

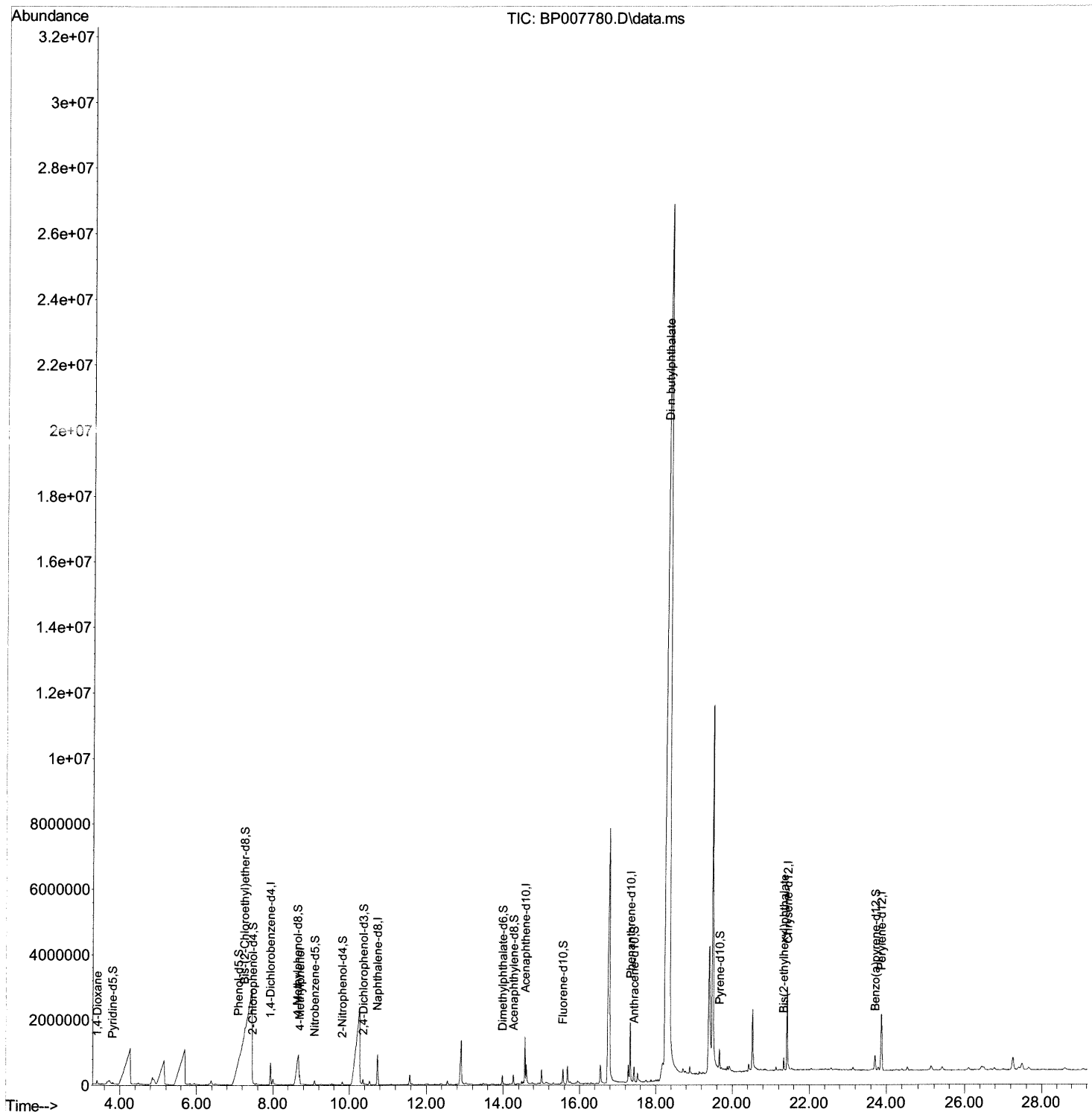
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007780.D
Acq On : 29 Oct 2021 11:52
Operator : CG/JU
Sample : M4281-09DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY65DL

Manual Integrations
APPROVED

mohammad
11/1/2021 1:36:21 PM

Quant Time: Oct 29 12:21:08 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 27 15:37:43 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

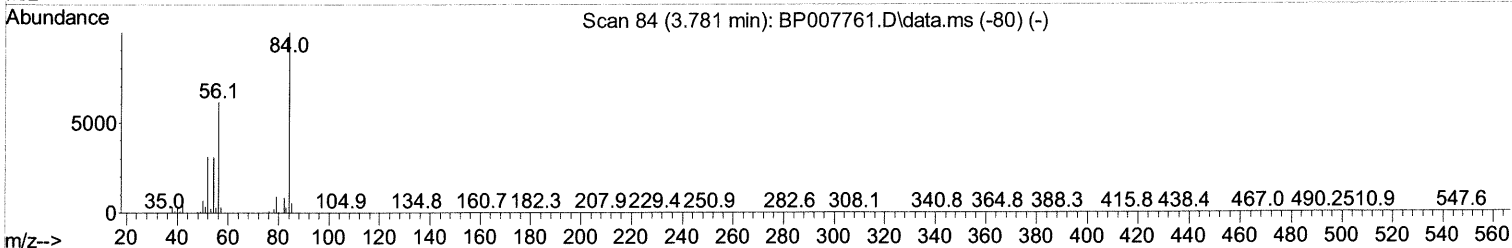
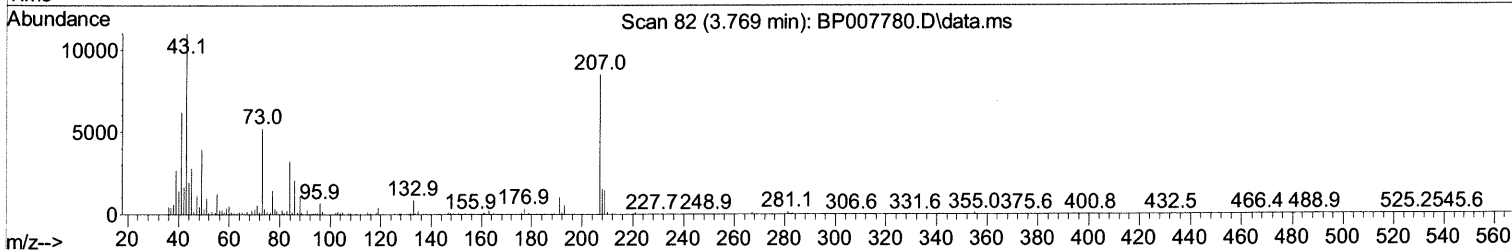
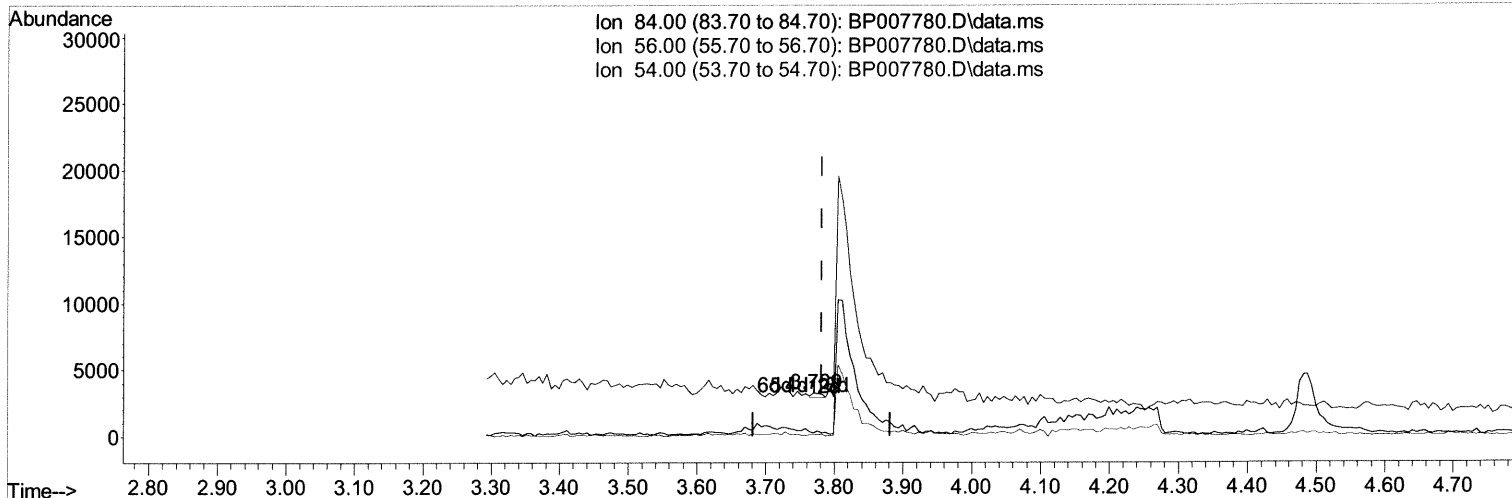
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TIC: BP007780.D\data.ms

(4) Pyridine-d5 (S)

3.769min (-0.012) 0.02 ng/ul

response 302

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	8.37#
54.00	31.70	3.27#
0.00	0.00	0.00

Quantitation Report (Qedit)

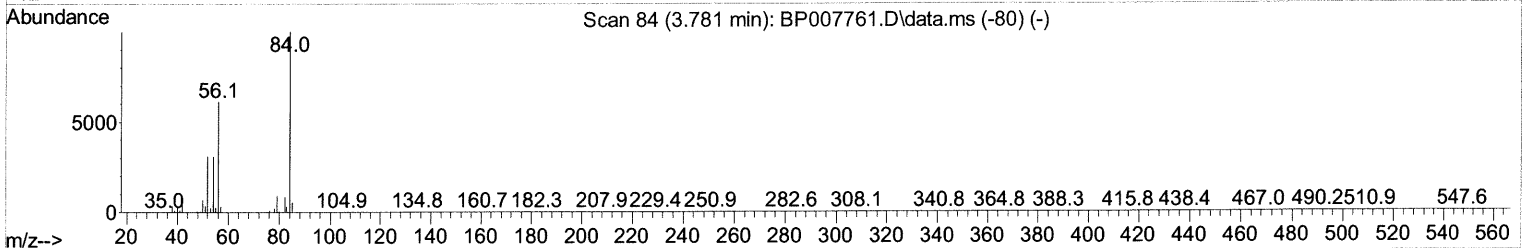
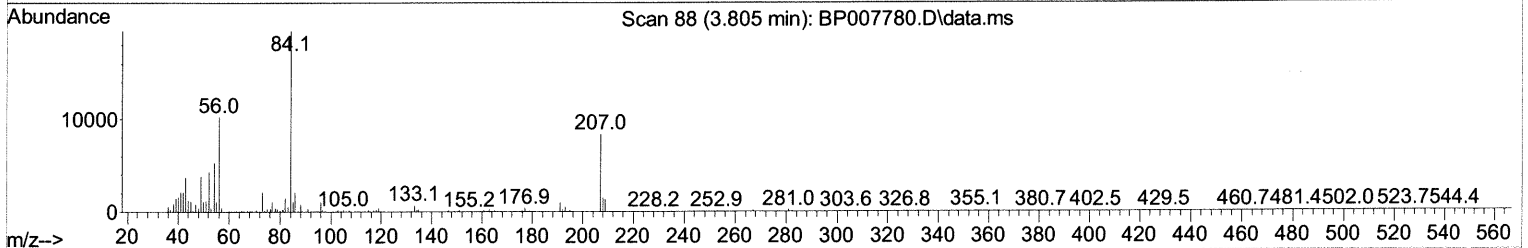
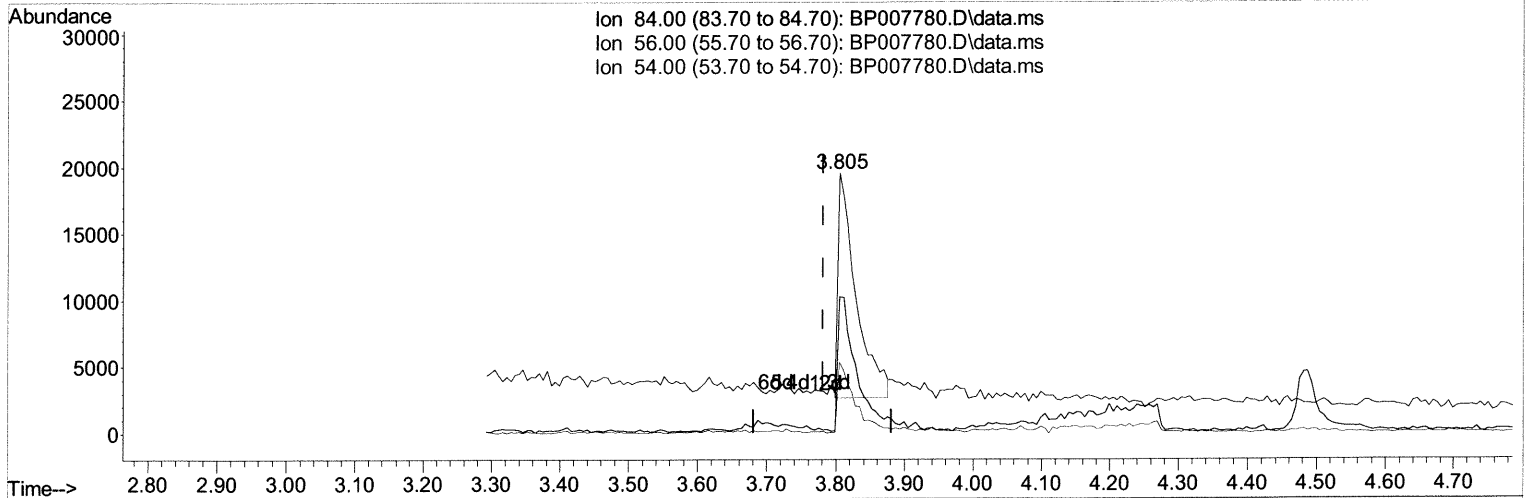
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Manual Integrations
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TIC: BP007780.D\data.ms

(4) Pyridine-d5 (S)

3.805min (+ 0.024) 2.42 ng/ul m 11/02/21 JU

response 30477

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	52.42#
54.00	31.70	27.56
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
 Data File : BP007780.D
 Acq On : 29 Oct 2021 11:52
 Operator : CG/JU
 Sample : M4281-09DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampled :
 BFY65DL

Manual Integrations
 APPROVED

mohammad
 11/1/2021 1:36:21 PM

Quant Time: Oct 29 12:21:08 2021
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	189140	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	754657	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	498911	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1069162	20.000	ng/ul	0.00
79) Chrysene-d12	21.421	240	1237225	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1362583	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	2167	0.520	ng/uL	0.00
4) Pyridine-d5	3.805	84	30477m	2.420	ng/ul	0.02 11/02/21 JV
7) Phenol-d5	7.099	99	20627	1.436	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	47271	5.927	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	48364	4.328	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	41206	3.663	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	29274	5.692	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	29347	5.225	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	55413	4.847	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	193716	5.732	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	217898	5.261	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	0.00
60) Fluorene-d10	15.563	176	174425	6.122	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	0.00
73) Anthracene-d10	17.422	188	263869	6.005	ng/ul	0.00
81) Pyrene-d10	19.663	212	320308	5.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	344075	5.140	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.405	88	59480	12.778	ng/uL	89
18) 4-Methylphenol	8.699	108	96745	8.274	ng/ul	84
78) Di-n-butylphthalate	18.292	149	96890	1.783	ng/ul#	91
86) Bis(2-ethylhexyl)phtha...	21.333	149	137090	3.319	ng/ul	98

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