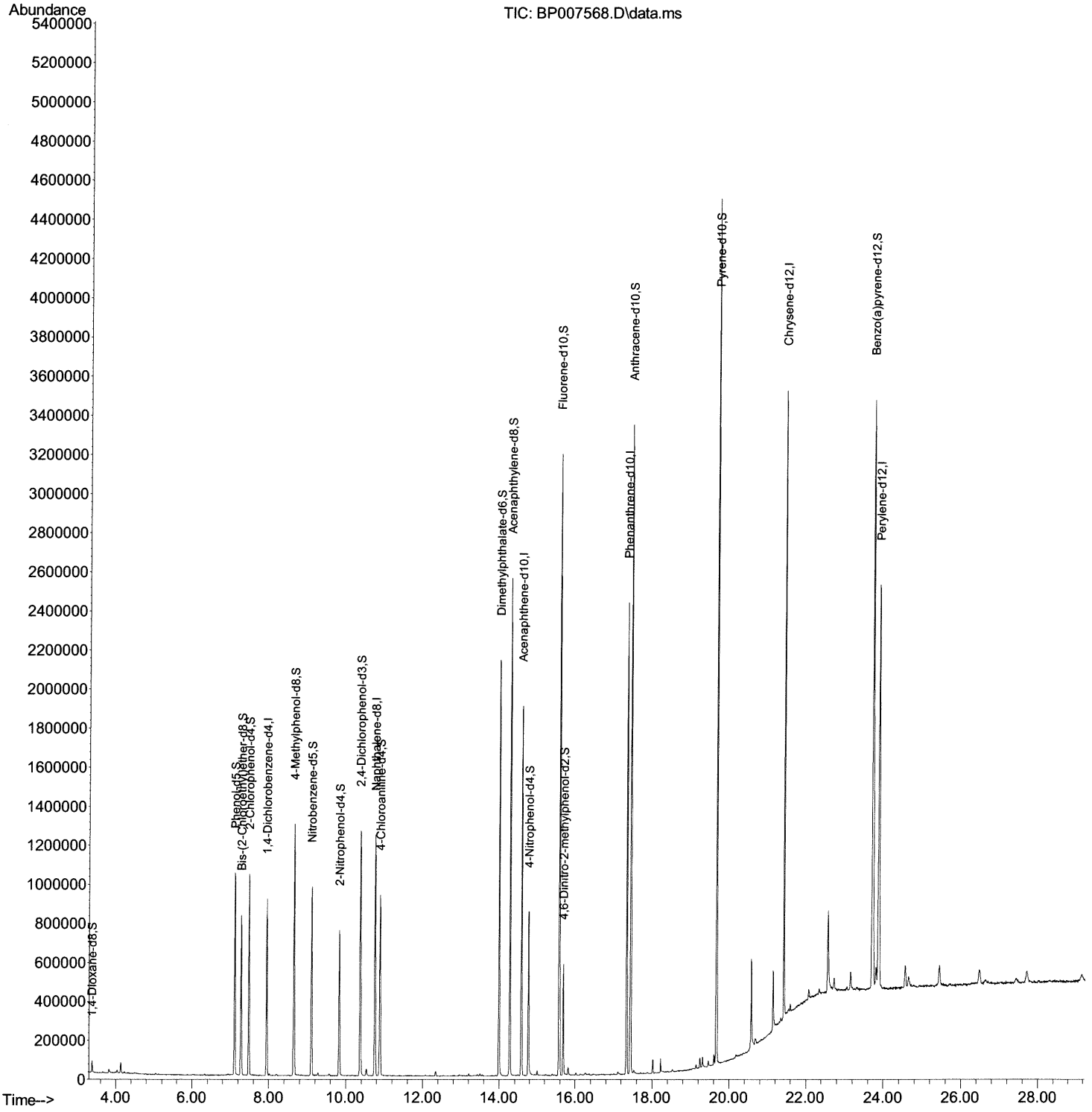
Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007568.D
Acq On : 21 Oct 2021 23:45
Operator : CG/JU
Sample : M4217-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Quant Time: Oct 22 01:50:52 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 :
JDGB6

Manual Integrations APPROVED
mohammad
10/22/2021 2:55:00 PM



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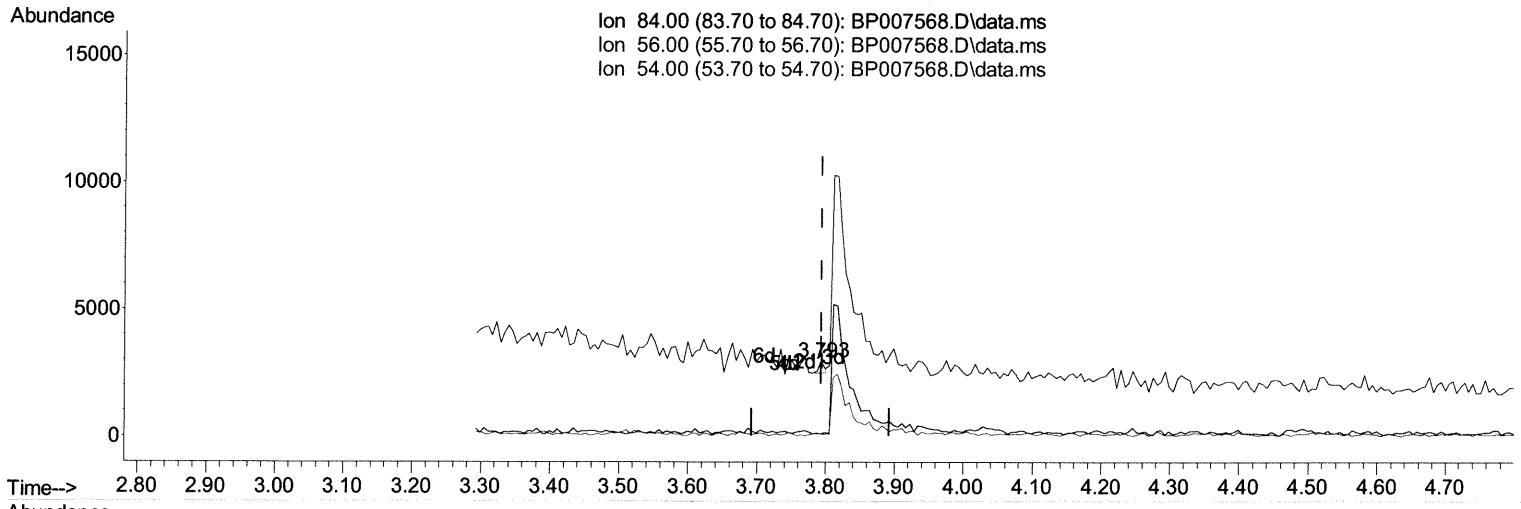

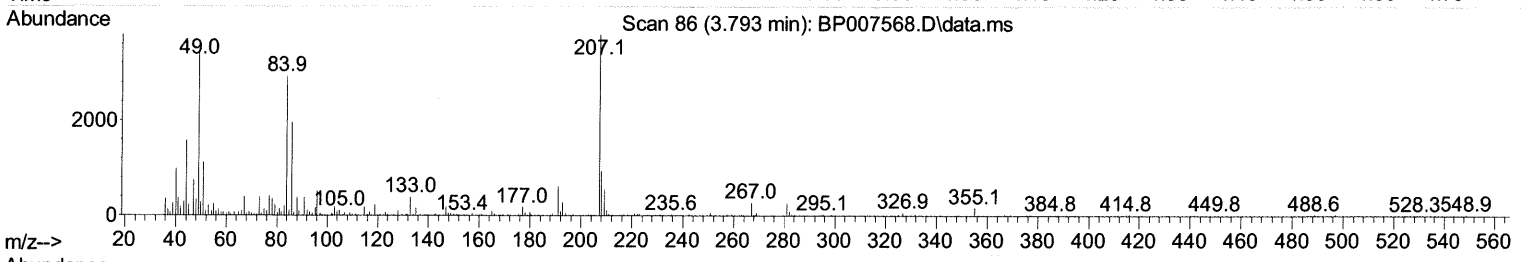

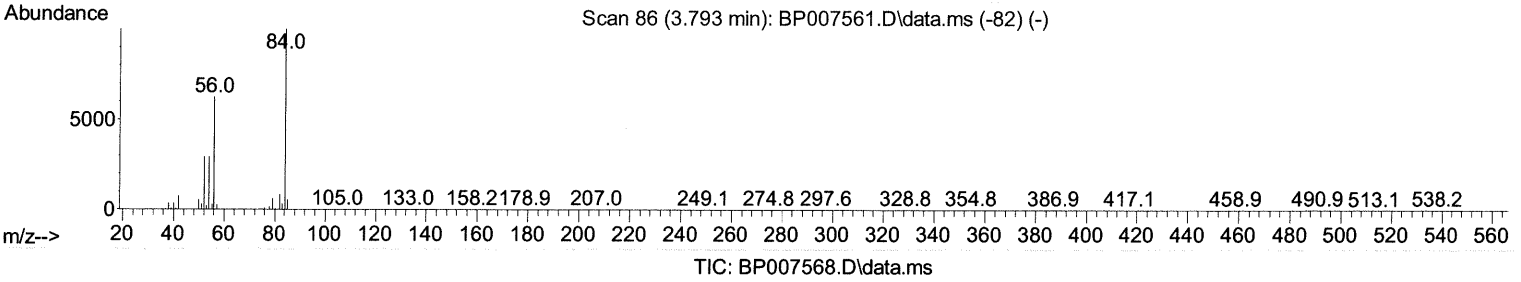

TIC: BP007568.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.000) 0.01 ng/ul

response 239

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	3.18#
54.00	31.70	3.25#
0.00	0.00	0.00

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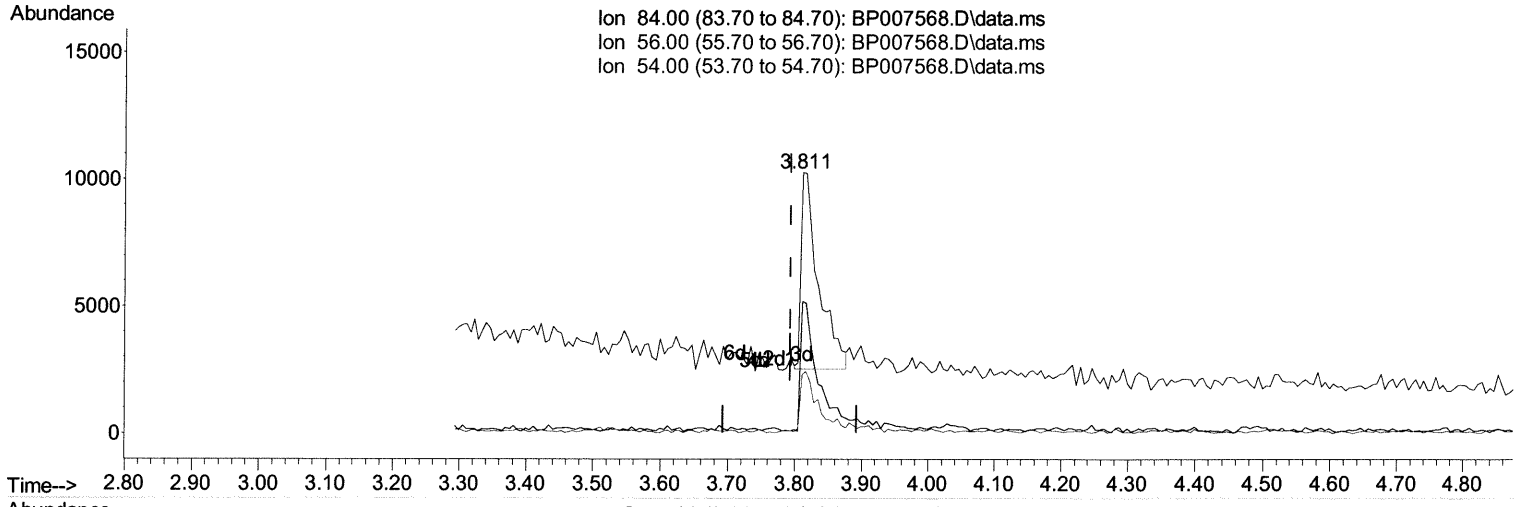

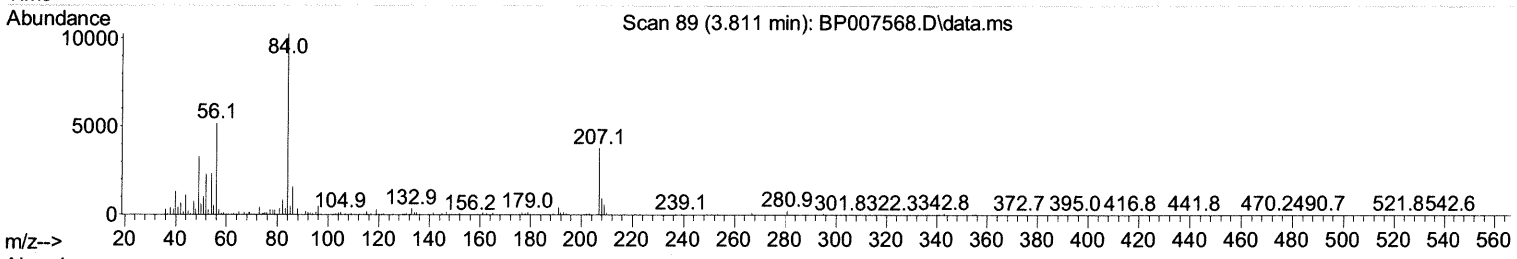

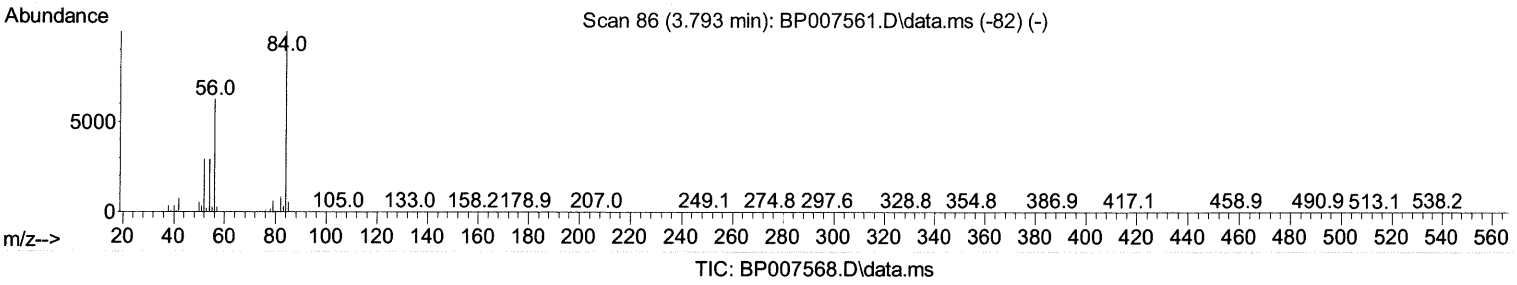

TIC: BP007568.D\data.ms

(4) Pyridine-d5 (S)

3.811min (+ 0.018) 0.81 ng/ul m 10/26/21 JJ

response 13829

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	50.61#
54.00	31.70	22.56#
0.00	0.00	0.00

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Manual Integrations APPROVED
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10/22/2021 2:55:00 PM

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	247169	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	1018324	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	663076	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.334	188	1435932	20.000	ng/ul	0.00
79) Chrysene-d12	21.422	240	1545628	20.000	ng/ul	0.00
88) Perylene-d12	23.869	264	1594537	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	28082	4.487	ng/uL	0.00
4) Pyridine-d5	3.811	84	13829m	> 0.807	ng/ul	> 0.02 10/22/2021
7) Phenol-d5	7.105	99	601127	31.142	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	346062	30.059	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	475622	31.322	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	474211	31.894	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128	228934	35.524	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	237983	39.535	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	472453	32.179	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	492806	24.821	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1386537	30.225	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1767620	30.660	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	200979	27.304	ng/ul	0.00
60) Fluorene-d10	15.575	176	1239411	32.285	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	124824	21.695	ng/ul	0.00
73) Anthracene-d10	17.434	188	1957802	32.105	ng/ul	0.00
81) Pyrene-d10	19.669	212	2359770	30.540	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	2376716	29.956	ng/ul	0.00

Target Compounds Qvalue

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