

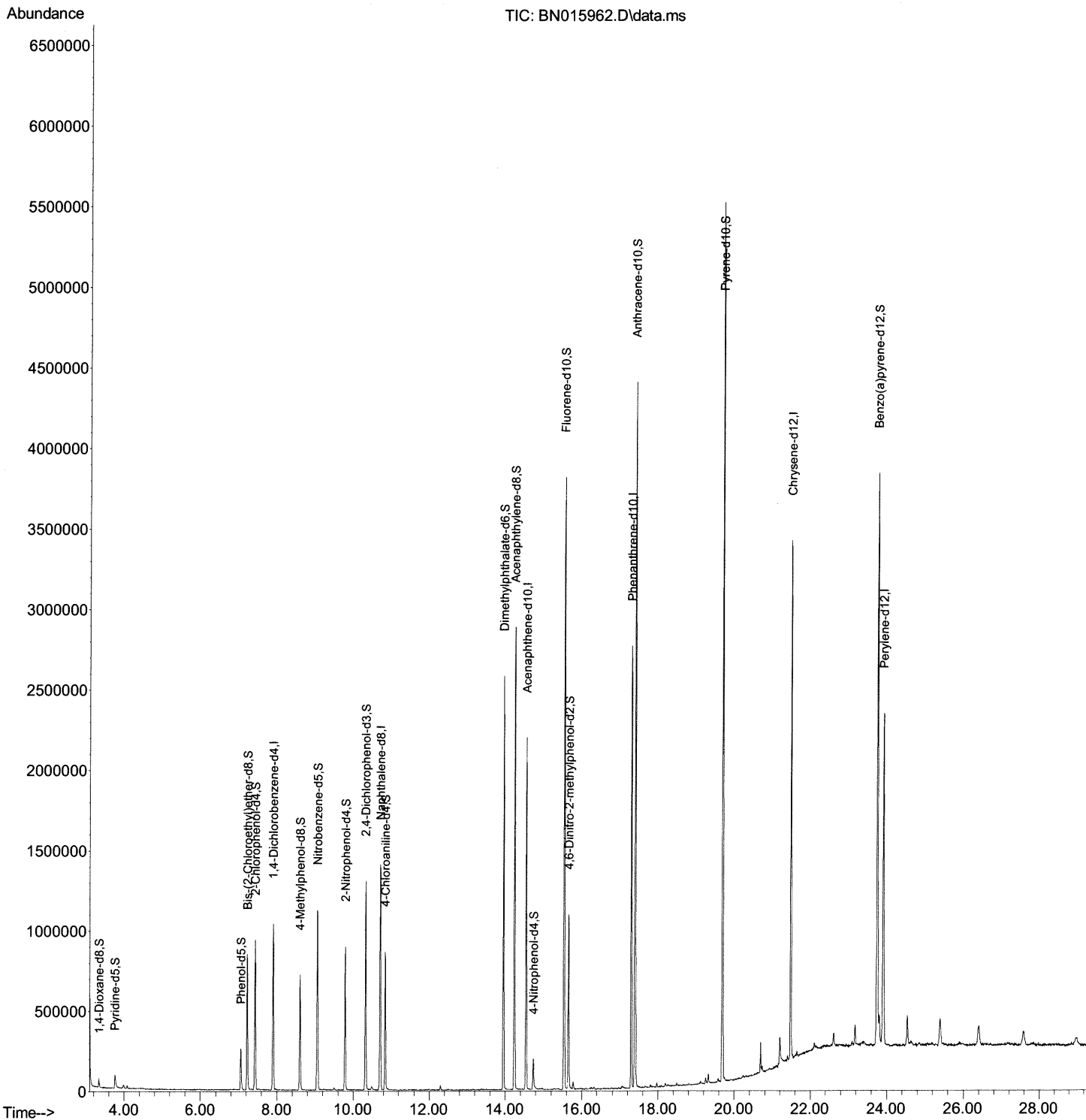
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
Data File : BN015962.D
Acq On : 19 Aug 2021 07:31
Operator : CG/JU
Sample : M3399-20
Misc :
ALS Vial : 68 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBKD2

Manual Integrations
APPROVED

mohammad
8/19/2021 1:26:17 PM

Quant Time: Aug 19 09:01:17 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 17 09:43:53 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

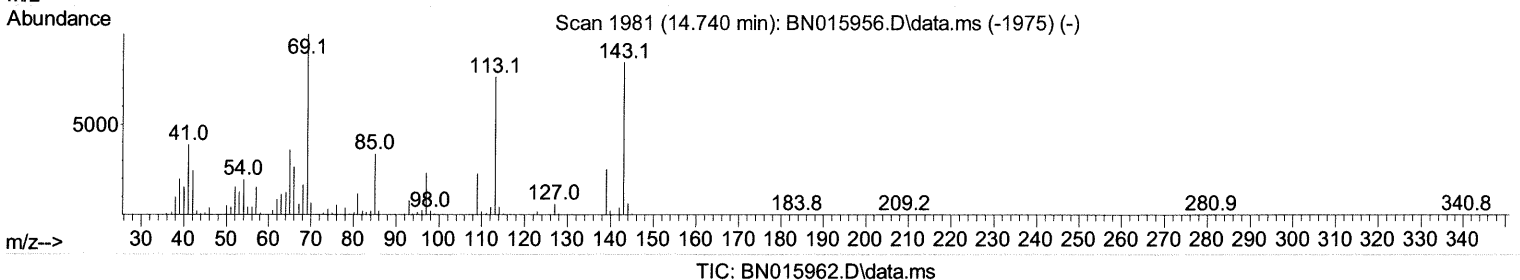
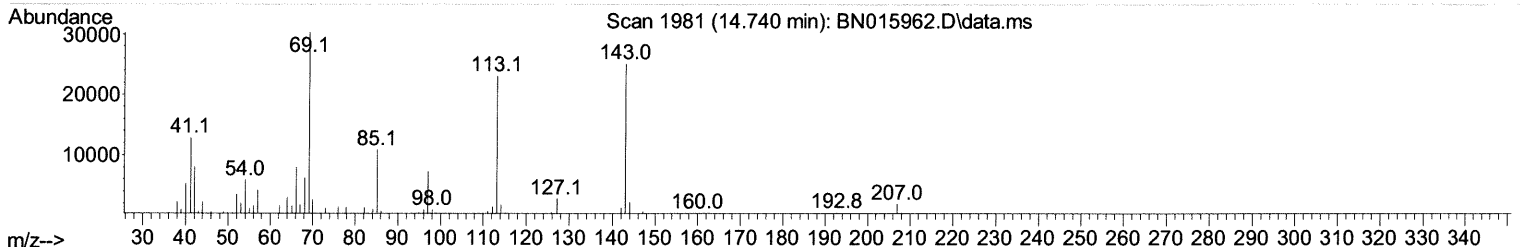
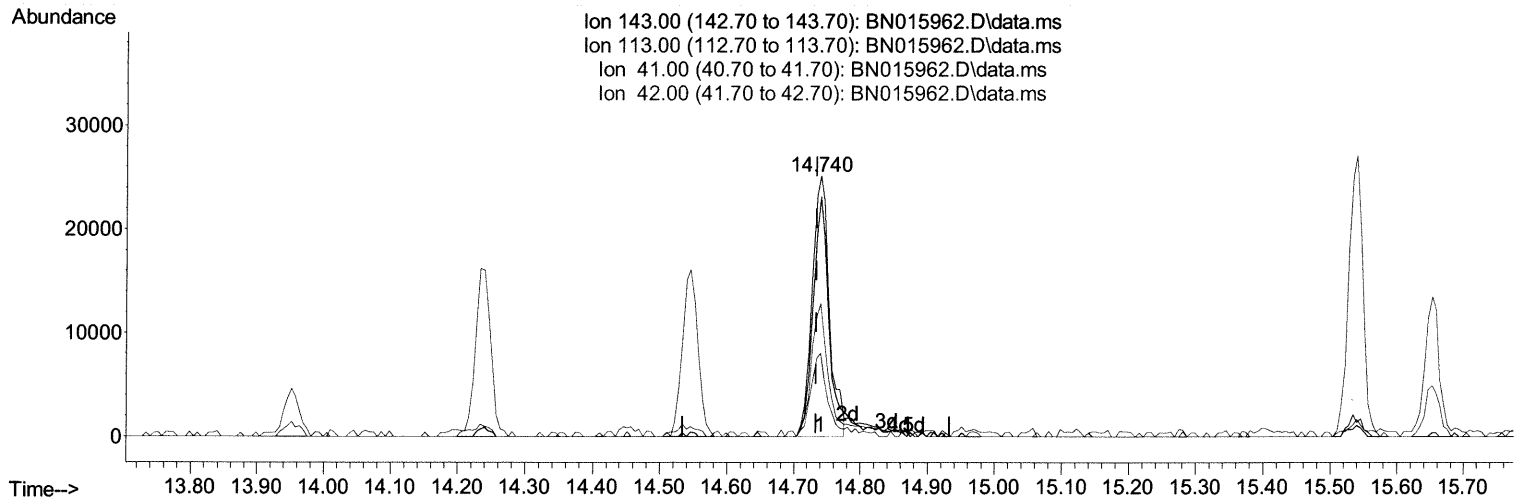
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(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.006) 4.66 ng/u1

response 45688

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	92.00	92.00
41.00	42.20	51.17#
42.00	29.90	32.08

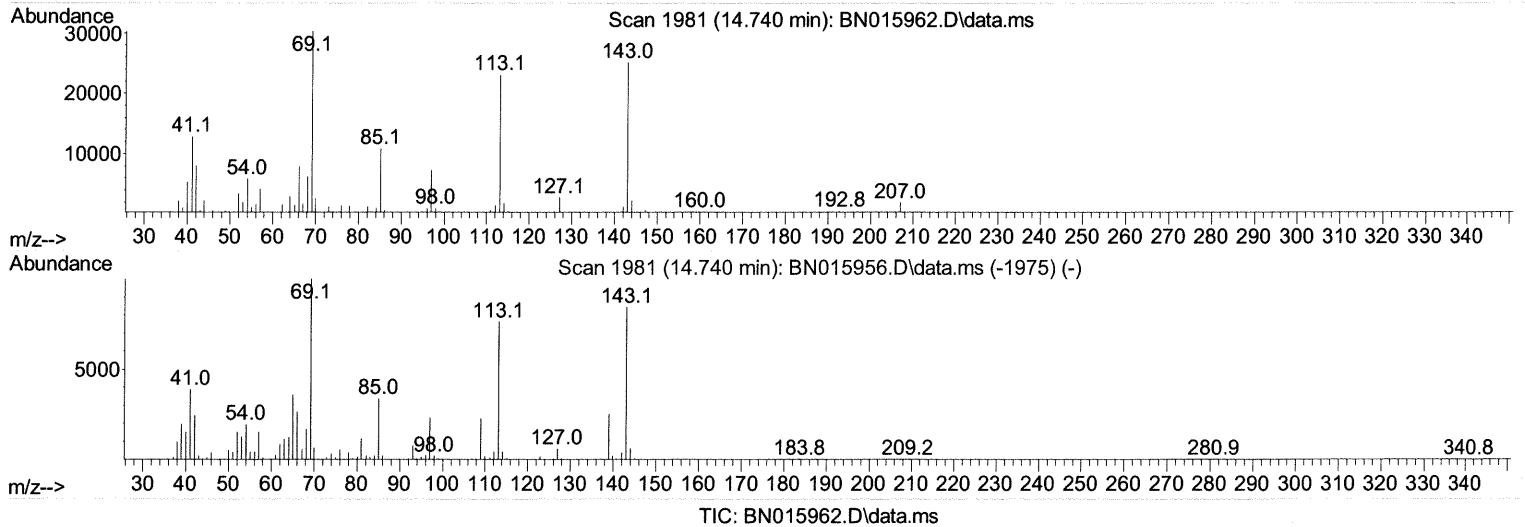
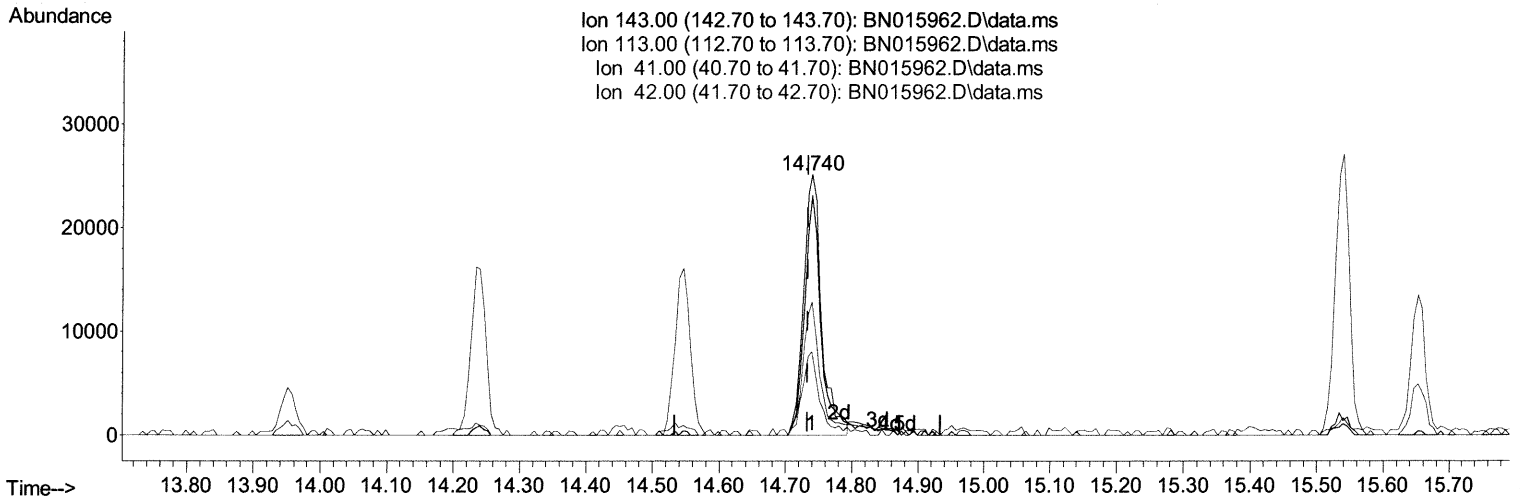
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Data File : BN015962.D
Acq On    : 19 Aug 2021   07:31
Operator  : CG/JU
Sample    : M3399-20
Misc      :
ALS Vial  : 68   Sample Multiplier: 1
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Instrument :
BNA_N
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DBKD2

Manual Integrations

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Response via : Initial Calibration



(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.006) 4.83 ng/ul m 08/20/21 JU

response 47353

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	92.00	92.00
41.00	42.20	51.17#
42.00	29.90	32.08

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081721\
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 Sample : M3399-20
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 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBKD2

Manual Integrations
 APPROVED

mohammad
 8/19/2021 1:26:17 PM

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	277012	20.000	ng/ul	0.00
20) Naphthalene-d8	10.705	136	1168030	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.546	164	755599	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	1608228	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	1550239	20.000	ng/ul	0.00
88) Perylene-d12	23.904	264	1539594	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	32672	4.594	ng/ul	0.00
4) Pyridine-d5	3.758	84	53783	2.706	ng/ul	0.01
7) Phenol-d5	7.058	99	142740	5.697	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	397270	25.750	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	395922	22.205	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	256568	13.165	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128	250836	29.178	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	264509	32.875	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	447942	25.728	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	436740	16.479	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	1662305	30.019	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1983742	28.605	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	47353m	4.835	ng/ul	0.00
60) Fluorene-d10	15.540	176	1439665	29.805	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	213957	24.911	ng/ul	0.00
73) Anthracene-d10	17.392	188	2519933	33.430	ng/ul	0.00
81) Pyrene-d10	19.686	212	2863670	34.295	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	2726551	32.754	ng/ul	0.00

Target Compounds Qvalue

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