

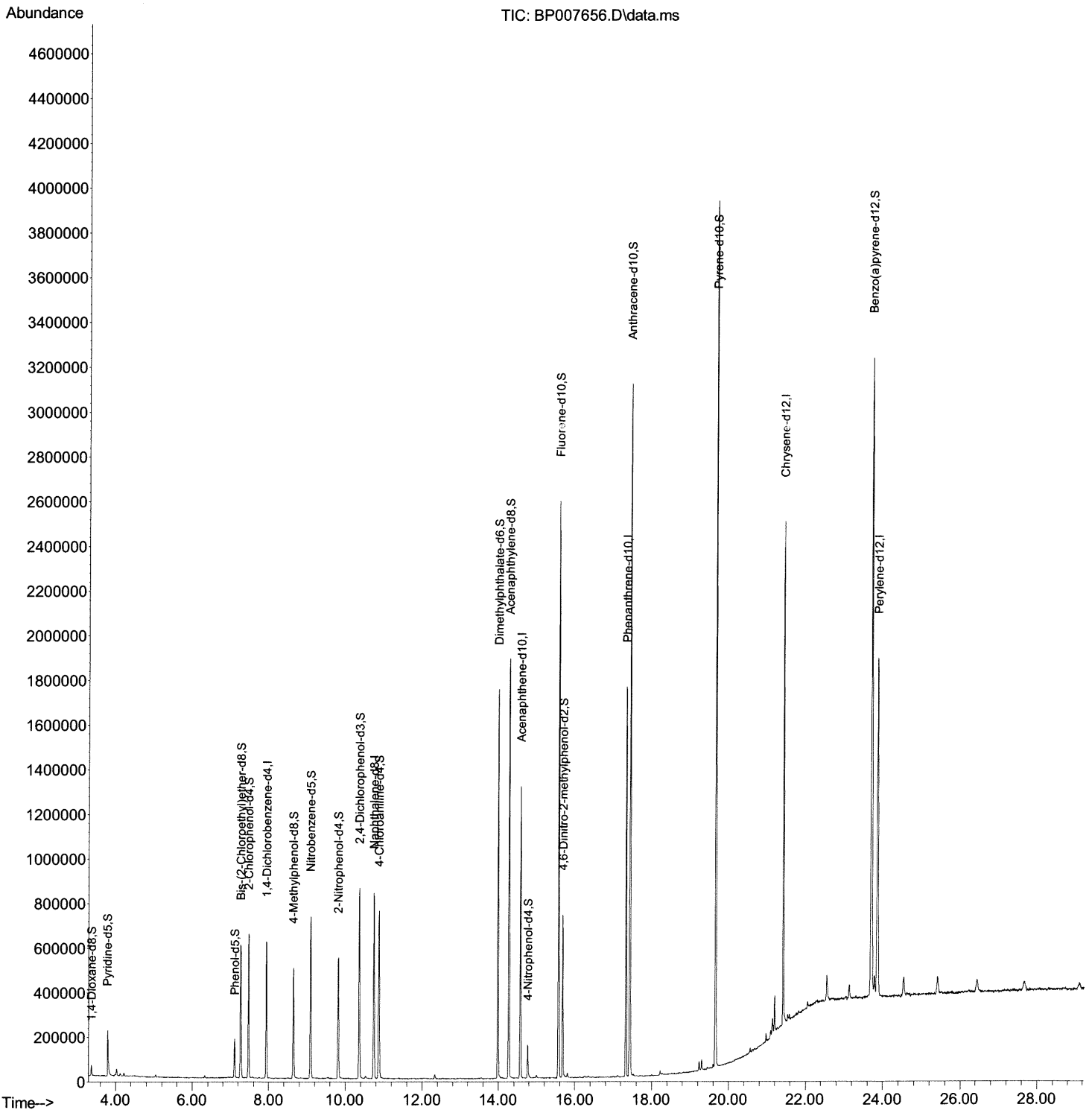
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007656.D
Acq On : 25 Oct 2021 23:55
Operator : CG/JU
Sample : M4303-21
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P055-TW001-01

Manual Integrations
APPROVED

mohammad
10/26/2021 11:34:46 AM

Quant Time: Oct 26 02:39:52 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)

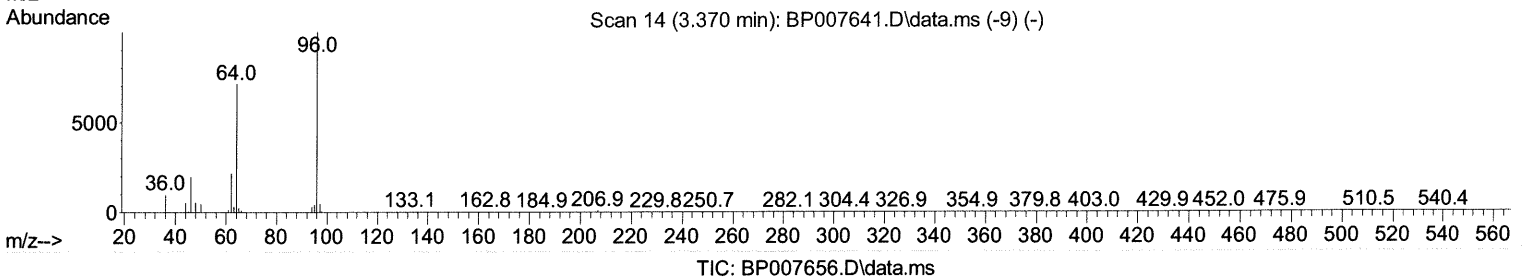
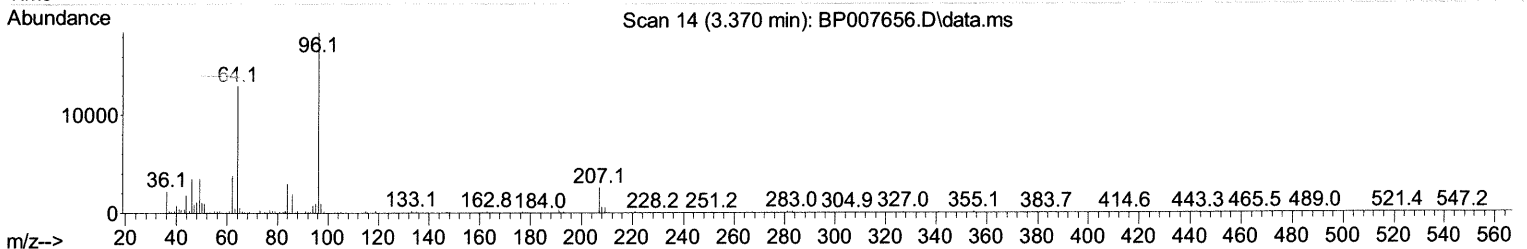
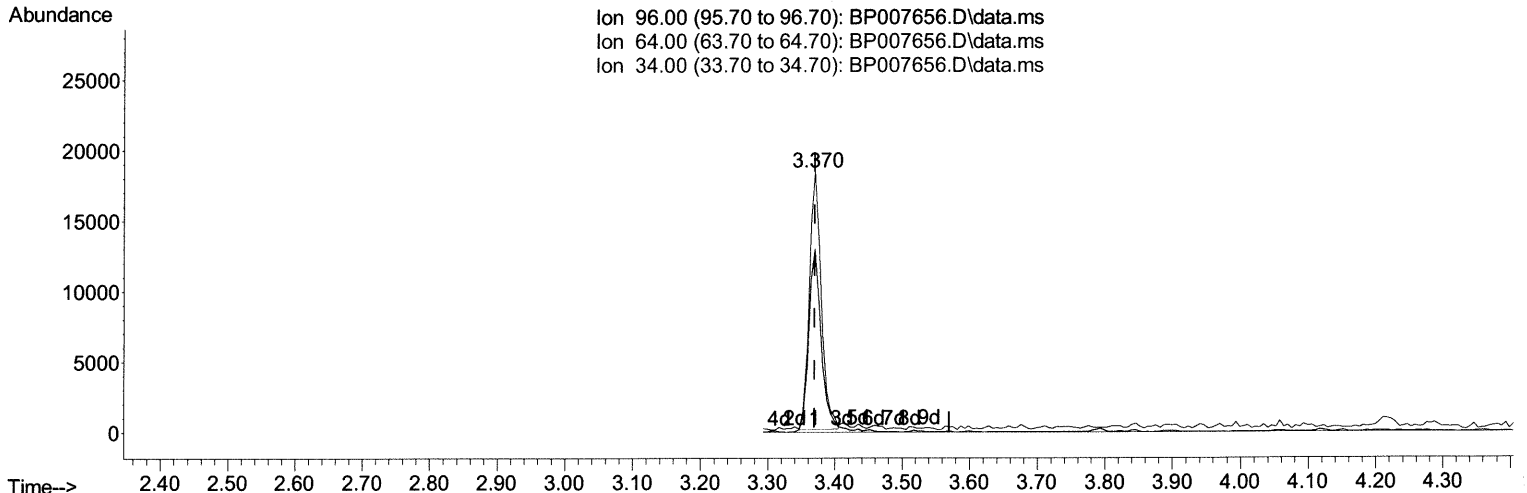
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 6.08 ng/uL

response 22967

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	70.43
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

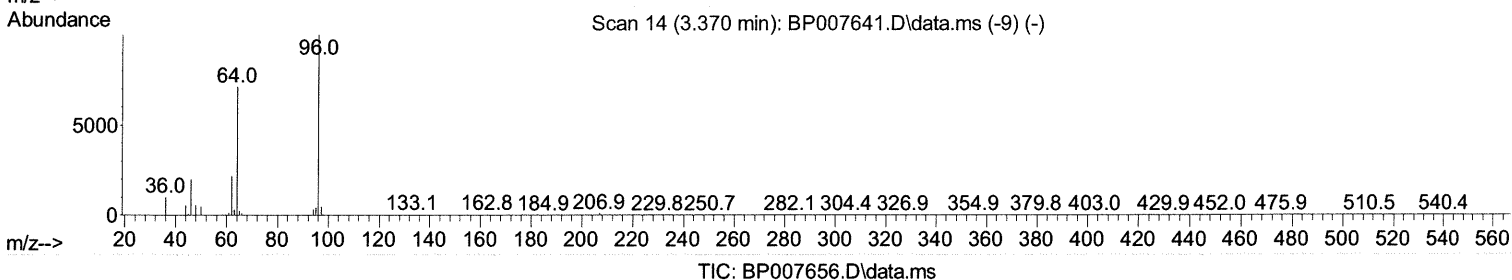
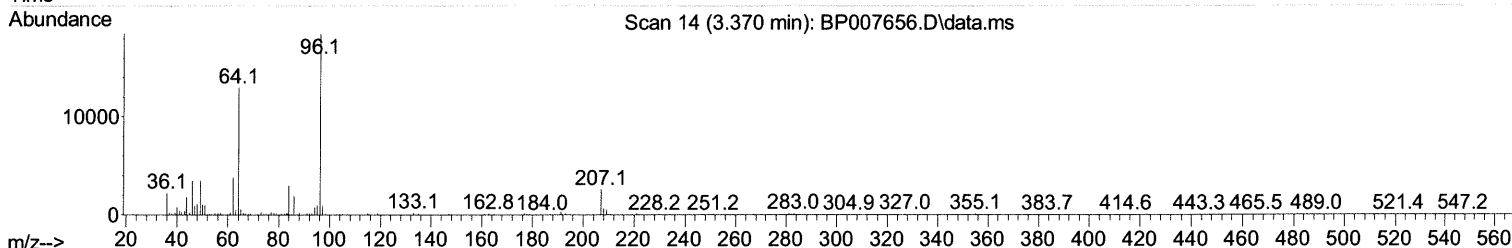
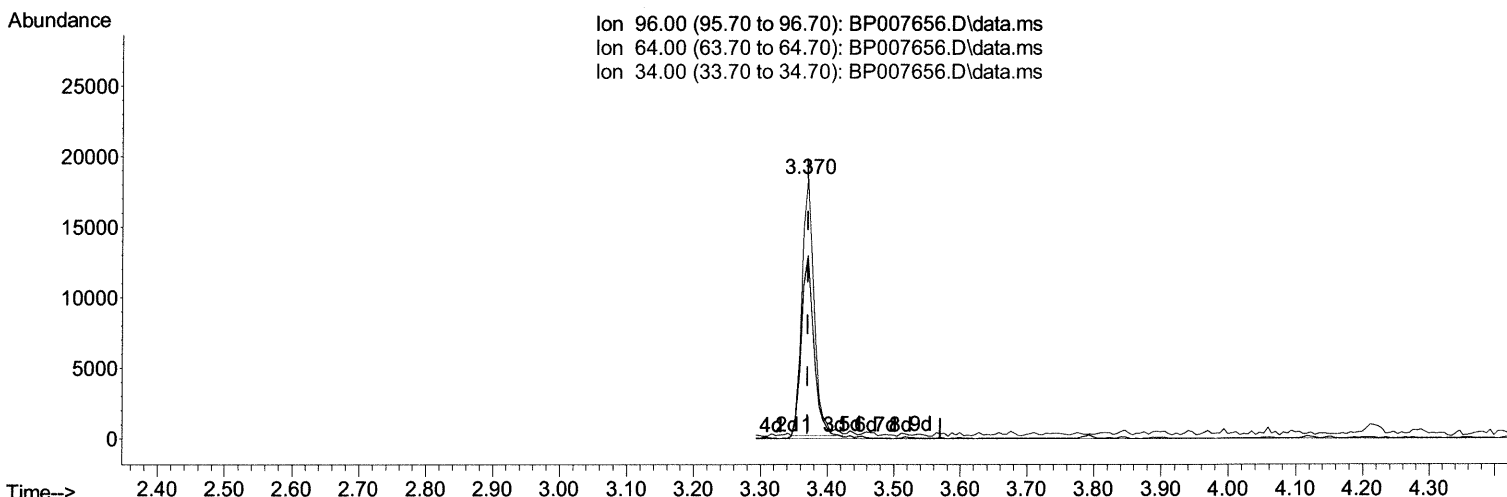
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 6.16 ng/uL m 10/27/21 JU

response 23256

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	70.43
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
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Manual Integrations
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	171543	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	695233	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	460644	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1000993	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1114134	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	1166950	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	23256m>	6.158	ng/ul >	0.00 10/27/21 JU
4) Pyridine-d5	3.787	84	117420	10.282	ng/ul	0.00
7) Phenol-d5	7.105	99	103060	7.910	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	262327	36.263	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	289880	28.605	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	183802	18.015	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	173024	36.520	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	179326	34.655	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	318637	30.252	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	408396	29.324	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	1190091	38.141	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1367223	35.750	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	38130	6.666	ng/ul	-0.01
60) Fluorene-d10	15.569	176	1031937	39.230	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	162708	30.388	ng/ul	0.00
73) Anthracene-d10	17.428	188	1734706	42.169	ng/ul	0.00
81) Pyrene-d10	19.663	212	2143877	41.331	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	2257523	39.381	ng/ul	-0.01

Target Compounds Qvalue

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