

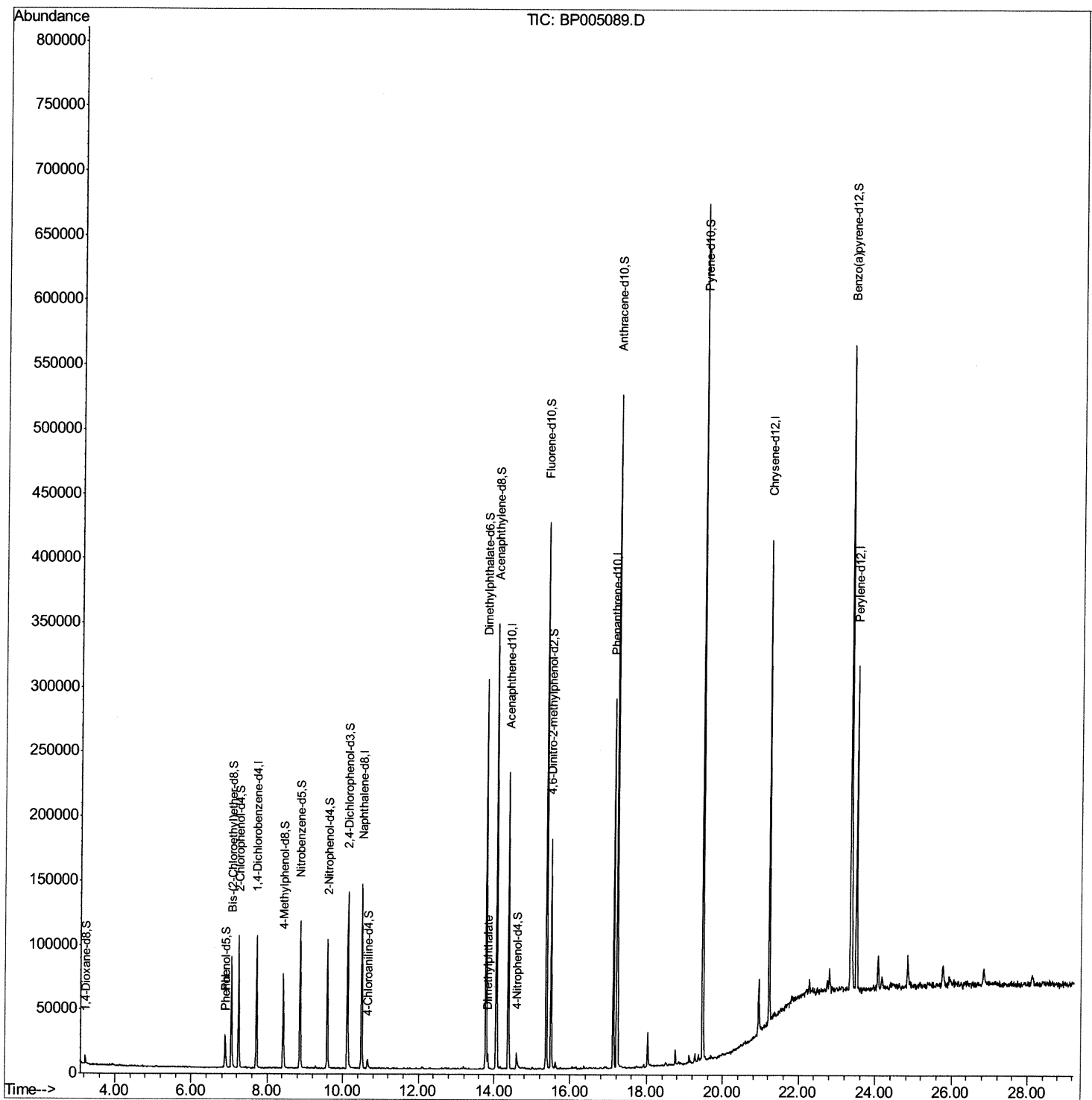
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032921\
Data File : BP005089.D
Acq On : 29 Mar 2021 23:36
Operator : CG/JU
Sample : M1818-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AA0

Manual Integrations
APPROVED

mohammad
3/30/2021 4:12:44 PM

Quant Time: Mar 30 02:36:41 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP032921MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 30 00:27:26 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

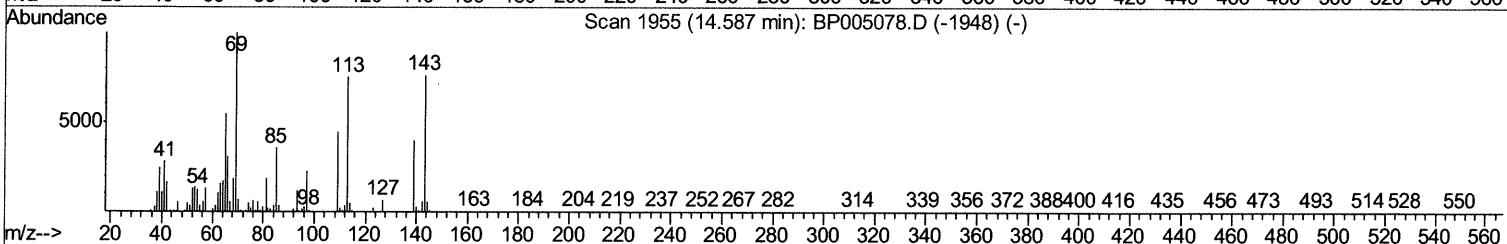
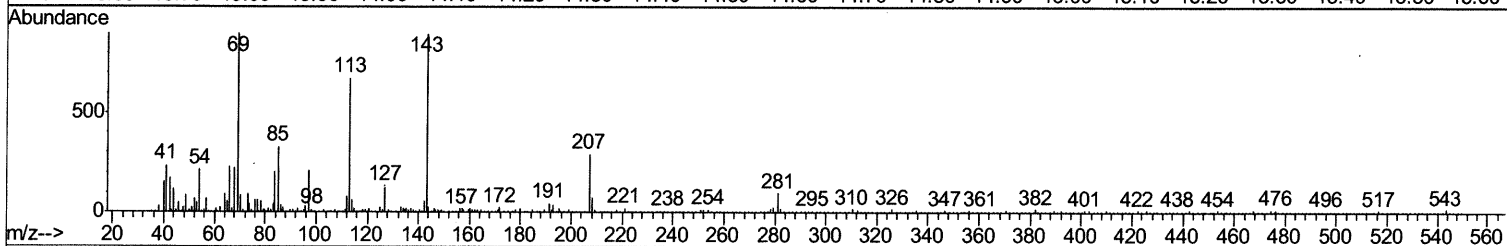
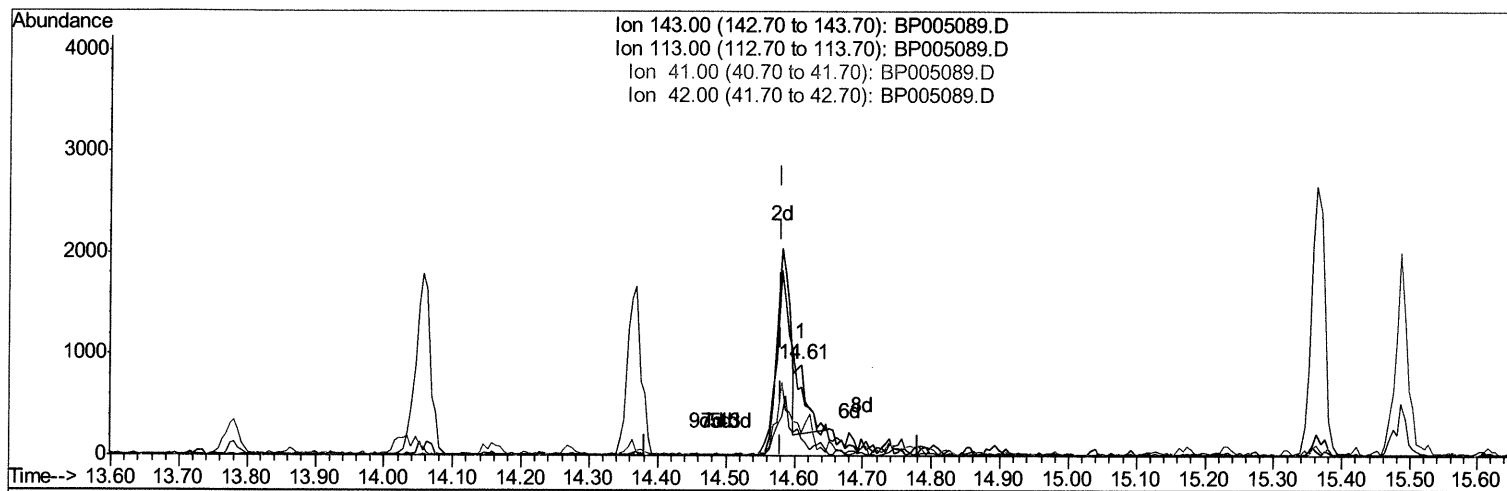
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TIC: BP005089.D

(52) 4-Nitrophenol-d4 (S)

14.610min (+0.029) 0.86ng/ul

response 771

Ion	Exp%	Act%
143.00	100	100
113.00	81.90	75.11
41.00	27.10	26.34
42.00	18.80	19.42

Quantitation Report (Qedit)

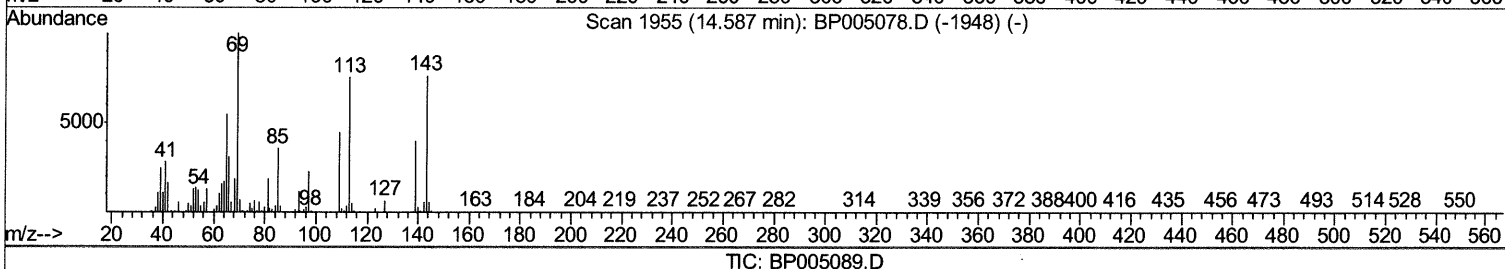
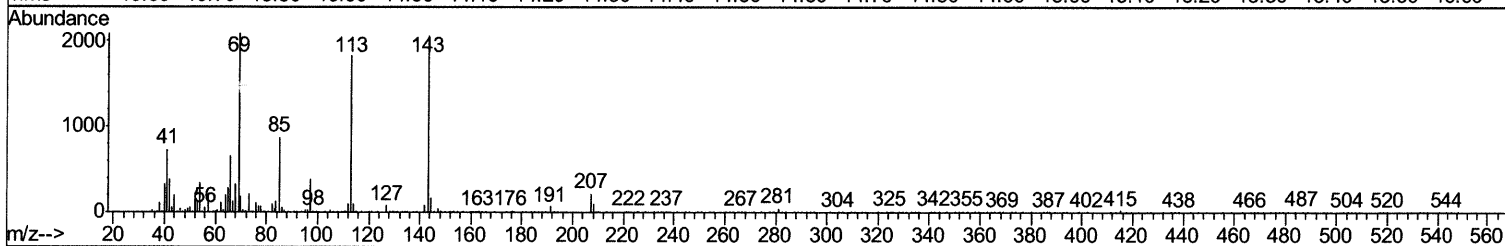
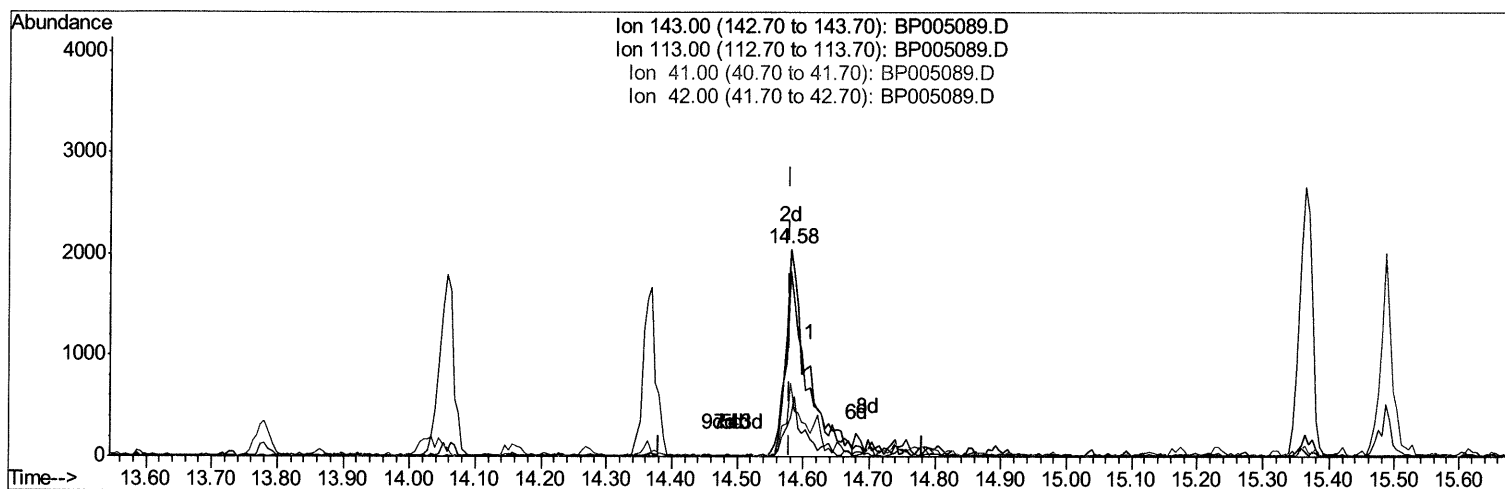
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Manual Integrations
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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BP005089.D

(52) 4-Nitrophenol-d4 (S)

14.580min (-0.000) 5.09ng/ul m 04/01/2020

response 4569

Ion	Exp%	Act%
143.00	100	100
113.00	81.90	89.13
41.00	27.10	35.62#
42.00	18.80	18.96

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032921\
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 Acq On : 29 Mar 2021 23:36
 Operator : CG/JU
 Sample : M1818-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA0

Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 30 00:27:26 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	27359	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	114745	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	75748	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.13	188	167828	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	178496	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	177305	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	3047	4.21	ng/uL	0.00
5) Phenol-d5	6.89	99	14718	6.25	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	37557	30.45	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	39463	23.77	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	26487	14.72	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	25145	29.68	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	28959	30.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	47533	26.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	4379	2.18	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	181856	33.06	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	218544	32.06	ng/ul	0.00
52) 4-Nitrophenol-d4	14.58	143	4569m >	5.09	ng/ul >	0.00
58) Fluorene-d10	15.36	176	161708	34.00	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	36927	33.31	ng/ul	0.00
71) Anthracene-d10	17.22	188	287201	38.01	ng/ul	0.00
79) Pyrene-d10	19.47	212	329689	38.61	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	359560	39.48	ng/ul	0.00

04/01/21 JU

Target Compounds

					Ovalue	
6) Phenol	6.91	94	2560	1.027	ng/ul#	95
45) Dimethylphthalate	13.83	163	6395	1.137	ng/ul	98

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