

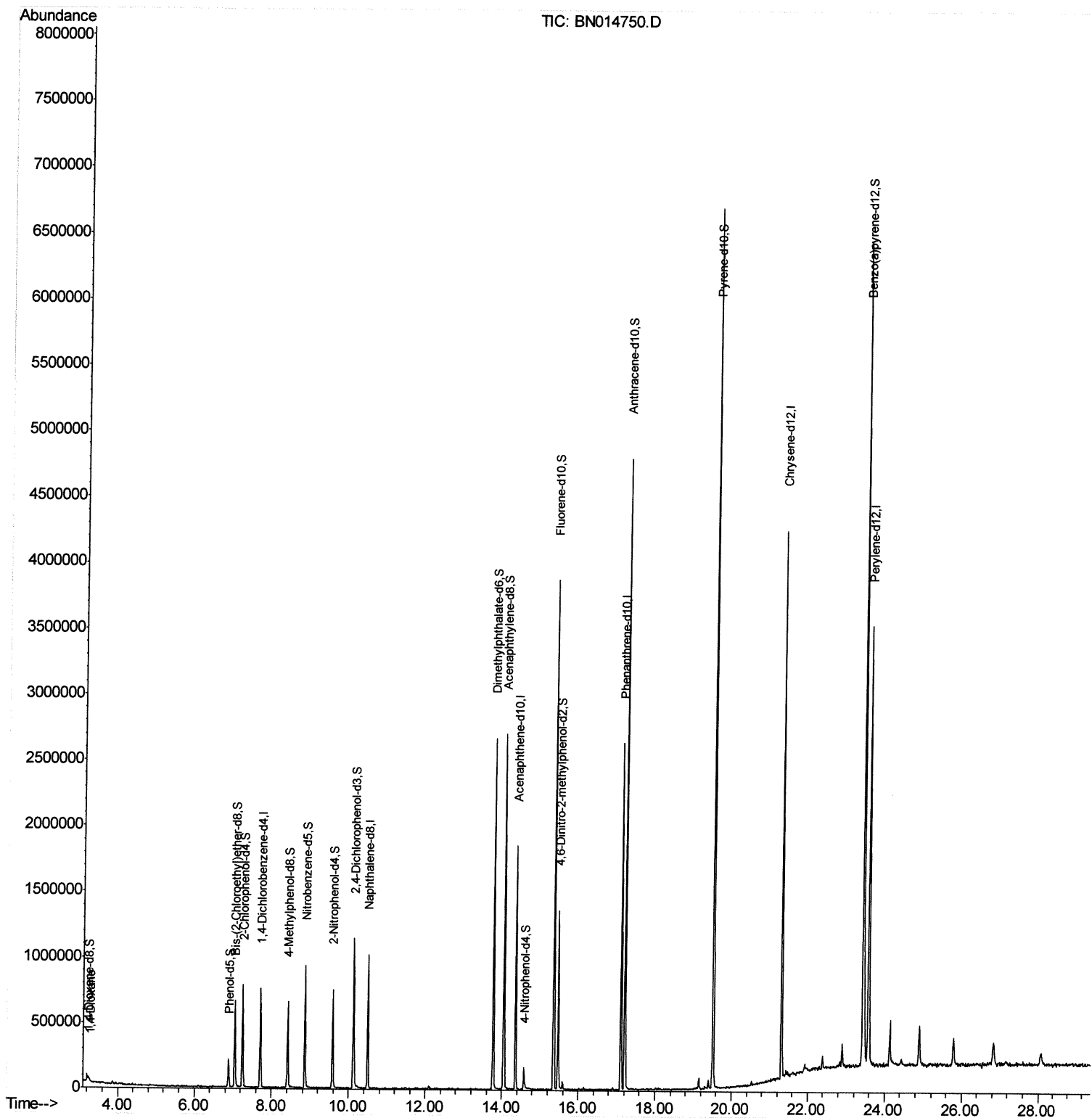
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN052521\
Data File : BN014750.D
Acq On : 26 May 2021 00:12
Operator : CG/JU
Sample : M2493-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GB417

Quant Time: May 26 03:23:22 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 25 10:33:04 2021
Response via : Initial Calibration

Manual Integrations
APPROVED

mohammad
5/26/2021 4:27:46 PM



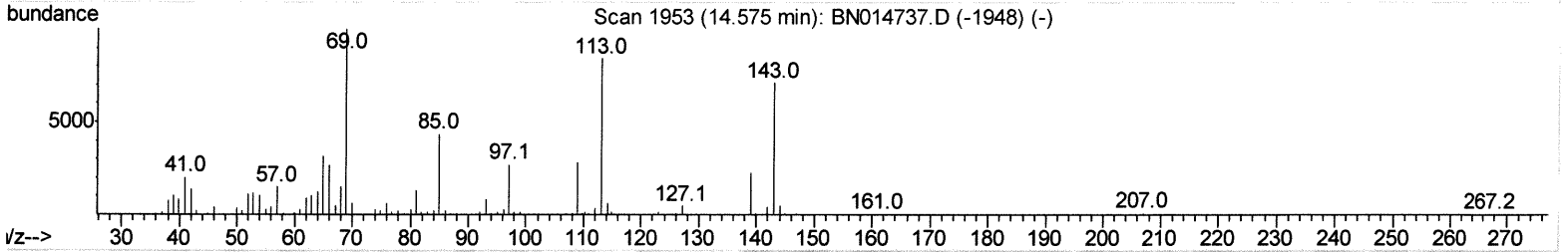
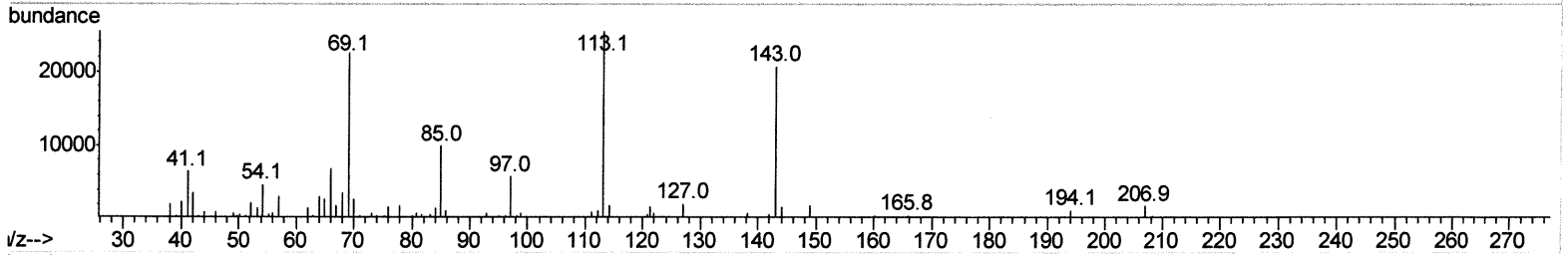
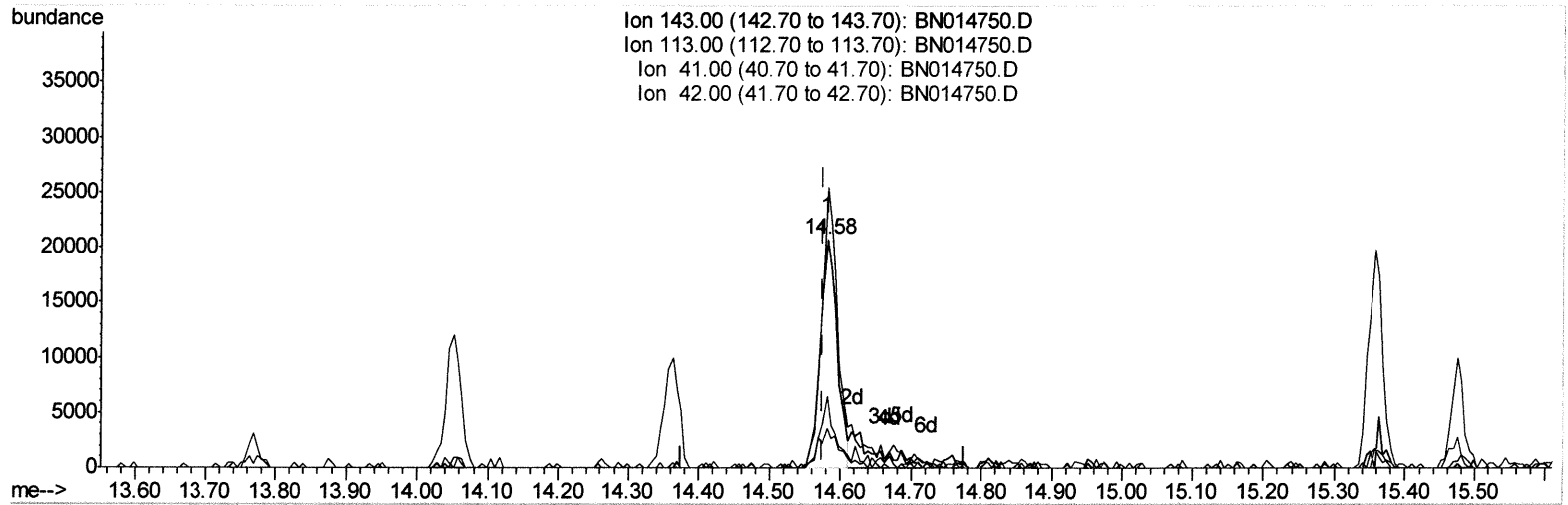
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(54) 4-Nitrophenol-d4 (S)

14.581min (+0.006) 4.57ng/ul

response 34123

Ion	Exp%	Act%
143.00	100	100
113.00	115.60	122.87
41.00	25.00	31.16#
42.00	19.00	17.54

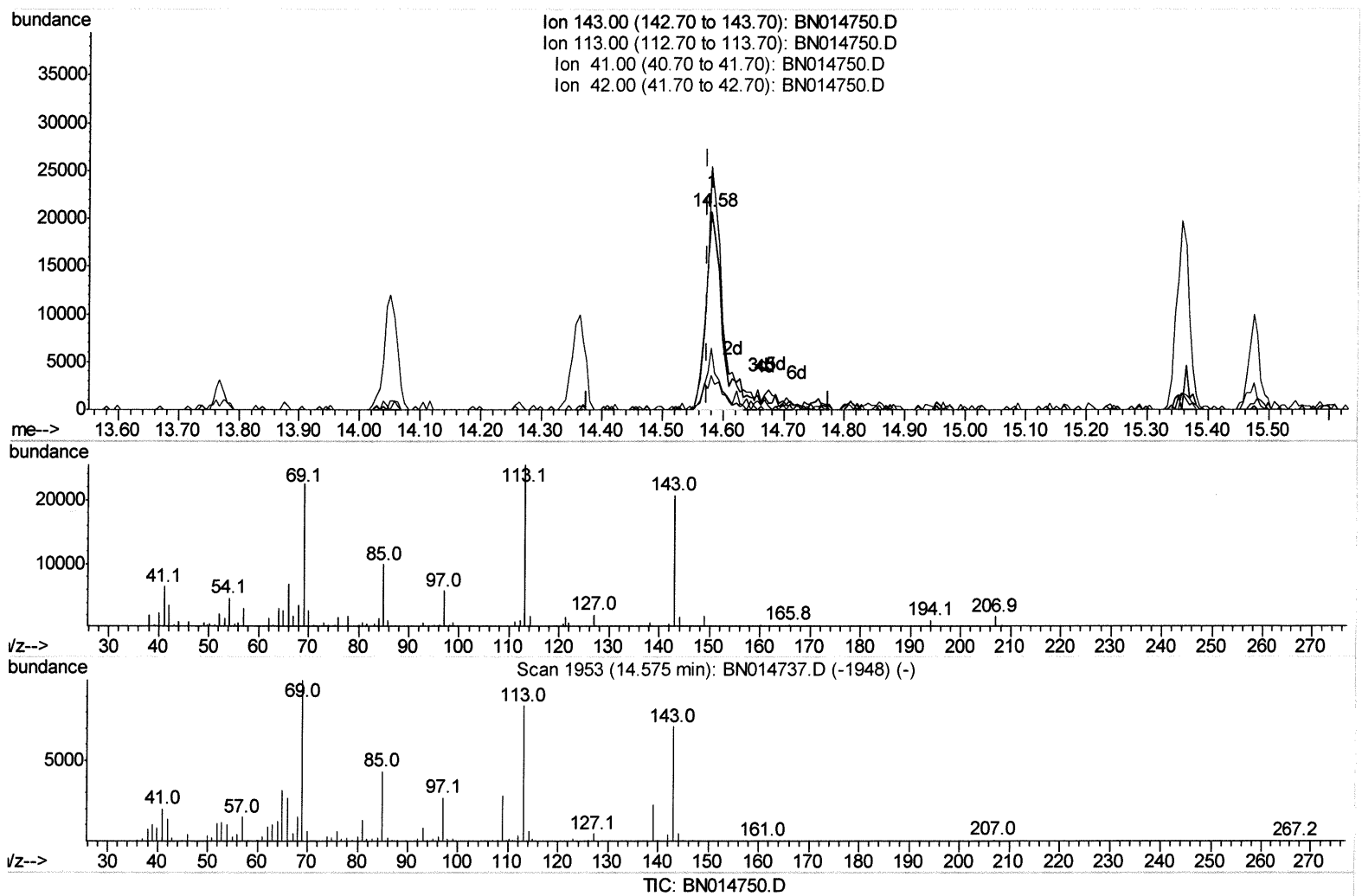
Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052521\
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 GB417

Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 25 10:33:04 2021
 Response via : Initial Calibration



(54) 4-Nitrophenol-d4 (S)

14.581min (+0.006) 5.08ng/ul m

response 37926

Ion	Exp%	Act%
143.00	100	100
113.00	115.60	122.87
41.00	25.00	31.16#
42.00	19.00	17.54

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN052521\
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Manual Integrations
 APPROVED

mohammad
 5/26/2021 4:27:46 PM

Quant Time: Mav 26 03:23:22 2021
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 25 10:33:04 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	187072	20.00	ng/ul	0.00
20) Naphthalene-d8	10.50	136	740458	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.36	164	591758	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.12	188	1483503	20.00	ng/ul	0.00
79) Chrysene-d12	21.31	240	2080684	20.00	ng/ul	0.00
88) Perylene-d12	23.57	264	2421191	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.23	96	15653	3.41	ng/uL	0.00
4) Pividine-d5	0.00	84	0d	0.00	ng/ul	0.00
7) Phenol-d5	6.89	99	126466	6.90	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.05	67	275647	30.26	ng/ul	0.00
11) 2-Chlorophenol-d4	7.25	132	300481	26.49	ng/ul	0.00
15) 4-Methylphenol-d8	8.42	113	224026	15.73	ng/ul	0.00
21) Nitrobenzene-d5	8.86	128	185441	33.25	ng/ul	0.00
24) 2-Nitrophenol-d4	9.58	143	190206	31.73	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.12	165	386169	28.25	ng/ul	0.00
31) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
46) Dimethylphthalate-d6	13.77	166	1636876	34.82	ng/ul	0.00
49) Acenaphthylene-d8	14.05	160	1781771	32.77	ng/ul	0.00
54) 4-Nitrophenol-d4	14.58	143	37926m	5.08	ng/ul	0.00
60) Fluorene-d10	15.36	176	1509234	37.26	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.47	200	260855	32.15	ng/ul	0.00
73) Anthracene-d10	17.22	188	2727569	39.36	ng/ul	0.00
81) Pyrene-d10	19.52	212	3587519	36.57	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.43	264	4849712	37.60	ng/ul	0.00

Target Compounds				Ovalue	
2) 1,4-Dioxane	3.26	88	12062	2.236	ng/uL# 81

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