

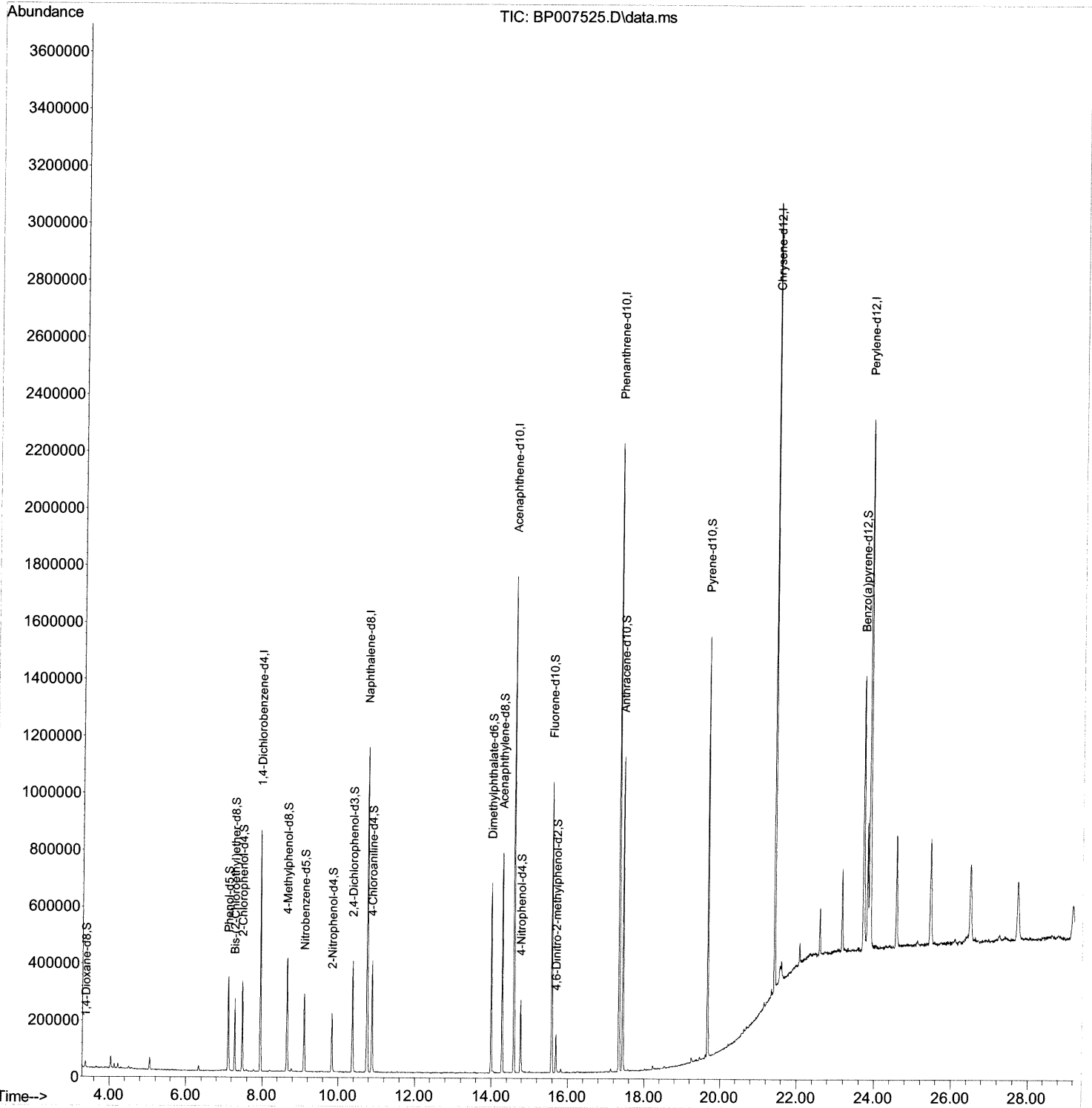
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007525.D
Acq On : 20 Oct 2021 13:01
Operator : CG/JU
Sample : M4196-14
Misc :
ALS Vial : 80 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGD0

Manual Integrations
APPROVED

mohammad
10/21/2021 11:15:40 AM

Quant Time: Oct 20 14:06:01 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 20 10:49:33 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

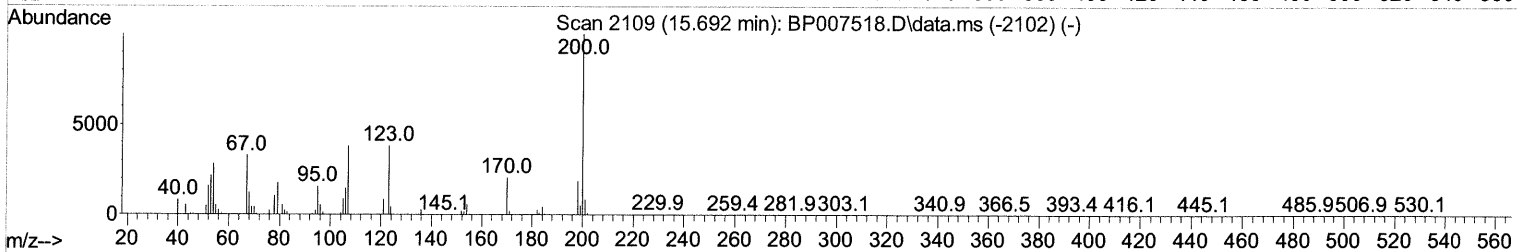
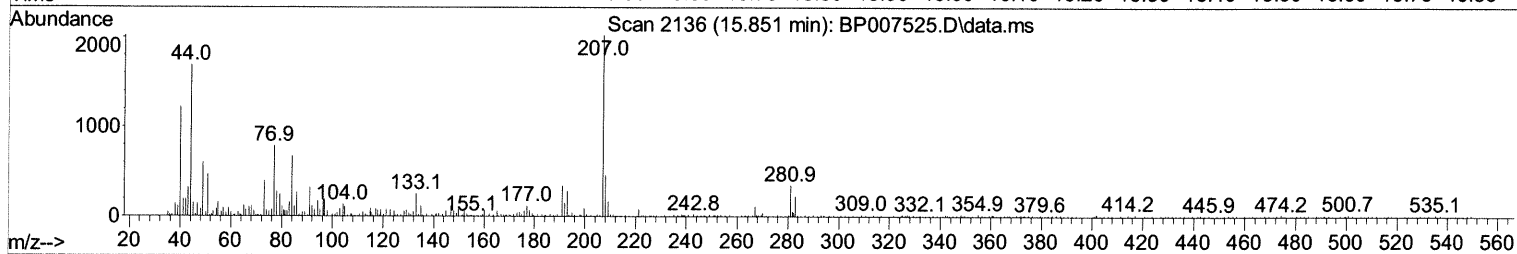
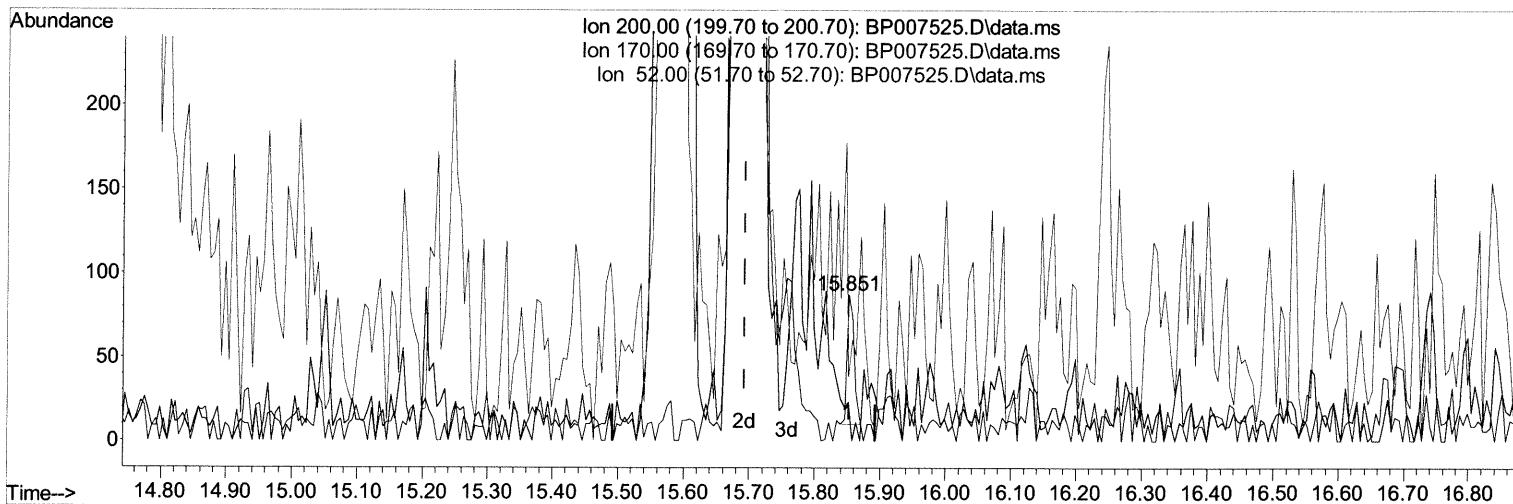
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TIC: BP007525.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.851min (+ 0.159) 0.01 ng/ul

response 58

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	20.60	21.30
52.00	40.30	35.19
0.00	0.00	0.00

Quantitation Report (Qedit)

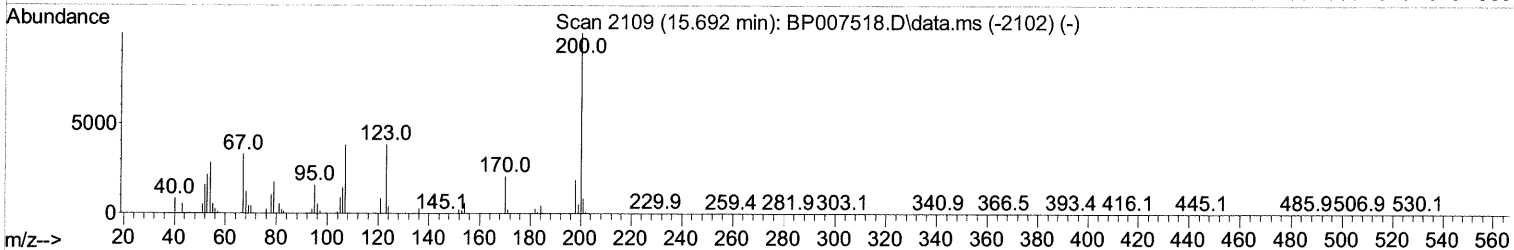
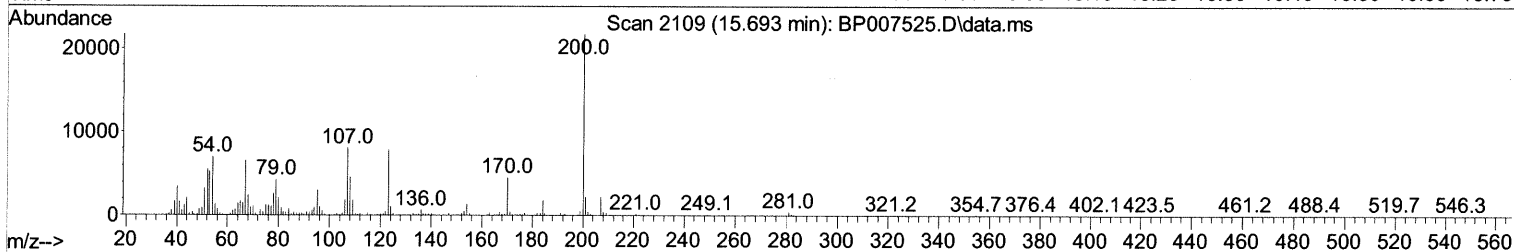
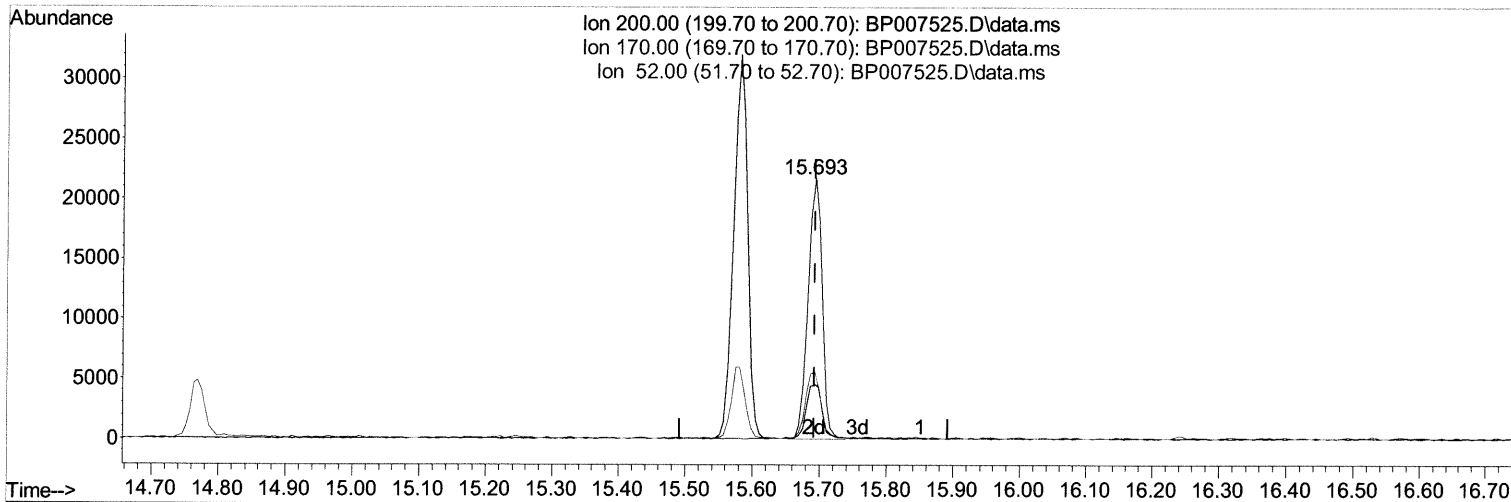
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TIC: BP007525.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.693min (+ 0.000) 5.86 ng/ul m 10/22/21 JU

response 31187

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	20.60	20.79
52.00	40.30	25.58#
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : M4196-14
 Misc :
 ALS Vial : 80 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual Integrations
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 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	233288	20.000	ng/ul	0.00
20) Naphthalene-d8	10.758	136	945574	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	615965	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.340	188	1327931	20.000	ng/ul	0.00
79) Chrysene-d12	21.422	240	1399349	20.000	ng/ul	0.00
88) Perylene-d12	23.874	264	1460260	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	10739	1.818	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	0.00
7) Phenol-d5	7.111	99	188312	10.336	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	108435	9.979	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	146091	10.193	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	151430	10.791	ng/ul	0.00
21) Nitrobenzene-d5	9.111	128	65788	10.994	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	65090	11.645	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	142575	10.458	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	205153	11.128	ng/ul	0.00
46) Dimethylphthalate-d6	13.999	166	437674	10.270	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	538075	10.047	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	59094	8.642	ng/ul	0.00
60) Fluorene-d10	15.581	176	391533	10.979	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.693	200	31187m	5.861	ng/ul	0.00
73) Anthracene-d10	17.439	188	611916	10.851	ng/ul	0.00
81) Pyrene-d10	19.669	212	741655	10.602	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	707687	9.740	ng/ul	0.00

10/22/21 JU

Target Compounds Qvalue

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