

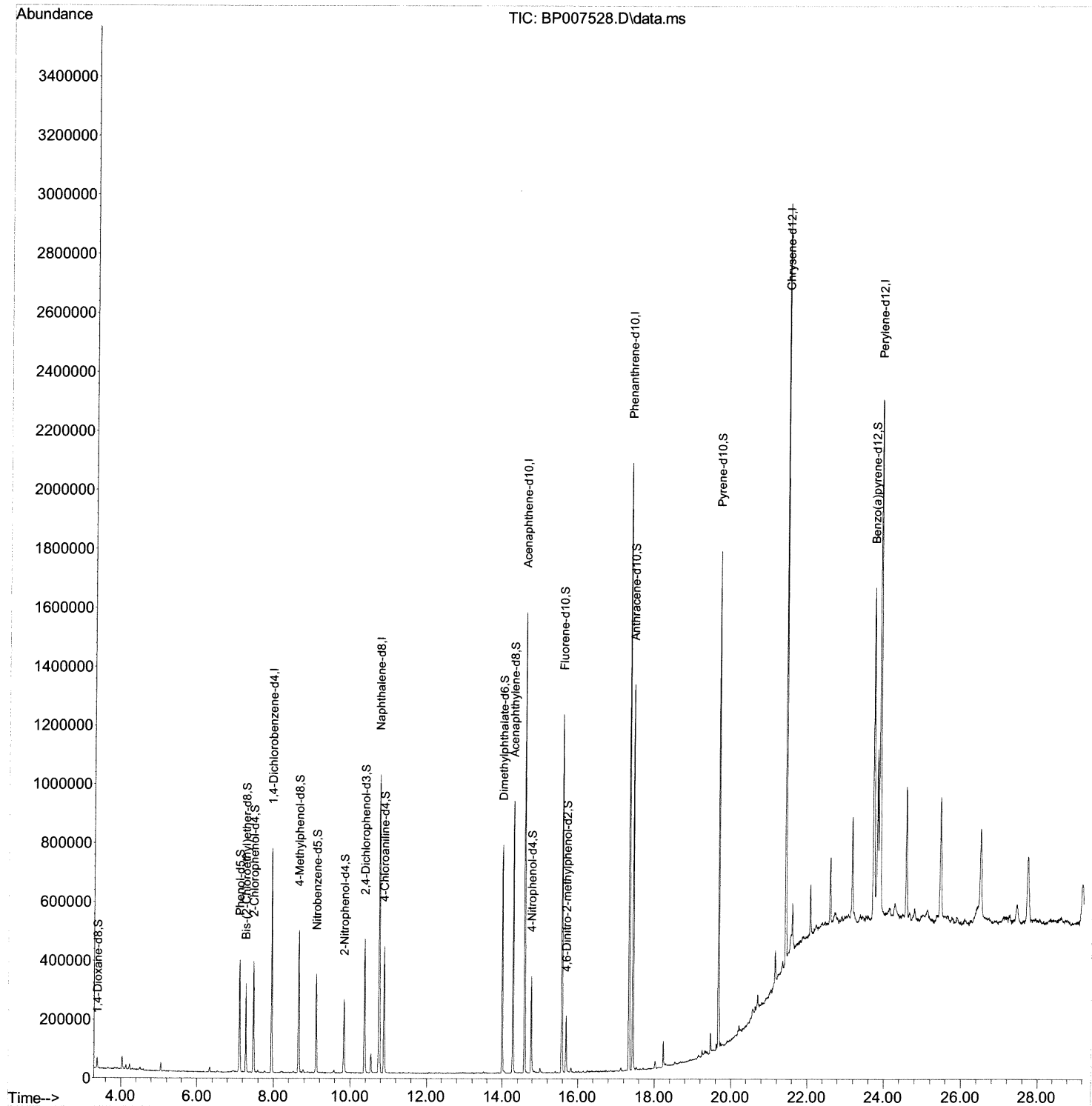
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007528.D
Acq On : 20 Oct 2021 14:43
Operator : CG/JU
Sample : M4218-05
Misc :
ALS Vial : 83 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGF2

Manual Integrations
APPROVED

mohammad
10/21/2021 11:15:48 AM

Quant Time: Oct 20 15:09:35 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)

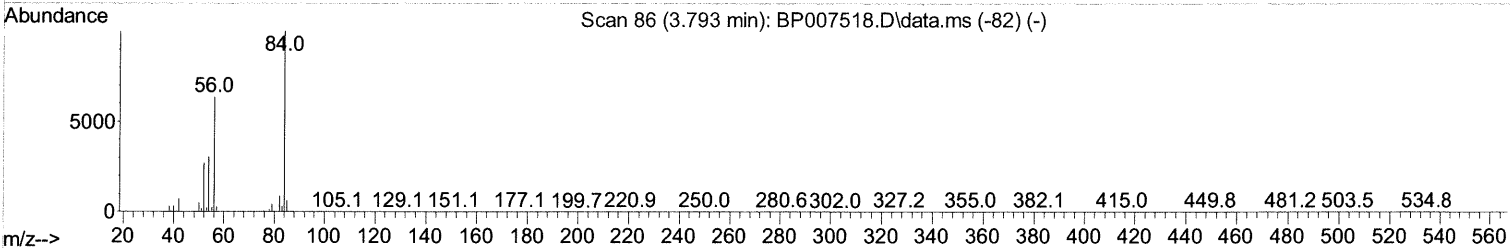
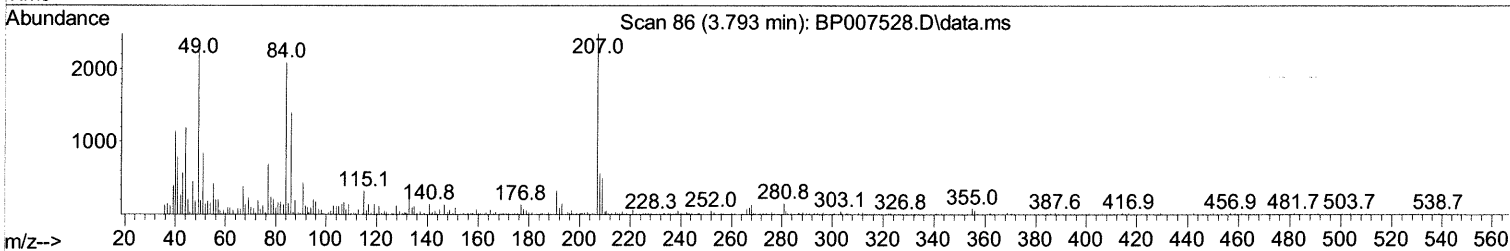
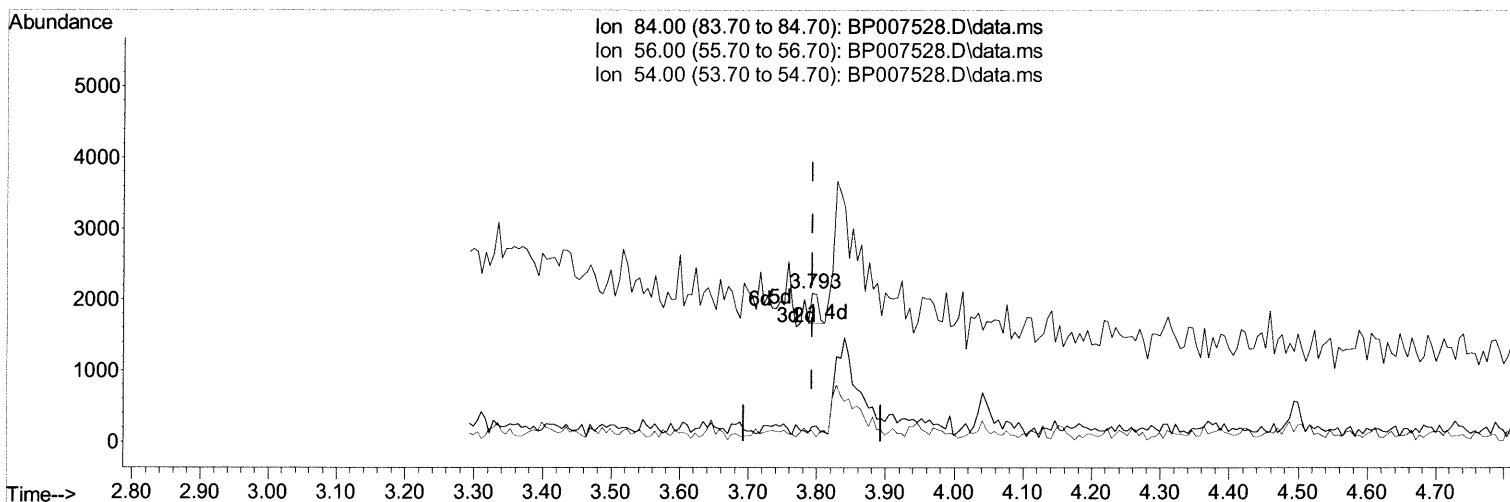
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TIC: BP007528.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.000) 0.02 ng/ul

response 313

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	9.66#
54.00	31.70	7.75#
0.00	0.00	0.00

Quantitation Report (Qedit)

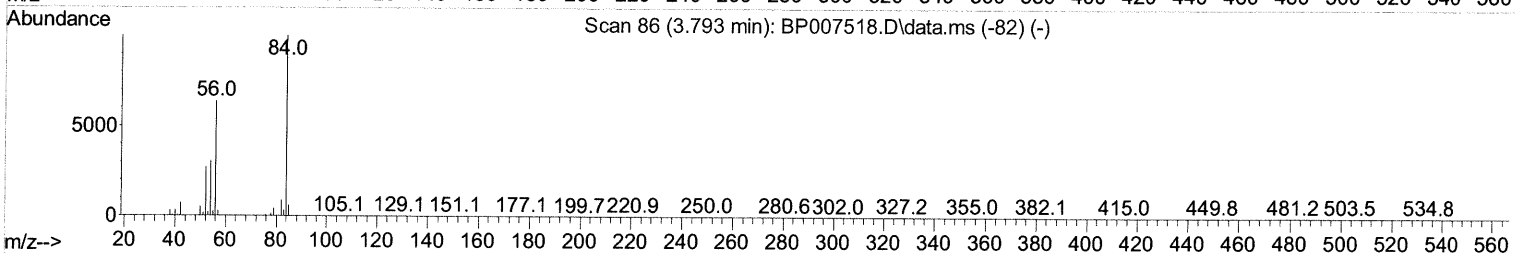
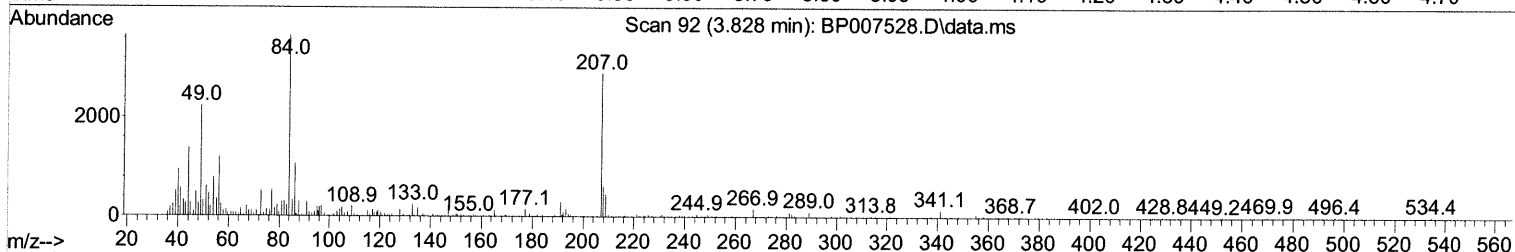
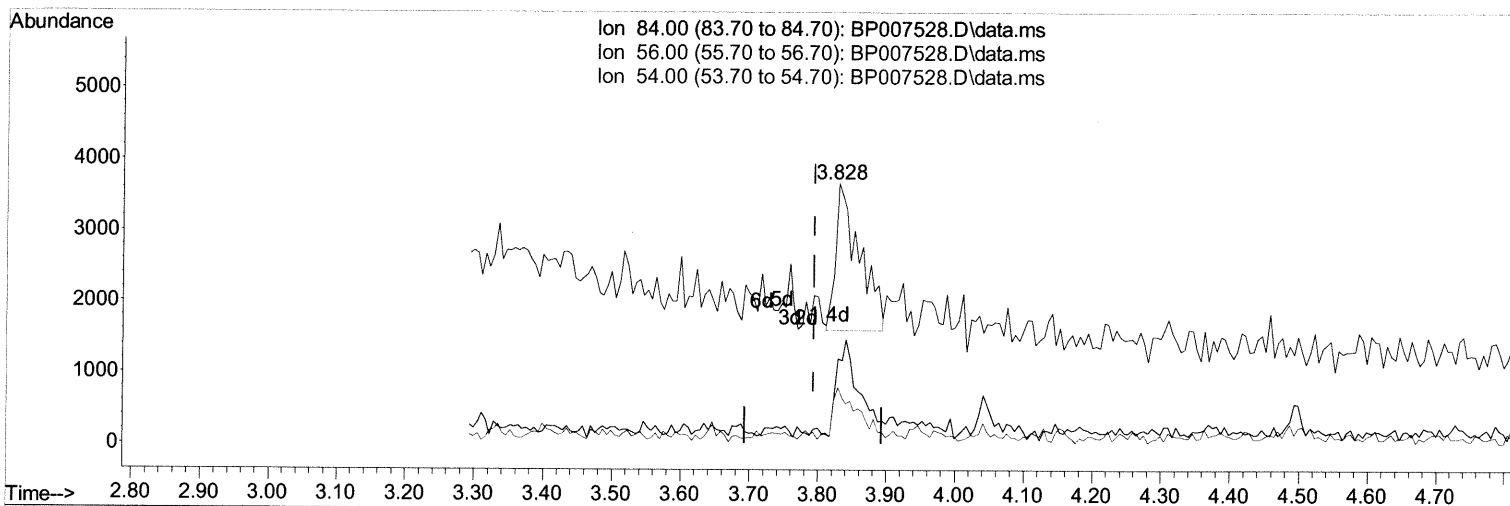
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TIC: BP007528.D\data.ms

(4) Pyridine-d5 (S)

3.828min (+ 0.035) 0.34 ng/ul m 10/20/21 JU

response 5025

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	32.72#
54.00	31.70	21.52#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007528.D
 Acq On : 20 Oct 2021 14:43
 Operator : CG/JU
 Sample : M4218-05
 Misc :
 ALS Vial : 83 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGF2

Manual Integrations
 APPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	211655	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	854244	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	555103	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1216640	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	1285074	20.000	ng/ul	0.00
88) Perylene-d12	23.880	264	1336833	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	18725	3.494	ng/ul	0.00
4) Pyridine-d5	3.828	84	5025m	0.343	ng/ul	0.04
7) Phenol-d5	7.111	99	219305	13.268	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	127282	12.911	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	176086	13.542	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	175995	13.823	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	79502	14.706	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	80744	15.990	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	169499	13.762	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	227177	13.640	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	527167	13.727	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	650288	13.474	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	73117	11.865	ng/ul	0.00
60) Fluorene-d10	15.581	176	464288	14.446	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	42458	8.709	ng/ul	0.00
73) Anthracene-d10	17.439	188	744091	14.401	ng/ul	0.00
81) Pyrene-d10	19.669	212	913005	14.212	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.715	264	892377	13.416	ng/ul	0.00

Target Compounds Qvalue

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