

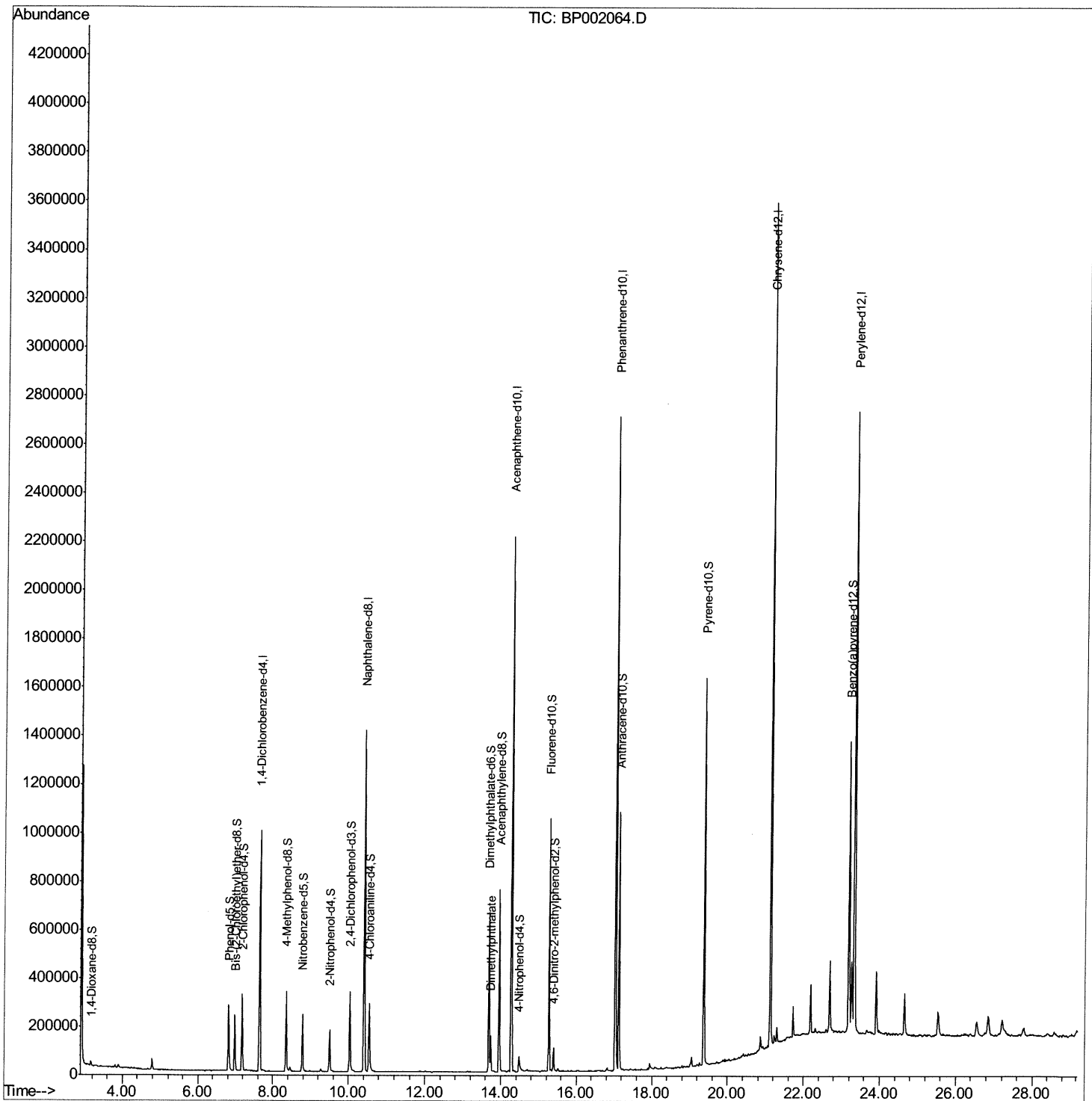
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP030420\
Data File : BP002064.D
Acq On : 04 Mar 2020 23:40
Operator : CG/JU
Sample : L1677-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CD8

Manual Integrations
APPROVED

mohammad
3/5/2020 3:55:11 PM

Quant Time: Mar 05 11:29:20 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 05 06:21:29 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

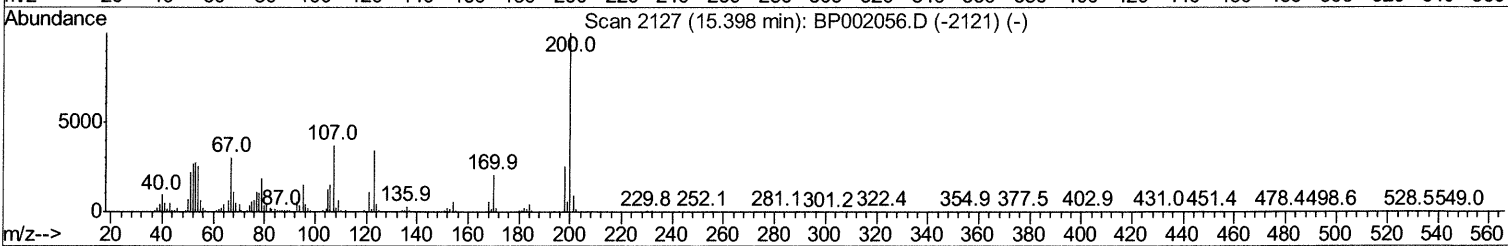
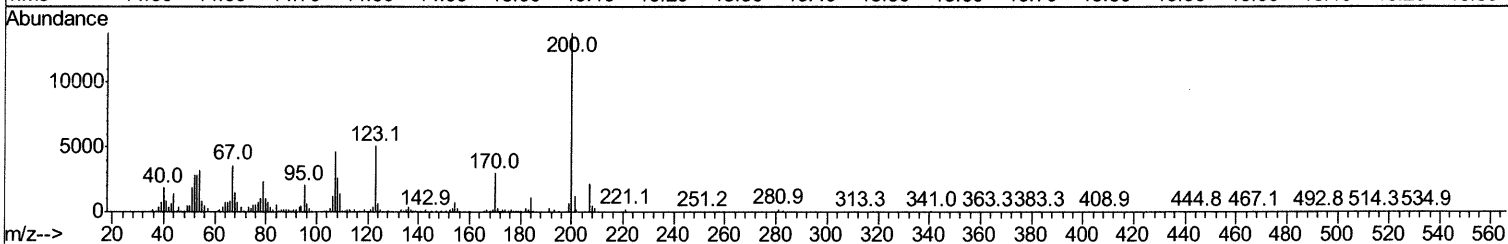
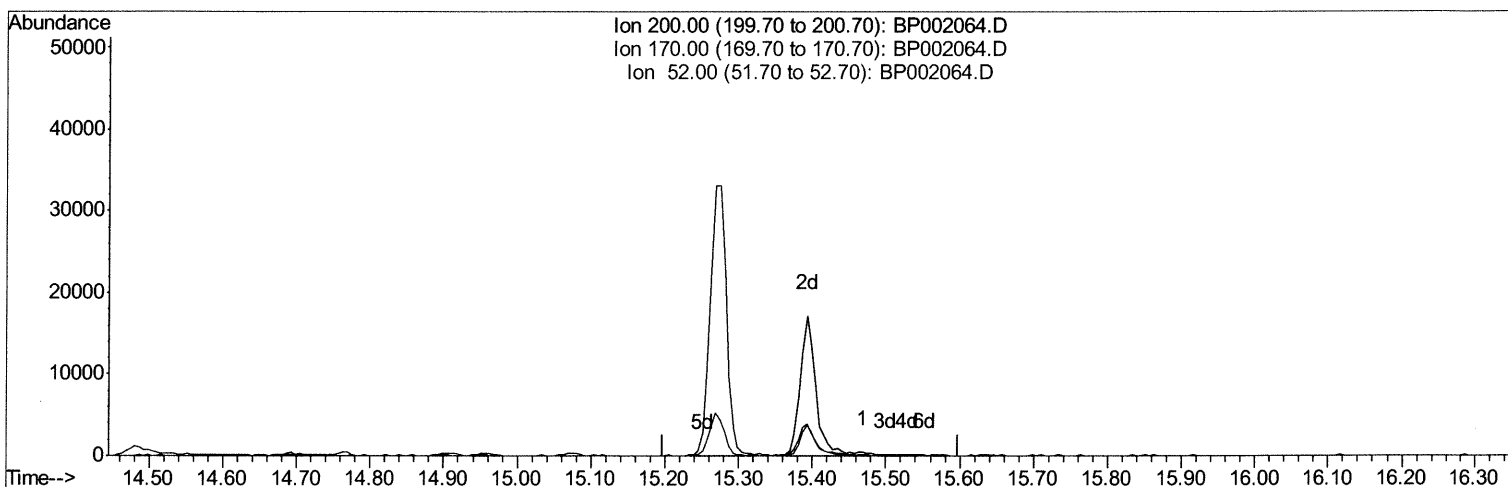
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TIC: BP002064.D

(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.398min (-15.398) 0.00ng/ul d

response 0

Ion	Exp%	Act%
200.00	100	0.00
170.00	19.70	0.00
52.00	29.70	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

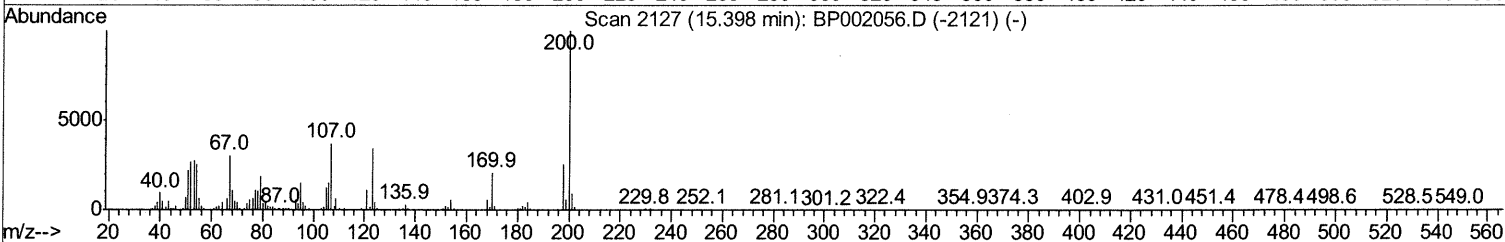
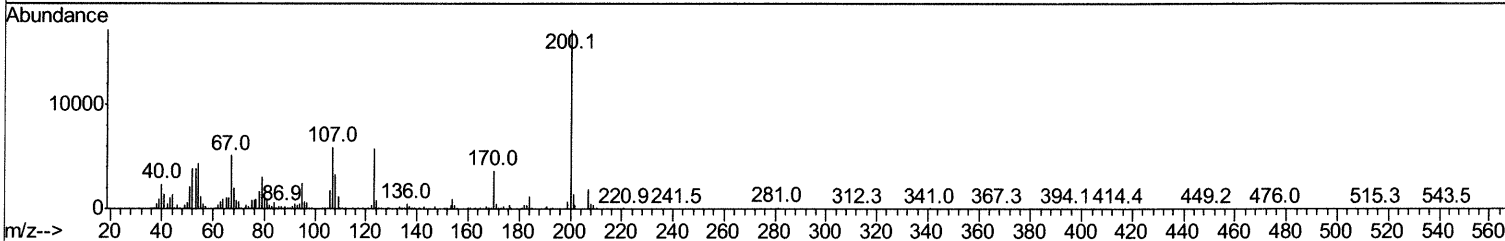
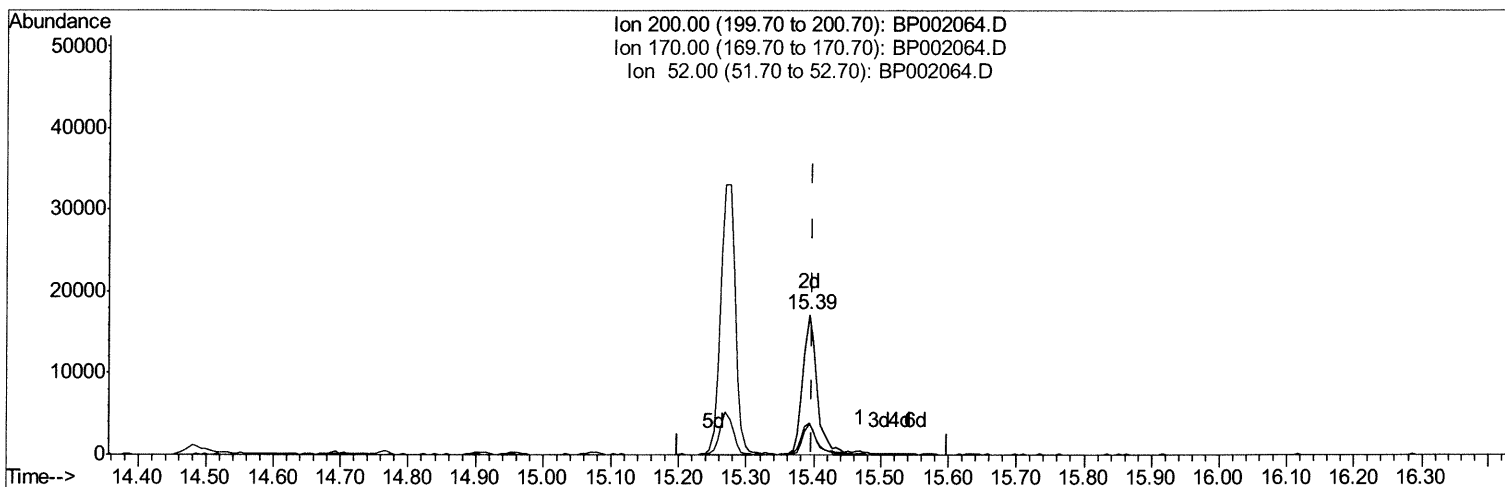
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Manual Integrations
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Response via : Initial Calibration



TIC: BP002064.D

(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.393min (-0.006) 2.72ng/ul m 2403/05/20

response 24957

Ion	Exp%	Act%
200.00	100	100
170.00	19.70	21.52
52.00	29.70	22.53#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP030420\
 Data File : BP002064.D
 Acq On : 04 Mar 2020 23:40
 Operator : CG/JU
 Sample : L1677-16
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CD8

Manual Integrations
 APPROVED

mohammad
 3/5/2020 3:55:11 PM

Quant Time: Mar 05 11:29:20 2020
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	293631	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1215711	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	774047	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1662245	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1636266	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1855547	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	9647	1.41	ng/uL	0.00
5) Phenol-d5	6.82	99	167089	7.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	100058	8.44	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	154857	7.99	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	134714	7.32	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	64806	6.94	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	64890	6.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	145375	7.18	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	186153	8.39	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	462338	7.89	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	547015	7.81	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	28750	2.57	ng/ul	0.00
58) Fluorene-d10	15.27	176	428437	8.45	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	24957m>	2.72	ng/ul>	0.00
71) Anthracene-d10	17.12	188	645864	8.65	ng/ul	0.00
79) Pyrene-d10	19.37	212	816383	9.53	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	879196	9.26	ng/ul	-0.01

8403/05/20

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.73	163	102121	1.689	ng/ul	99

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