

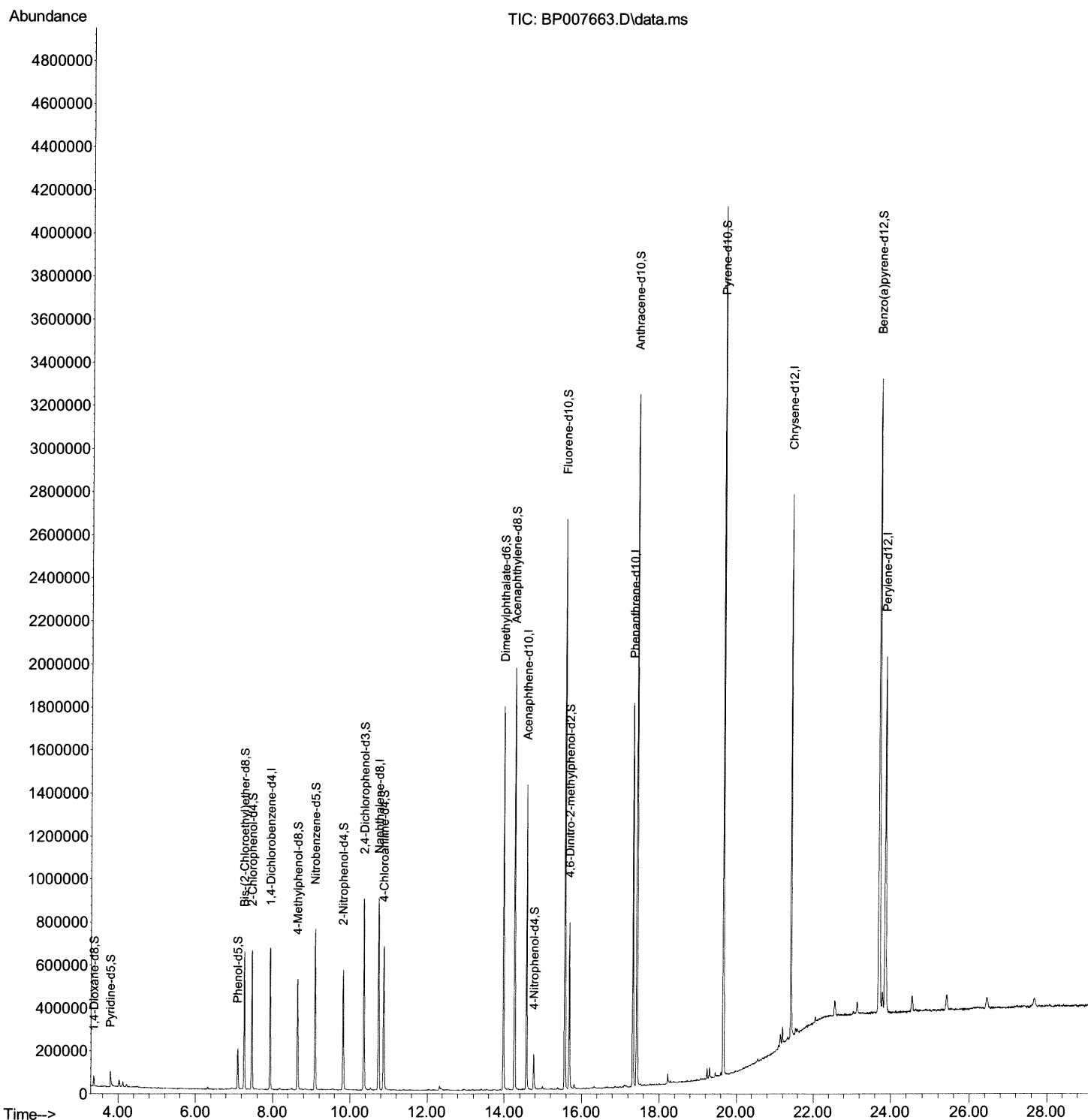
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007663.D
Acq On : 26 Oct 2021 03:52
Operator : CG/JU
Sample : M4291-08
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P009-TW001-02

Manual Integrations
APPROVED

mohammad
10/26/2021 11:34:55 AM

Quant Time: Oct 26 04:36:15 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 25 16:57:16 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

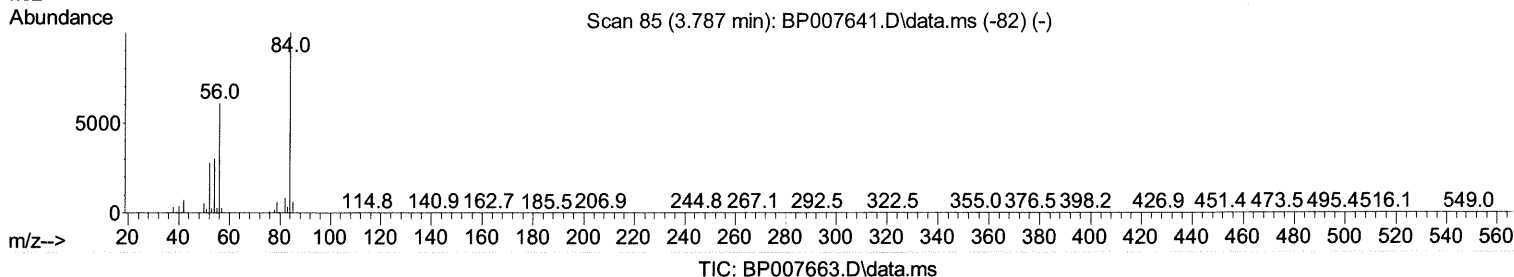
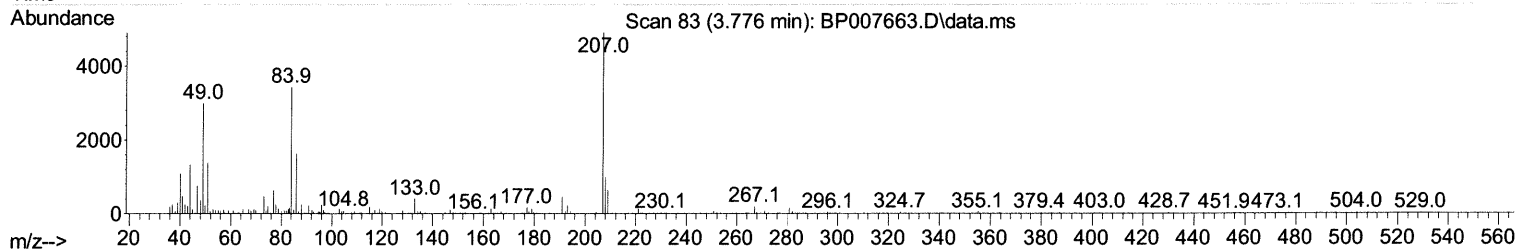
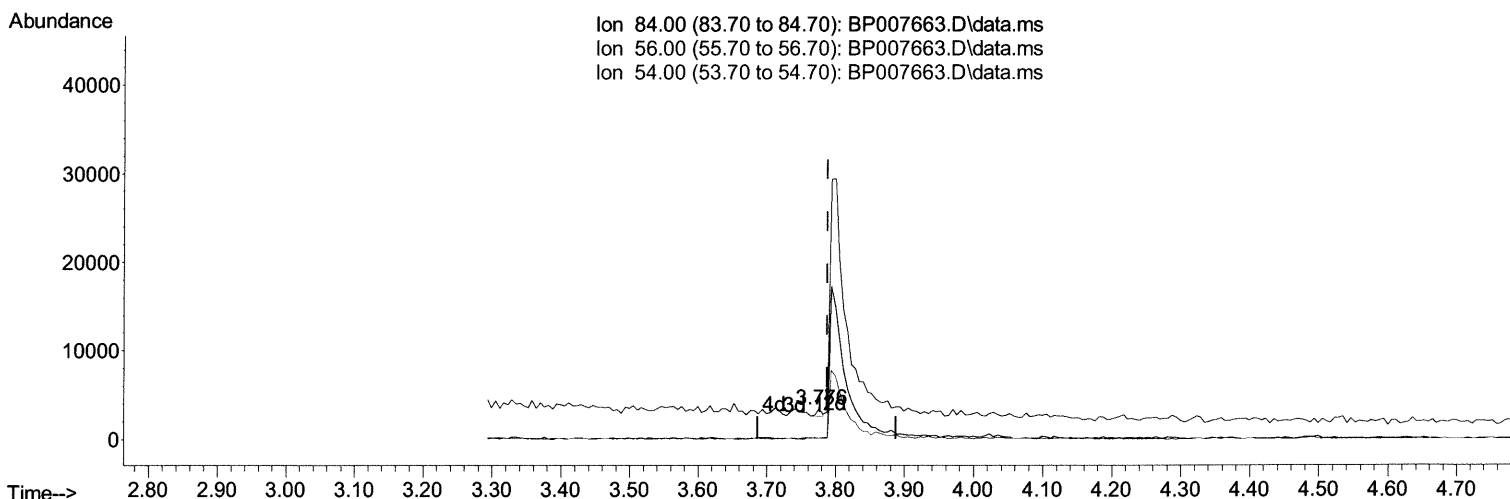
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(4) Pyridine-d5 (S)

3.776min (-0.012) 0.04 ng/ul

response 442

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	2.36#
54.00	31.70	3.09#
0.00	0.00	0.00

Quantitation Report (Qedit)

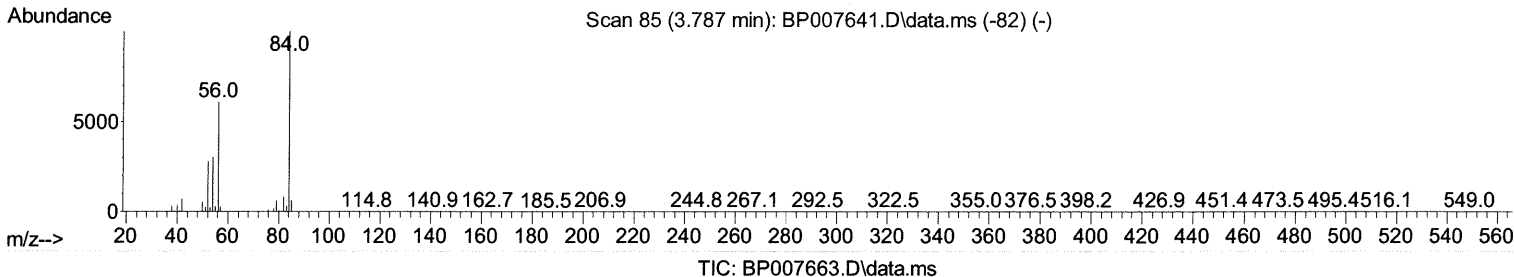
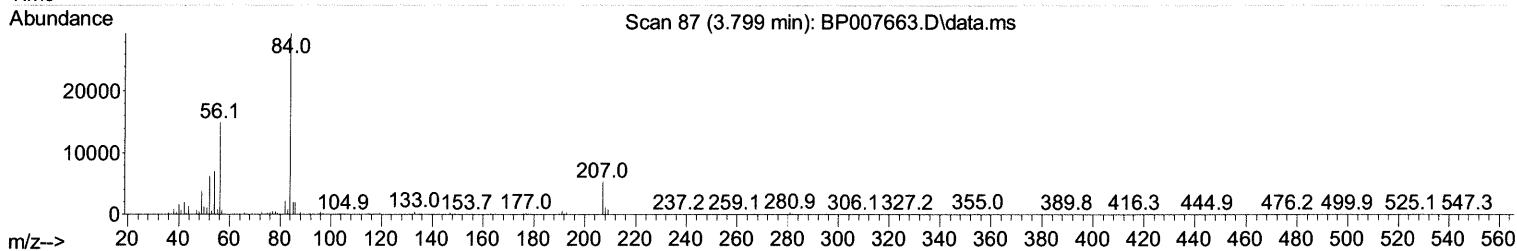
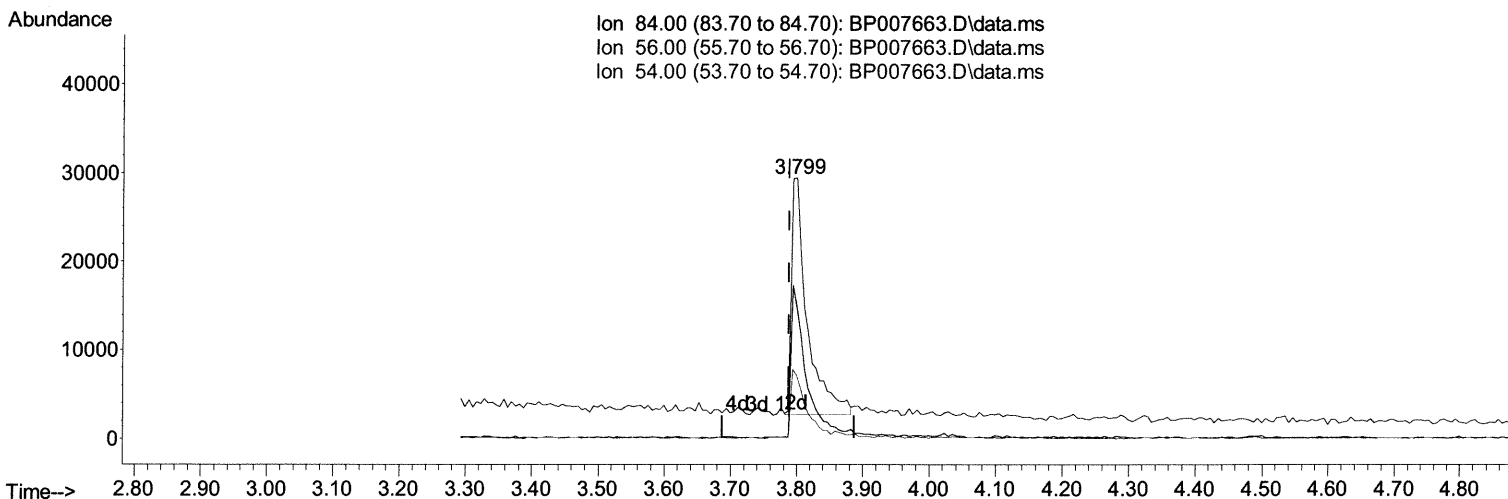
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Manual Integrations
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Response via : Initial Calibration



(4) Pyridine-d5 (S)

3.799min (+ 0.012) 3.54 ng/ul m 10/27/21 JU

response 43683

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	50.73#
54.00	31.70	24.19#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007663.D
 Acq On : 26 Oct 2021 03:52
 Operator : CG/JU
 Sample : M4291-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P009-TW001-02

Manual Integrations
 APPROVED

mohammad
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Quant Time: Oct 26 04:36:15 2021
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 25 16:57:16 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	185248	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	738388	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	495574	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1082420	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1200784	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	1272978	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	24521	6.013	ng/ul	0.00
4) Pyridine-d5	3.799	84	43683m	3.542	ng/ul	0.01
7) Phenol-d5	7.099	99	107449	7.636	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	272923	34.936	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	299233	27.343	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	193165	17.532	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	178605	35.495	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	181864	33.091	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.358	165	334793	29.928	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.875	131	362223	24.489	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1221345	36.383	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1397308	33.961	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	40755	6.623	ng/ul	-0.01
60) Fluorene-d10	15.569	176	1053298	37.219	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	172874	29.858	ng/ul	0.00
73) Anthracene-d10	17.428	188	1847657	41.536	ng/ul	0.00
81) Pyrene-d10	19.657	212	2274901	40.692	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	2391175	38.238	ng/ul	0.00

Target Compounds Qvalue

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