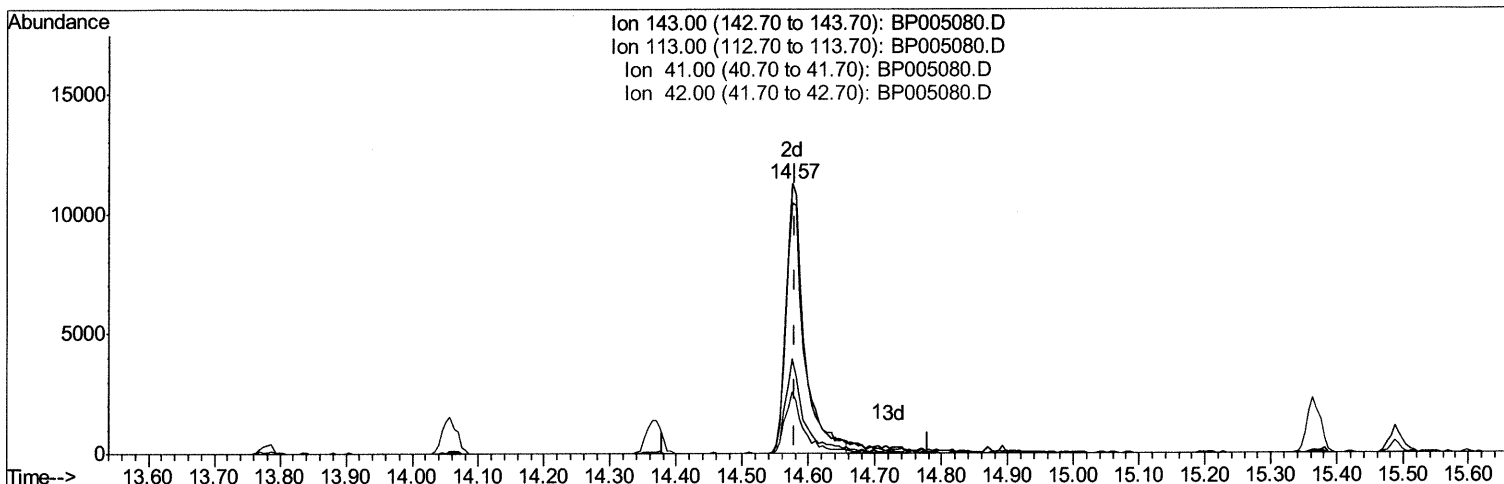
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 ClientSampleId :
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Manual Integrations
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 Response via : Initial Calibration



TIC: BP005080.D

(52) 4-Nitrophenol-d4 (S)

14.575min (-0.006) 26.28ng/ul m 04/01/2021

response 21542

Ion	Exp%	Act%
143.00	100	100
113.00	81.90	92.64
41.00	27.10	35.03#
42.00	18.80	22.73#

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP032921\
 Data File : BP005080.D
 Acq On : 29 Mar 2021 18:31
 Operator : CG/JU
 Sample : PB135393BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK93

Manual Integrations
 APPROVED

mohammad
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	26086	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	103790	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	69120	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.13	188	153194	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	159921	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	158123	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	4204	6.09	ng/uL	0.00
5) Phenol-d5	6.89	99	66416	29.56	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	38189	32.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	51017	32.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	51776	30.18	ng/ul	0.00
19) Nitrobenzene-d5	8.88	128	23940	31.24	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	27817	32.83	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	47784	28.90	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	65394	35.92	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	163870	32.65	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	190750	30.66	ng/ul	0.00
52) 4-Nitrophenol-d4	14.57	143	21542m>	26.28	ng/ul >	0.00
58) Fluorene-d10	15.37	176	136695	31.50	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	27378	27.06	ng/ul	0.00
71) Anthracene-d10	17.23	188	221529	32.12	ng/ul	0.00
79) Pyrene-d10	19.47	212	256874	33.57	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	274886	33.84	ng/ul	0.00

04/01/21

Target Compounds Ovalue

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