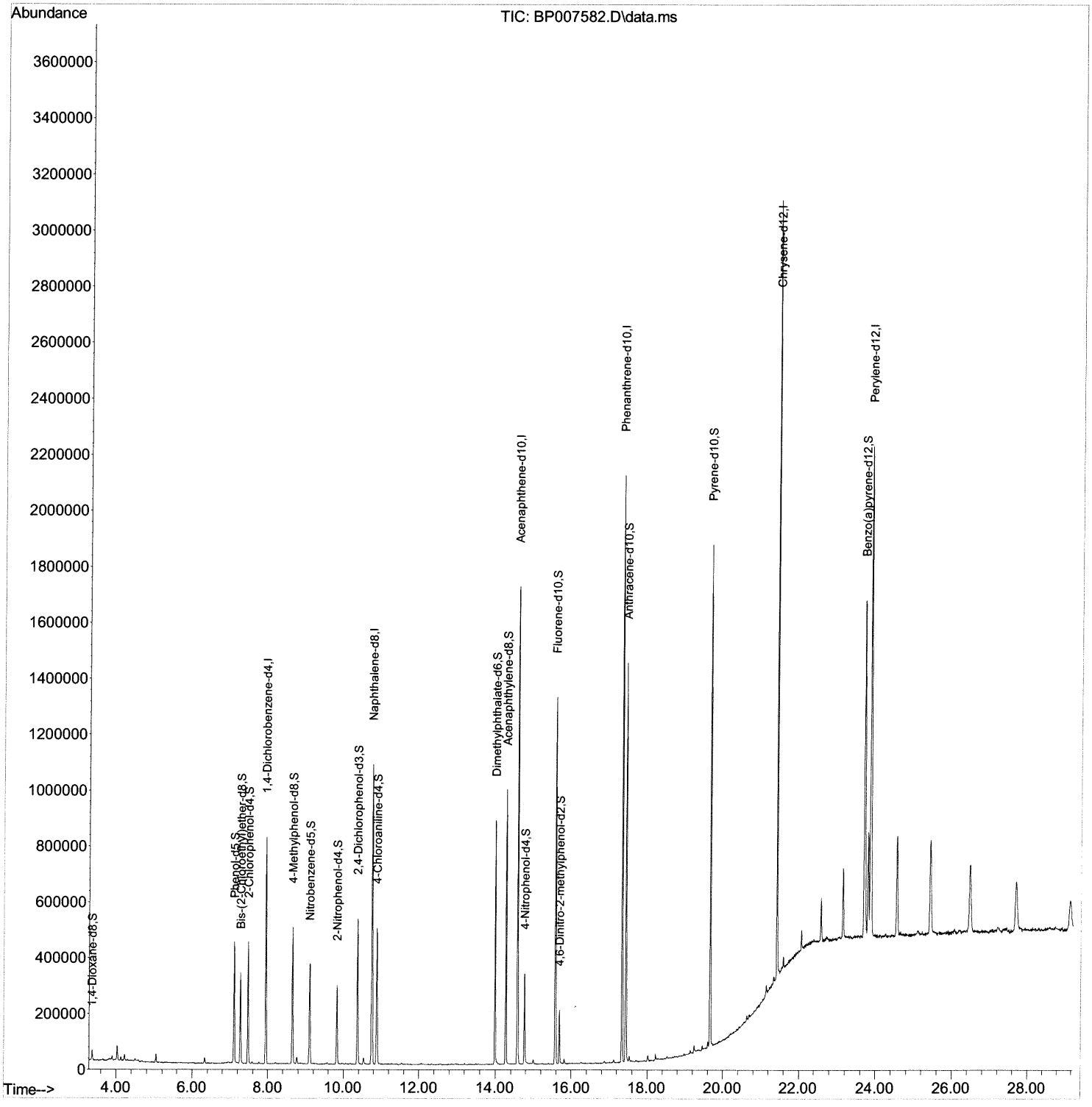
Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007582.D
Acq On : 22 Oct 2021 10:35
Operator : CG/JU
Sample : M4217-13
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Oct 22 13:00:58 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

Instrument :
BNA_P
ClientSampleId :
JDGE7

Manual Integrations APPROVED
mohammad
10/22/2021 2:55:14 PM



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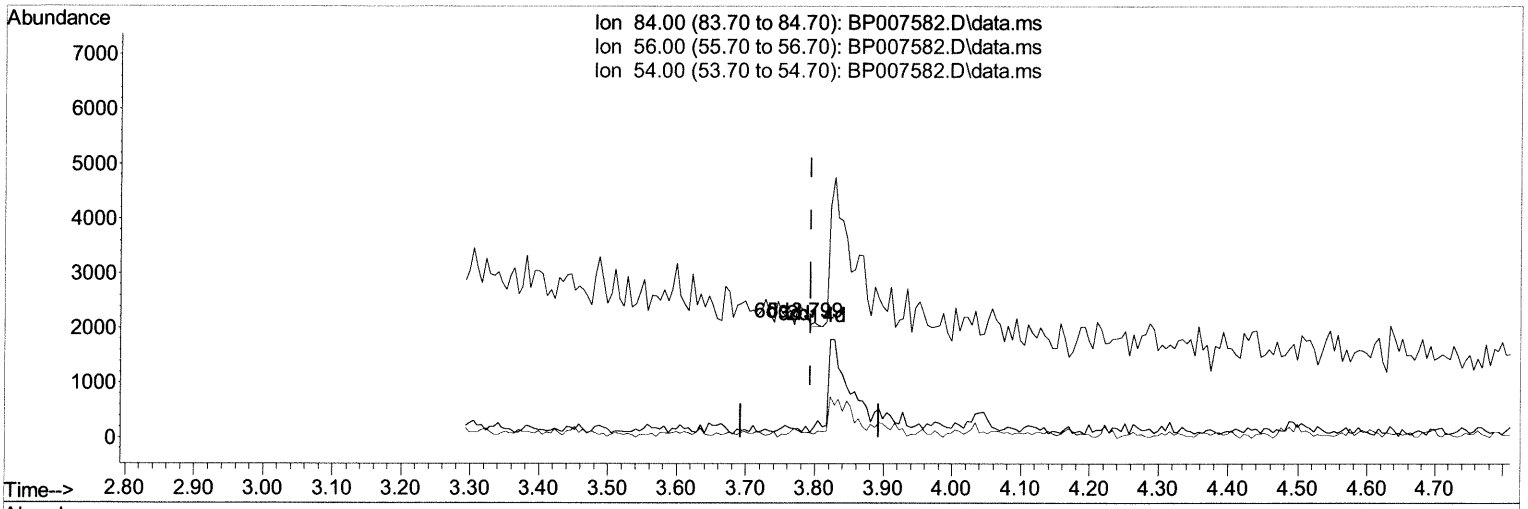

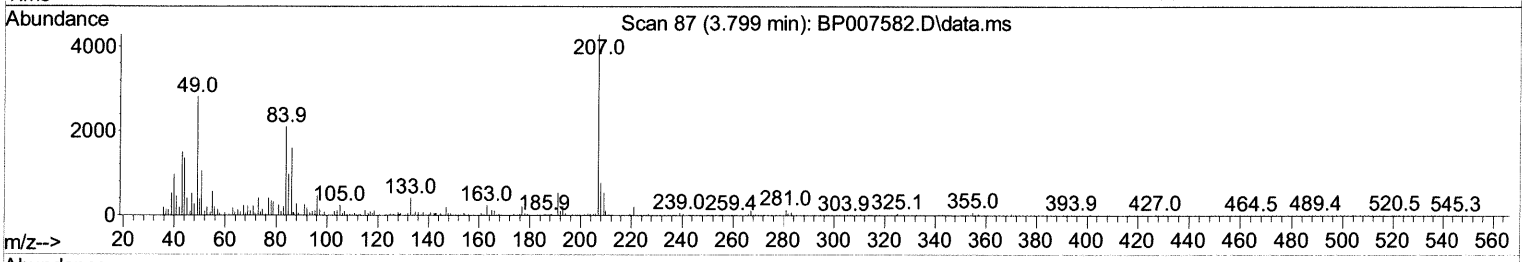

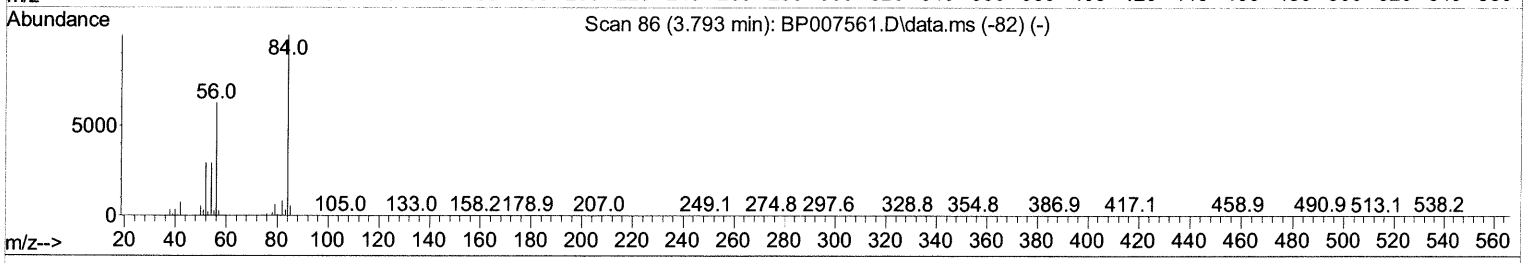

TIC: BP007582.D\data.ms

(4) Pyridine-d5 (S)

3.799min (+ 0.006) 0.00 ng/ul

response 36

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	9.54#
54.00	31.70	3.18#
0.00	0.00	0.00

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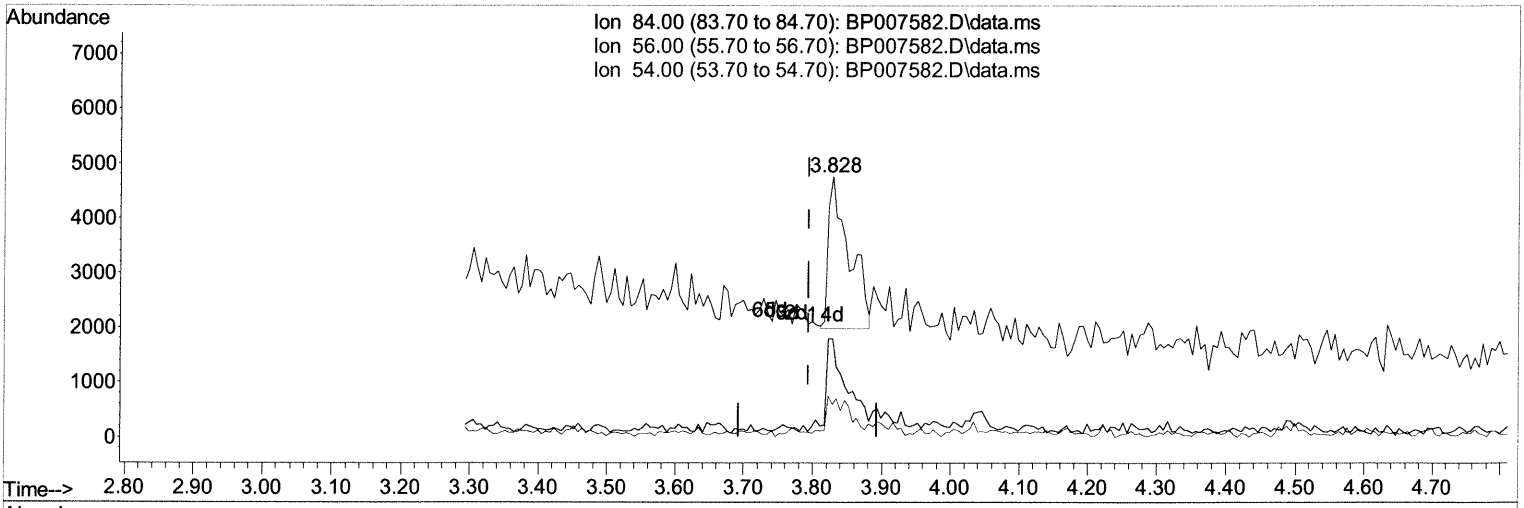

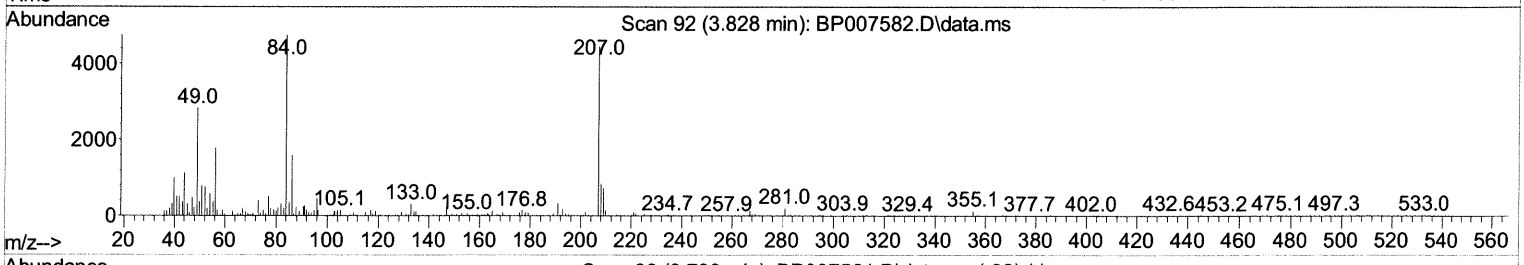

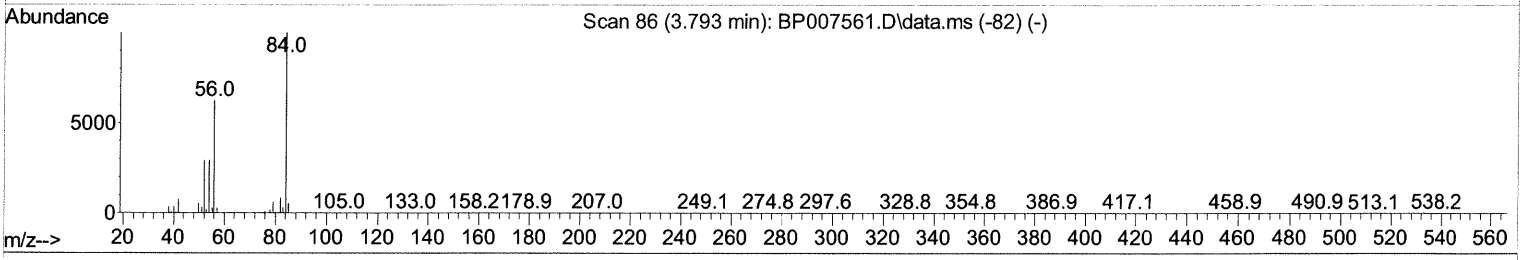

TIC: BP007582.D\data.ms

(4) Pyridine-d5 (S)

3.828min (+ 0.035) 0.38 ng/ul m 10/26/21JU

response 5815

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	37.54#
54.00	31.70	12.32#
0.00	0.00	0.00

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Manual Integrations APPROVED
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10/22/2021 2:55:14 PM

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	221510	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	899855	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	586454	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	1246667	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	1304487	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	1343164	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	17213	3.069	ng/uL	0.00
4) Pyridine-d5	3.828	84	5815m	> 0.379	ng/ul	> 0.04 10/26/21 JU
7) Phenol-d5	7.105	99	244625	14.141	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	140226	13.591	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	195730	14.383	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	181851	13.648	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	90278	15.853	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	92001	17.296	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	189016	14.569	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	256990	14.648	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	592099	14.593	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	714605	14.015	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	77706	11.936	ng/ul	-0.01
60) Fluorene-d10	15.575	176	524210	15.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	43341	8.676	ng/ul	0.00
73) Anthracene-d10	17.433	188	816269	15.418	ng/ul	0.00
81) Pyrene-d10	19.663	212	985026	15.105	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	940315	14.070	ng/ul	-0.01

Target Compounds Qvalue

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